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> DEC 8 2000 EC 18 2000

December 18, 2000

Mr. Charles Schwer Ms. Maria Stadlmayer Hazardous Materials Management Division Agency of Natural Resources 103 South Main Street Waterbury, VT 05671-0404

OMYA Florence Plant, Spill Investigation RE:

Dear Chuck and Maria:

Enclosed please find Heindel and Noyes' report of the spill at OMYA's Verpol plant in Florence.

If you have comments or questions, please feel welcome to call me at (802) 658-0820.

Sincerely,

Jeffrey Noyes

Chief Hydrogeologist

Neal Jordan CC:

Enclosure

OMYA, INC. VERPOL PLANT

CONTAMINANT SPILL INVESTIGATION

December 18, 2000

HEINDEL AND NOYES

Consulting Hydrogeologists, Engineers, and Environmental Scientists



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• Engineers

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OMYA, Inc. Verpol Plant Florence, Vermont

CONTAMINANT SPILL INVESTIGATION

November 2000

Prepared by:

Heindel and Noyes

Prepared for:





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OMYA, Inc. Verpol Plant

Florence, Vermont

CONTAMINANT SPILL INVESTIGATION

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EXECUTIVE SUMMARY

- A release of 4,500 gallons of Biocide occurred on November 18, 2000, due to an equipment failure at OMYA's Verpol plant in Pittsford, VT.
- The release entered the Dogleg Quarry, an open rock quarry currently used for settling of solids from wash water.
- OMYA and Heindel & Noyes have been investigating the chemistry of the biocide and studying its transport and distribution at the site.
- The released chemical, ortho-phenylphenol (OPP), degrades naturally into carbon dioxide and water. During the breakdown process, intermediate compounds may be formed.
- Water contaminated with OPP from the Dogleg Quarry has circulated to two other quarries used for solids settling and water storage: the Dolomite Quarry and the Pittsford-Italian Quarry (PIQ or East Quarry).
- OPP and possible intermediate breakdown substances have been detected in the Dogleg Quarry and the Dolomite Quarry. OPP alone has been detected in the Pittsford-Italian Quarry.
- Declining OPP levels in all quarries indicate degradation of OPP, and possible formation of temporary byproducts in very low levels.
- Groundwater monitoring has detected no evidence to date of the release impacting groundwater.
- Three private wells and a public community well are being monitored; no impact has been discerned in these wells.
- Additional monitoring is proposed, including installation of monitor wells in locations where no groundwater data exists, and continued analysis of surface water, groundwater, private wells, and the public community well.
- A site wide remediation plan has been developed in the event surface water or groundwater treatment becomes necessary.



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CONTAMINANT SPILL INVESTIGATION

I. BACKGROUND

A. OMYA, Inc. Plant Overview

The OMYA, Inc. Verpol plant in Florence, Vermont produces fine ground calcium carbonate. Quarried marble ore is delivered to the plant in tractor trailer trucks; it is ground into powder and purified. The finished product is shipped off-site as a water-based slurry and as a dry powder in rail cars and trucks. See Site Location Map included in Appendix 1, page 1.

OMYA, Inc. uses water in the purification process. Impurities are floated out of a mixture of water and rock powder, with a flotation reagent. Water from this flotation process and from plant wash water is conveyed through a series of settling ponds and former quarries. Water from these quarries is reclaimed for use in the flotation process, and as the process water that makes up the slurry product.

B. History of Release

On Saturday, November 18, 2000, OMYA's Verpol plant experienced an equipment failure that allowed approximated 4,500 gallons of a calcium carbonate slurry product preservative (P-1 Preservative 20% solution) to enter an on-site wash water quarry (Dogleg Quarry). None of the material or water left the OMYA site via surface water routes.

OMYA personnel immediately investigated and responded to the event. The State of Vermont was notified at about 2:00 a.m. the morning of November 19th via telephone, as soon as the details of the situation were determined. Later that same morning, representatives of the State Agency of Natural Resources and the Town of Pittsford met on site with OMYA representatives to review the situation. OMYA representatives have since been in regular contact with both state and town officials.

Subsequent to the release event, in-plant use of the Dogleg Quarry water was suspended until the night of November 21, 2000. On November 22, Dogleg Quarry water levels had risen such that it was necessary to transfer water into the west settling cell via the flotation launderers. Water balance issues arose as a result of this strategy, and on November 28, the flotation waste (partially consisting of Dogleg Quarry water) was transferred from the west settling cell to the Dolomite Quarry and from there to the Pittsford-Italian Quarry. The switch back from the Dolomite Quarry to the west settling cell occurred on December 4, 2000. Quarry locations are shown on the maps included in Appendix 1.

Plant biocide piping was audited on November 20, 2000 and non-standard piping replaced. Additional change-in-tank-level detection and alarming was added to the plant control system. The biocide storage area sump pump control strategy was modified to incorporate the additional tank level detection. Other system changes designed to enhance environmental protection and safety have been started. Further enhancement of the existing monitor systems is occurring with a project budget request expected to be solidified before the end of December 2000.

In addition to conducting its own investigation, OMYA retained Heindel & Noyes to conduct a hydrogeologic investigation. The Heindel & Noyes (H&N) study included the following:

- Analyze the occurrence and distribution of contaminant release;
- Assess the degradation pathways of contaminant (aerobic and anaerobic conditions);
- Assess the movement of contaminant in the environment (surface water and groundwater);
- Develop a groundwater monitoring plan; and

 Develop a backup remediation strategy, should corrective measures become necessary.

Heindel & Noyes toured the OMYA facility and reviewed engineering drawings to determine the flow path of the contaminant through the plant and surrounding quarries. Water and sediment samples were collected from various quarries and wells on site to determine the extent of contamination resulting from the spill event. Extensive research and laboratory analysis were conducted to determine degradation pathways of the preservative (ortho-phenylphenol, or OPP). Regional surface water and groundwater flow paths were determined to assess the potential movement of the contaminant and to develop a groundwater monitoring program. Finally, a groundwater remediation strategy was developed.

The following sections detail the work conducted and conclusions and recommendations drawn from that work.

II. GEOLOGY AND ENVIRONMENTAL SETTING

A. Location and Topography

The OMYA plant is situated in the community of Florence, within the town of Pittsford, Vermont. The USGS map (page 1 of Appendix 1) shows the site location.

OMYA is located in the foothills at the eastern margin of the Taconic Mountain Range. This region is considered the northern end of the Vermont Valley, which runs north to south between the Green Mountains and the Taconics. The Otter Creek, flowing north, is east of the site. The topography generally slopes east to the Creek, although due to the numerous hills and knolls, slope and topography are highly varied. The OMYA plant itself is located on a promontory of bedrock, surrounded on the west, north, and east by lowlands, swamp, and streams.

B. Geology

Surficial Geology

Surficial materials vary between the upper locations at the plant site, and the valleys on the west, north, and east. At the plant proper, thin layers of glacial till cover bedrock, and at many locations, the bedrock surface is exposed. Drilling logs from water wells in this area indicate till thicknesses from 1 to 3 feet on the higher terrain. In the valleys

surrounding the OMYA site, gravel wells have been drilled into a confined aquifer which is covered by 60 to 70 feet of clay, peat, and silty clay. The confined gravel aquifers are underlain by glacial till.

The glacial till at the OMYA plant and beneath the gravel aquifers would have been deposited during glaciation. As glaciers began to melt, the gravel was deposited by the meltwater which flowed through the valleys. Later, as glacial lakes developed, the valleys were submerged, and the clays and silty clays were deposited. The higher terrain at the OMYA plant proper was above the level of the lakes, and had not received any depositional materials following glaciation. (See Surficial Map, page 2 of appendix 1, and well logs, appendix 2, pages 1-25).

Bedrock Geology

Bedrock is close to the ground surface at the OMYA plant, so structure will determine the flow of contaminants from the spill. Carbonate bedrock with karst features underlies the site, and fractures run sub-parallel with the valley, in a north-northwest to southsoutheast orientation.

Because of the use of the site for bedrock mining, the geology has been well studied. According to reports by Geomapping Associates, the carbonate rocks have solution voids (karst features), which have been filled with clay minerals and sediments. The solution voids are the results of erosion and dissolution of the rock material by groundwater.

Although several bedrock formations are mapped at the site, the rock types in all of these are carbonates: dolomite, limestone, and marble. Beneath the spill location and the plant proper, Ordovician Shelburne limestone has been mapped. This "limestone" has technically been metamorphosed into marble; but the degree of metamorphism is so minor that it resembles the sedimentary rock limestone. A formation of Cambrian-age Clarendon Springs dolomite lies to the east, with Cambrian-age Danby quartzite/dolomite beyond. West of the plant, the Ordovician-age Bascom formation is mapped. This formation is reported to consist of dolomite, marble, and quartzite. (See Bedrock Map and Legend, pages 3-4 of Appendix 1).

To the east of the plant, wells #4, #1, #3, #2, and a second #2 are shown on the map (from north to south). Geomapping reported that these wells penetrate the actual location of the bedrock contact between the Shelburne marble and the Clarendon Springs dolomite. Well #5 lies just east of this contact zone. Along the contact, bedrock

fractures are apparently well developed, as wells #1, #2, and #3 have high yields and interfere with one another.

Generally, the carbonate bedrock can result in high yielding wells and wide-reaching, well-developed fracture networks due to the presence of karst features. However, two of the wells drilled at the site (96-1, 96-1A) have low yields (<5 gpm), suggesting that the fractured zones are not pervasive. The analysis of the existing wells indicates that the fracture zones at bedrock contacts are well developed, while other areas are only slightly fractured. Generalized geologic cross-sections (Centennial Geologic Map of Vermont, 1961) of the area show shallow bedding angles to the rock strata, meaning that fractures should not dip significantly away from the plane of the land surface.

Following the bedrock contacts, major fractures run in a north-northwest to south-southeast direction. We have evaluated two fracture trace analyses: one conducted by Geomapping Associates from unknown photographs, and one performed by Heindel and Noyes using three series of aerial photographs.¹

Fractures in the Geomapping Associates analysis are scattered and show no particular orientation. Because wells drilled at fracture intersections (i.e. 96-1, 96-1A) yielded very low amounts of water, the accuracy of the Geomapping fracture trace cannot be verified. From the Heindel and Noyes fracture trace analysis, the dominant fracture pattern is clearly north-northwest to south-southeast, following the trend of the valley and of the bedrock formations. Some weaker transverse fractures cross the valley.

Although this fracture orientation would be a primary direction for groundwater flow, the topography slopes perpendicular to this orientation. Thus under non-pumping conditions, groundwater flow would be preferentially to the northeast, at a restricted rate due to the lack of fractures trending in that direction. The disparity between fracture orientation and topography creates a highly anisotropic condition.

Due to the use of the plant, the regional groundwater flow trend may have been disrupted. Significant volumes of plant water are recycled through the Dogleg, Dolomite, and Pittsford-Italian Quarries. However, these quarries do not likely discharge or draw water from the aquifer, due to the presence of dozens of feet of impermeable sediment, deposited from settling of waste rock fines. Water overflows from the Dolomite Quarry to the Pittsford-Italian Quarry (PIQ), which has no overland outlet, and is pumped to supply

¹ US Army COE, 1977, Infrared 1:80,000 series. Flown at 40,000 ft. Photos 1-165, 1-164 VT Mapping Program, 1962, VT-62-H 1:18,000 series. Flown at 9,000 ft. Photos 21-92, 21-91 NASA, 1974, VT 7420 1:20,000 series. Flown at 10,000 ft. Photos 7-214, 7-215, 7-216

the plant with process water. As a result, this quarry could act either as a collection basin or a recharge zone for the aquifer, depending on the balance of pumping versus inflow. More monitoring points have been proposed (see Section VI below) to determine the water balance between the quarries and the aquifer.

C. The OMYA Plant

The OMYA plant itself is an important factor in local environmental conditions because of its scale, and uses of rock and water. The orthophoto map (page 5 of appendix 1) shows the plant.

Water Circulation Processes

The two gravel wells (Gravel well #1 and Gravel well #2) and town water are used to supply the plant with water. The gravel wells are located beyond the high ridgeline to the west of the spill site. Because of this geographic separation, and because the wells tap into a gravel aquifer not present at the plant site, these wells are unlikely to be affected by conditions near the spill site.

Water from the Dogleg Quarry, where the spilled OPP collected, is used for the flotation system and for washing plant floors. Floor wash water returns to the Dogleg quarry. Water from the Pittsford-Italian Quarry is pumped to the plant and used for making the slurry product. Some of the water which flows into the Pittsford-Italian Quarry is the overflow of the Dolomite Quarry.

The flotation system separates the heavier carbonate product from the unwanted mica and schist. The ground marble slurry is mixed with a flotation reagent, and air bubbles in a tank. The waste material from the flotation system is sent either to the Dolomite Quarry, or to the two settling cells adjacent to the Dogleg Quarry.

A boom across the Dolomite quarry facilitates settling and separation of floating debris. When the Dolomite Quarry is receiving floation waste, its outflow discharges over a weir, to a gravity culvert/penstock, to the Pittsford-Italian Quarry. When it is not receiving floation waste, its outflow discharges over a weir, following its natural course to a gravity drainage swale, where it mixes with stormwater for eventual flow off-site.

Immediately following the spill, the Dogleg Quarry was not used for flotation water or process water; this water was instead obtained from the gravel wells, the town water, and the PIQ. Use of the Dogleg Quarry water was resumed on November 21, 2000.

There is also a storm water system on site; a settling pond north of the plant receives the water and discharges to a stream. Sanitary wastes are collected in a septic system and discharged to a leachfield located northwest of the main plant. Neither of these systems influence the distribution of the OPP.

Wells

Of the many wells shown on maps, only a few are still serviceable. Due to silty and ochre-colored water, many wells have been abandoned in accordance with the Vermont Water Supply Rule procedures for well abandonment.

Well 96-1A has been filled in with gravel, and capped with bentonite. Wells #2 (from CV104) and #3 are buried and cannot be located. Well #1 has also been filled in. Well #4 cannot be located. Well #2 (the southern #2) exists, but a soft bottom was encountered at 22' while the well is reported as 150 feet deep.² This well was reported to have produced silty, ochre-colored water during production. The silt originated from one of the karst cavities which had filled with clay minerals and sediment; the well may have silted in naturally.

Existing and Abandoned Well Details at OMYA Property

Well	Type	Total	Yield.	Notes
		Depth	gpm	
96-1	Bedrock	500	5	Unused, sampled 12/4/2000
96-1A	Bedrock	300	5,14.51	Abandoned; muddy water
96-2	Bedrock	590	15	Not examined, 2300 feet from Dogleg. Unused.
Gravel #1	gravel	71	20	Used as plant water supply. Drillers yield = 62 gpm
Gravel #2	gravel	82	70	Used as plant water supply. Drillers yield = 124 gpm
#1	bedrock	7	70 2	Abandoned
#2 (CV104)	bedrock	?	?	Abandoned
#2 (S)	bedrock	150 ?	50 ?	Unused, buried. Sampled 12/4/2000. Measured depth of 22' BTC.
#3	bedrock	7	30.7	Abandoned
#4	gravel 🦠	76	75 ?	Abandoned
#5	bedrock	?	30 ?	Supplies a contractor's trailer. Sampled by OMYA.

Shaded wells = abandoned

^{? =} Information unknown or uncertain

² Geomapping Associates, ltd. Verpol Plant Process Water Supply Groundwater Development: Phase I Evaluation. January 7, 1997.

D. Sensitive Environmental Receptors

A survey of sensitive receptors within a 2 ½ mile radius of the spill site was performed. The map on page 6 of Appendix 1 shows mapped Class II wetlands, rivers and streams, and water supply wells. Upon OMYA's initiative, the three nearest private wells are being sampled. Wells serving the Pittsford-Florence Public Community Water System are also being tested.

Bedrock Aquifers

The quarries where the contamination has been detected are excavated into bedrock. Water may discharge from the quarries into the aquifers, depending on water use patterns at the plant. Therefore the potential for contamination of bedrock groundwater exists. At all three quarries, the presence of a thick, impermeable layer of settled mineral fines should inhibit the exfiltration of contaminated water. Permeability testing conducted on sediment cores indicated a very low permeability (see Section IV.3 below). Possible flows from the quarries into the bedrock aquifer could occur where the sediment layer does not extend up the sidewalls to the water surface. Accordingly, no OPP or breakdown products have been found to date in groundwater. Additionally, because of the dilution in the Pittsford-Italian Quarry, contaminant levels within it are fairly low.

Because the possibility of groundwater contamination remains, we are proposing a monitoring program, including the installation of more wells, to verify that the impermeable sediment layers, dilution, and contaminant breakdown are preventing migration of contaminants to the bedrock aquifer. Additional wells will also be used to determine whether groundwater flows into or out from each of the quarries. Additional discussion of the monitoring system is contained in Section VI of the report.

Confined Gravel Aquifers

The confined gravel aquifers are distinct from the bedrock aquifer. A layer of glacial till beneath the gravel is likely to inhibit communication between the two aquifers under natural conditions. However, because both aquifers are confined, it is possible that significant pumping of gravel wells could depressurize the gravel aquifer to the extent that the bedrock aquifer would discharge to it. As the gravel aquifer is located in the low valleys, it is at a lower elevation than the portions of the rock aquifer at the OMYA plant. The proposed monitoring plan will evaluate the potential for this connection (Section VI).

Public Community Water System

The Pittsford-Florence Water System (WSID #5226) operates a pair of gravel wells near the Otter Creek, east of the plant. The main and emergency backup well are located adjacent to each other. According to the water system operator, the wells produce up to 600,000 gallons on peak days. Average production is 200,000 gallons to 350,000 gallons per day. Demand depends upon OMYA's use of municipal water to supplement its own sources.

The wells are approximately 3,300 feet northeast from the Dogleg Quarry, 3,400 feet northeast from the Dolomite Quarry, and 2000 feet east from the Pittsford-Italian Quarry. The wellhead protection area (WHPA) extends to the ridgeline east of the Pittsford-Italian Quarry; none of the three quarries nor the spill site are within the WHPA.

The Public Community Wells are not believed to be at risk from the spill. Because the primary fracture trend runs northwest, contaminant flow is unlikely to move towards the wells. The presence of a ridgeline between the OMYA site and the wells should also direct groundwater flow away from the wells. To date, no OPP has been detected in these wells, or in any other groundwater. Additionally, the wells are located in a gravel aquifer, which is separate from the bedrock aquifer present at the spill area. Because the wells are adjacent to the Otter Creek in a gravel aquifer, recharge water to this aquifer may come from the Creek.

However, because the regional topography slopes from the OMYA site towards these wells, groundwater is likely to flow slowly in this direction. It is also possible that the gravel aquifer could draw water under pumping conditions, from the bedrock aquifer. Therefore, the monitoring program will investigate these wells and the gravel aquifer.

Private Wells

Three private bedrock wells exist in the vicinity of the spill site. These are shown on the sensitive receptor map (page 6 of Appendix 1). The wells are between 2,500 and 2,900 feet from the Dogleg Quarry, at lower elevations. No fracture traces were observed connecting the quarries and the private wells. Sampling to date has found no contamination (see Section IV.6 below).

Mapped Class II Wetlands

Numerous Class II wetlands exist on and off of the OMYA property within a 2 ½ mile radius. Some of these are located northwest from the spill area and the quarries. These wetlands could potentially receive contaminated groundwater discharge from the quarries if such contamination were to enter groundwater. However, no groundwater contamination has been detected to date.

The Pittsford-Italian Quarry is the downstream-most quarry in the series. It has no overland outflow and is instead pumped into the plant for use as process water, or is pumped to downstream settling ponds for offsite discharge in accordance with the OMYA East Plant's NPDES permit. This offsite discharge is conducted infrequently, only when necessitated by spring runoff. Therefore, direct overland flow to wetlands is not possible. To date, only low levels of OPP contamination have been detected in this quarry. Therefore, it is not likely that significantly contaminated groundwater would emanate from the Quarry and move towards the wetlands.

Natural Heritage Areas

From the Vermont Center for Geographic Information files, we have located one site which has been listed as a state natural heritage area. A rare/threatened/or endangered vascular dicot plant species has been identified near the village of Pittsford. Because this location is approximately 2 miles upgradient from the spill site, it is unlikely to be impacted.

Surface waters

The Otter Creek and its tributary streams are present near the OMYA plant. These surface waters could potentially receive contaminated groundwater discharge from the quarries if such contamination were to enter groundwater. However, no groundwater contamination has been detected to date.

Because the contamination has been contaminated in the settling ponds and quarries, and because the Pittsford-Italian Quarry, the downstream-most quarry, does not currently outflow, it is unlikely that contaminated surface waters could flow from the OMYA site to the Otter Creek or its tributaries.

CONTAMINANT CHARACTERIZATION 111.

The industrial process undertaken at OMYA involves the use of P-1 Preservative (20% solution) with the active ingredient ortho-phenylphenol (99.9%) and inert ingredients (0.1%). P-1 is a product of Verichem Inc. of Pittsburgh, Pennsylvania. P-1 is effective against a wide variety of mold fungi and bacteria. P-1 is used for the preservation of polymer emulsions, (coating, PVA systems and rubber), thickeners (starch, cellulose, guar), paper, starches, clay slurries and slurries (See MSDS included in Appendix 3, pages 1-15).

Normally, P-1 preservative is added to the OMYA product as it is loaded into rail tankers; it is not introduced to the water stream anywhere on site. The preservative is used to prevent bacterial growth in the slurry; bacteria cause the slurry to separate and form a concrete-like mass inside the rail cars.

Ortho-Phenylphenol

Physical and chemical properties of ortho-phenylphenol (OPP) include the following:

Appearance and odor:

white flakes; mild, distinct odor

Boiling point:

547° F/ 286° C 134° F/ 56-58° C

Melting point: Specific gravity:

1.26 at 68° F/ 20° C

pH:

6.1 of a saturated solution 2,000 mg/l @ 68°F/ 20°

Solubility in water:

C12H10O

Molecular formula: Reactivity:

oxidizing materials

pH saturated water solution: 12.0-13.5 @25°

Biodegradation:

After 3 weeks of adaptation of 10-40mg/l at 22° C, 100% degradation under aerobic and anaerobic conditions, as sole carbon source or with synthetic

sewage

OPP is soluble in fixed alkali hydroxide solutions and most organic solvents, and is nearly insoluble in water.3 OPP fact sheets are included in Appendix 3, pages 16-28. Diagrams of the molecular structure of OPP and breakdown products are provided on page 29 of appendix 3.

³ Windholz et al, 1983.

OPP Breakdown Products

Given the highly reactive characteristics of OPP, much attention has been given to the potential breakdown pathways of OPP and potential intermediate degradation. Biodegradation depends on many factors including physical and chemical properties, concentrations, diffusion rates, bioavailability, nutrient availability, other chemicals in the matrix, pH, temperature and other seasonal effects.⁴

There are a number of naturally occurring microorganisms capable of degrading OPP through a myriad of potential pathways and intermediates. Generally aerobic breakdown of phenolic compounds results in the production of catechols which are easily degraded via mineralization to carbon dioxide and water. The most likely intermediate breakdown products include oxygenated hydrocarbons (carboxylic acids, ethers, esters, ketones and etc.) of 12 carbons and less. The breakdown pathways are facilitated by the bacterial enzymes used to transfer electrons from the contaminant to an "electron acceptor". The most common electron acceptors are oxygen, and dissolved metals species such as iron and manganese. Dissolved nitrate, sulfates and carbon dioxide can also act as electron acceptors.

H&N's research also indicates that anaerobic degradation of phenolic compounds is common. Anaerobic degradation depends on the presence of an electron donor for enzymes to facilitate reduction of the contaminant. The presence of calcium carbonate in the process waters provides a direct source of electron donor via the negatively charged carbonate ion.

All of the degradation pathways evaluated for phenolic compounds indicate complete degradation of the parent compound without the creation of significant hazardous end products. All of the intermediate compounds are relatively unstable and quickly degraded. Due to rapid decay, the intermediate compounds are unlikely to exist in significant concentrations.

Estimated degradation half life's for OPP are presented in the Handbook of Environmental Degradation Rates⁵ for unacclimated aerobic and anaerobic aqueous systems ranging from 1 to 28 days. These rapid degradation rates are based on field data acquired from a spill into a river. Research on OPP and metabolic products

⁴ Sutherland et al., 1995.

⁵ Howard et al, Lewis Pub., 1991.

continues and will be used to specify the test methods to detect both aerobic and anaerobic metabolics.

IV. DATA COLLECTON

A. Sampling Methods

OMYA commenced daily surface water monitoring of the Dogleg, Dolomite and Pittsford Italian Quarries on November 18. Groundwater monitoring of the on site "CDP" bedrock well (well #5 on map), the gravel well serving the Pittsford-Florence Water System, and three nearest residential wells commenced on November 20. Samples were submitted to Verichem, Inc. for analysis of OPP and pH. Data summary sheets compiled by OMYA are included in Appendix 4, pages 53-55.

H&N also collected surface water, sediment and groundwater samples at the OMYA facility. On November 21, surface water and sediment samples were collected from the Dogleg quarry. On December 4, water samples were collected from the Dolomite Quarry and on site groundwater wells 96-1 and well #2. Samples were submitted to Endyne, Inc. in Williston, Vermont. Individual laboratory data sheets are included in Appendix 4, pages 1-52.

Surface Water Sampling

Surface water samples were collected using a disposable bailer. Samples were preserved with ice prior to delivery to Endyne, Inc..

Groundwater Sampling

Bedrock wells were selected for on-site groundwater sampling. Wells were examined which were located downgradient from, and within 1,000 feet of the Dogleg Quarry. Well #5 ("CDP") is currently in use, and was sampled at the tap. Wells #2 (from CV104), #3, #1, and 96-1A have been abandoned (filled in), and could not be sampled although they were in the area of interest. Wells 96-1 and #2 (the southern well #2) were unused, and were sampled by purging with a portable sampling pump. Sample locations are presented on the Orthophoto Map in Appendix 1, page 5.

Test well 96-1 was sampled with a Redi-Flow pump set at a depth of 300 feet, which is the maximum length of power cable for the pump. The well was purged for 50 minutes, until temperature and specific conductance readings stabilized. This well is a modern bedrock well with an 8-inch casing and a 6-inch borehole, and is located by the

500,000-gallon water tank, at a site which is laterally downgradient from the Dogleg Quarry.

Well #2 (the southern of the two well's #2) was also sampled with the Redi-Flow pump. This well had been buried in a truck trailer drop area, and was reportedly a 150 foot deep bedrock well. It is located between the Dogleg Quarry and the Dolomite Quarry. A soft bottom was encountered in the well at only 22 feet below the top of the casing, and tan-brown sand particles were entrained in the initial purge water. Because the well was able to produce over 1 gpm with virtually no drawdown, it is likely that water from the bedrock fractures was able to flow up through the sand fill. A sample was collected after 55 minutes, when temperature and conductivity had stabilized. Water from this well had a faint odor of peat or muck.

A previous consultant's report⁶ indicated that well #2 produced silty, ochre-colored water during production. This water originated from one of the karst cavities which had filled with clay minerals and sediment. We suspect that the well may have silted in naturally because of this cavity.

Off-site groundwater sampling has been performed at the Pittsford-Florence gravel well, and at the three residential wells identified in the sensitive receptor survey. These wells have been sampled at the taps, using the existing pumps in the wells.

Sediment Coring

Sediment core samples were collected from four locations in the Dogleg quarry. Sample locations are indicated on the Site Investigation Map in Appendix 1, page 7. A clear acetate liner/tube (3ft long and 1.75" diameter) was connected to PVC pipes that were advanced by hand into the quarry sediment. Access to the sample points was by rowboat. The acetate liners were plugged and transported back to H&N's office in Burlington where samples were extracted from the tubes for laboratory analysis. Sample collection depths are presented in the following table.

Boring Location	Depth (below water surface)
SS#1	15-18'
SS#3	6,
SS#4	25-28'
SS#5	12-15'

⁶ Geomapping Associates, LTD. Verpol Plant Process Water Supply Groundwater Source Development: Phase I Evaluation. January 7, 1997.

B. Laboratory Methods

Samples have been analyzed by Verichem and by Endyne laboratories.

Endyne, Inc. of Williston, Vermont, an EPA and VTDEC certified analytical laboratory, analyzed all samples collected by H&N. In order to reasonably document orthophenylphenol (OPP) concentrations and potential OPP degradation products, several laboratory test methods were employed.

EPA Method 642 (an HPLC method) is applicable to the determination of OPP in municipal and industrial discharges. Ortho-phenylphenol is extracted at neutral pH into an organic solvent. After extraction, quantitation is performed by high performance liquid chromatography (HPLC) with an ultra-violet (UV) detector.

The OPP analytical method via EPA Method 642 has inherent limitations due to the specified neutral pH extraction protocols and the generally poor peak resolution. High calcium carbonate and emulsifier/flocculent concentrations in the sample matrix tend to interfere with the extraction efficiency and the ability for the instrumentation to adequately resolve individual compounds. Initial analytical results tentatively identified both phenol and OPP in the HPLC analysis of Dogleg Quarry water. See December 4, 2000 Memo from Endyne, Inc. included in Appendix 4, page 1.

Confirmation of the EPA Method 642 results was performed at Endyne, by using gas chromatography/mass spectroscopy techniques (EPA Method 8270C). EPA Method 8270C for semi-volatile organic compounds (SVOCs) includes OPP and phenol as target compounds. Method 8270C specifies both acidic and basic pH extraction with GC/MS analysis using mass spectral fingerprinting of resolved compound peaks. Gas chromatography/mass spectroscopy techniques (GC/MS) SVOC analysis confirmed the OPP in one of the samples (SS#1) but did not confirm the phenol tentatively identified in the HPLC analysis of the same sample. Based on this finding, identifications of phenol and other degradation products via EPA Method 642 were changed to non detect (see Endyne Memo dated 12/07/00 in Appendix 4, page 3-4). Therefore, EPA Method 8270C is the preferred method for detection of phenol and other potential degradation products in this sample matrix.

A final laboratory test method (EPA Method 8260) was employed for the volatile fraction of the potential OPP breakdown products. Since the microbial breakdowns of OPP into smaller potential intermediate compounds are likely, analysis of the more

volatile fraction of organic compounds was necessary to determine breakdown product characterization and concentration. EPA Method 8260 for volatile organic compounds (VOCs) specifies using a purge and trap sample concentration system and a GC/MS separation and detection system to measure VOCs in various matrices. The mass spectral detector can also estimate the identity non-target compounds ("unidentified peaks") by their spectral fingerprint based on a large compound database.

C. Laboratory Results

Dogleg Quarry: Surface Water

OMYA commenced surface water sampling at the north and south ends of the Dogleg Quarry on November 18, 2000. Samples were submitted to Verichem, Inc. for analysis of OPP and pH. The laboratory results are summarized in the following table.

Dogleg Quarry: Surface Water Sampling (Verichem)

Sample Date	OPP (ppm) Dogleg Quarry South End (Plant Discharge End)	pΗ	OPP (ppm) Dogleg Quarry North End (Plant intake End)	pH	Laboratory
44/40/00	650	9.81	950	10.04	Verichem
11/18/00	550	9.64	950	10.06	Verichem
11/19/00	108	7.76	391	9.14	Verichem
11/20/00		9.59	518	9.59	Verichem
11/21/00	360	7.77	167	9,27	Verichem
11/22/00	187	7.59	153	7.44	Verichem
11/23/00	100	7.36	145	7,36	Verichem
11/24/00	55	7.54	47	7.43	Verichem
11/25/00	123	7.52	57	7.53	Verichem
11/26/00	86	7.35	37	7.51	Verichem
11/27/00	19	7.42	51	7.9	Verichem
11/28/00	40		26	7.55	Verichem
11/29/00	<0.040	7.43	25	7.33	Verichem
11/30/00	11	7.45	14	7.31	Verichem
12/1/00	<0.040	7.53	7		Verichem
12/4/00	<0.040		<u> </u>		Endyne
	-		0.283		

Water samples were also collected for analysis of aerobic bacteria and pH. Laboratory results are summarized in the following table.

Bacteria and pH Data: Dogleg Quarry (Verichem)

Date		g Quarry ntake End		Quarry Discharge End
	ρН	Aerobic Bacteria	рH	Aerobic Bacteria
		(cfu/ml)		(cfu/ml)
11/20/00	8.22	510000	9.29	12000
11/21/00	8.09	2500000	9.73	900
11/22/00	8.19	40000	9.12	120000
11/23/00	8.03	80000	8.08	520000
11/24/00	7.97	370000	7.94	700000
11/25/00	7.78	400000	8.05	**
11/26/00	7.58	>1000000	7.86	800000
11/27/00	7.53	*	7.66	***
11/28/00	7.85	250000	7.82	***
11/29/00	7.39	650000	7.55	380000
11/30/00	7.91	800000	7.65	860000
12/1/00	7.79	8000000	7.67	1250000
12/2/00	7.73	4000000	7.67	2000000
12/3/00	7.54	1500000	7.5	600000

Notes: *= Data Questionable (10⁴ – 10⁶) **= Data Questionable (10⁵-10⁶)

SS#1: Surface Water

		OOM 1: Culture Trace.	
Sample Date	Test Method	Compound	Concentration (ppm)
11/21/00	EPA 642	OPP	97
11/21/00	EPA 8260	Acetone*	0.950
11/21/00	EPA 8270C	Benzoic Acid*	0.015 0.005
		1-Phenyl ethanone* 2-Methoxy-1, 1-Biphenyl*	0.003
		1-Methyl-4-1-(1-methylethenyl)- cyclohexanol*	0.002

^{*}Substance was detected as an Unidentified Peak in spectral analysis. Identification and quantitation of substance are based on a best match to library data and are not confirmed.

^{***=} Data Questionable (10⁵-10⁶)
****= Data Questionable (10⁵-10⁶)

Surface water samples collected at SS#1 and SS#2 on November 21, 2000 were submitted for laboratory analysis via EPA Method 8260 for VOCs, EPA Method 642 for OPP, and EPA Method 8270C for semi-volatile organic compounds. Compounds which have been detected via these analysis are presented in the following table. (See appendix 3 for information of the properties and structure of these substances.)

SS#2: Surface Water

Concentration (ppm)	7.0
	7.37
Tact Method Compound	<u> </u>
Test Method Compound Concentration (ppin)	. !
FDA 642 OPP 3.32	
EPA 642 OPP	_

Dogleg Quarry: Sediment Coring

Water separated from the top of the sediment core collected at SS#1 (15-18' below water surface) was submitted to Endyne, Inc. for analysis via EPA Method 8260 for VOCs, and EPA Method 642 for OPP. Compounds detected via these analysis are presented in the following table.

SS#1 (15-18' below water surface): Sediment Core Water

20#1(12-14	Delow water se	211400/: 00 ann	
Sample Date	Test Method	Compound	Detection (ppm)
11/21/00	EPA 642	OPP	168
11/21/00	EPA 8260	Acetone*	0.550
		Toluene	0.0023

^{*}Substance was detected as an Unidentified Peak in spectral analysis. Identification and quantitation of substance are based on a best match to library data and are not confirmed.

The top 4 inches (0-4") of the solid phase of the sediment core on SS#1 (15-18' below water surface) was sampled and analyzed for VOCs and OPP. This sample was approximately 56% solids. Compounds detected via these analysis are presented in the following table.

SS#1 (15-18' below water surface) 0-4" Sediment

	33#1(1	0-10 Delott trace	7, 00,100-7	
	Sample Date	Test Method	Compound	Detection
				(ppm)
			OPP	42.3
	11/21/00	EPA 642	1 011	0.470
-	11/21/00	EPA 8260	Acetone*	0.470
	11121100			

^{*}Substance was detected as an Unidentified Peak in spectral analysis. Identification and quantitation of substance are based on a best match to library data and are not confirmed.

Sediment Permeability Test (K Test): Dogleg Quarry

The permeability of a sediment core from the Dogleg quarry was tested with a constant-head permeability experiment. This test was conducted to assess the rate at which contaminated water could discharge through the quarry walls and floor into the rock aquifer. The permeability ranges from 0.01 to 0.08 ft/day, which is within the low end of the range reported for silt in hydrology texts (Freeze and Cherry, 1979). Calculation briefs are included in Appendix 4, pages 56-57.

The experiment was conducted on a core sample of sediment obtained from a depth of 25-28 feet. The sediment was compact and had a distinct odor. The core was contained

in a 3-foot long, 1.75" diameter (4.267 cm) clear acetate liner/tube. The tube was fitted with a filter fabric sleeve to allow water to permeate through the base of the column, while retaining the sediment. This soil column was vertically oriented in a constant-head basin. Water was then added to the core tube, and the level kept at a mark near the top, thus creating the constant head condition. Because the core was already saturated, the flow rate could be immediately measured at the outflow from the constant head basin. A schematic of the test device is shown in Appendix 4, page 56.

During the 22-hour test, water levels in the core tube, and discharge volumes in the beaker were checked periodically. Discharge volumes were accurately measured with a graduated cylinder, and an equal volume was added to the core tube to maintain constant head conditions.

Permeability was calculated using Darcy's law, rearranged to solve for permeability (K) when cross-sectional area (A), gradient (dH/dL), and flow (Q) are known. See the calculation sheet (Appendix 4, pages 56-57).

The flow rate through the sediment sample was not steady during the entire test; at its fastest, 3.5 mL were collected in about two hours, while over the following 15 hours, only 0.1 mL flowed through the core. Small reductions in the outflow rate are most likely related to gas formation (CO₂) in the soil column. Therefore, two analyses of the results were performed:

- 1) The maximum permeability was calculated from the fastest measured flow rate. This result was 0.080 ft/day.
- 2) The average permeability was calculated from the average flow rate over the entire test. This result was 0.0112 ft/day.

4. <u>Dolomite Quarry: Surface Water</u>

OMYA commenced surface water sampling of the Dolomite Quarry on November 18, 2000. Samples have been analyzed for OPP and pH by Verichem, Inc. Samples collected by Heindel and Noyes on December 4, 2000 were submitted to Endyne Inc. for laboratory analysis via EPA Methods 642, 8260 and 8270C. Laboratory results are summarized in the following table. Individual laboratory reports are included in Appendix 4, pages 1-52.

Dolomite Quarry: Surface Water

Sample	Test	Laboratory	Compound	Detection
Date 💮	Method			(ppm)
11/18/00	377-15-15-15-15-15-15-15-15-15-15-15-15-15-	Verichem	OPP	<0.040
			pH	8.02
11/19/00		Verichem	OPP	<0.040
			pH	7.94
11/20/00		Verichem	OPP	<0.040
			pH	7.77
11/21/00		Verichem	OPP	0.067
11/23/00		Verichem	OPP	<0.040
11/24/00		Verichem	OPP	<0.040
11121100			pH	7.54
11/27/00	<u> </u>	Verichem	OPP	<0.040
11,21,00			Hq	7.8
11/28/00	 	Verichem	OPP	<0.040
11720100			pH	7.37
11/29/00		Verichem	OPP	<0.040
1723700		10110112111	pH	7.8
11/30/00	<u> </u>	Verichem	OPP	10
11/00/00			PH	7.86
12/1/00	 	Verichem	OPP	<0.040
12.7700			pH	7.59
12/4/00		Verichem	OPP	<0.040
12/4/00	EPA 642	Endyne, Inc.	OPP	4.75
12/4/00	EPA 8270C	Endyne, Inc.	OPP	4.99
124700			Phenol	TBQ <0.005
			Crotonic Acid*	0.002
			2-Heptanone*	0.002
			2-Piperidinone*	0.003
			1-Phenyl ethanone*	0.001
			2-Dibenzofuranoi*	0.002
			3-Amino Phenol*	0.005
			Unknown*	0.010
12/4/00	EPA 8260	Endyne, Inc.	Chloroform	0.0019
100	/		Acetone*	0.230
			2-Butanone*	0.033
			Toluene*	TBQ <0.0009
12/6/00	EPA 642	Endyne, Inc.	OPP	3.57
12/12/00	EPA 8720C	Endyne, Inc.	OPP	5.36

Laboratory results from the 8270C analysis of samples collected on December 4 indicated a concentration of 4.99 ppm OPP. In addition, the EPA 8270C analysis indicated a trace below quantitation (TBQ) concentration (<5ppb) phenol and seven unidentified peaks. The unidentified peaks were tentatively characterized based on the best match of the sample peak and library spectral data. The seven peaks best match were crotonic acid, 2- heptanone, 2-piperidinone, 1-phenyl ethanone, 2-dibenzofuranol, 3-amino phenol, and an unknown with no reliable spectral match. Information on the structure and properties of these substances is contained in appendix 3.

Surface water samples from the Dolomite Quarry were also analyzed for aerobic bacteria levels. Laboratory results are summarized in the following table.

Dolomite Quarry Discharge: Bacteria and pH Data (Verichem)

Date	pH /	Aerobic Bacteria
		(cfu/ml)
11/20/00	8.13	720000
11/21/00	8.07	600000
11/22/00	8.01	630000
11/23/00	7.95	960000
11/24/00	7.91	250000
11/25/00	8.01	300000
11/26/00	7.9	270000
11/27/00	7.8	90000
11/28/00	7.95	160000
11/29/00	8.02	190000
11/30/00	7.96	90000
12/1/00	7.94	140000
12/2/00	7.95	330000
12/3/00	7.69	390000

5. Pittsford-Italian Quarry

OMYA commenced surface water sampling of the Pittsford-Italian Quarry on November 18, 2000. Samples have been analyzed for OPP and pH. Laboratory results are summarized in the following table.

Date Sampled	OPP	pH -	Laboratory
	(p pm)		
11/18/00	<0.040	7.92	Verichem
11/19/00	<0.040	8.0	Verichem
11/20/00	<0.040	7.84	Verichem
11/21/00	< 0.040	7.8	Verichem
11/22/00	<0.040	7.74	Verichem
11/23/00	<0.040	7.79	Verichem
11/23/00	< 0.040	7.9	Verichem
11/24/00	<0.040	7.71	Verichem
11/25/00	<0.040	7.78	Verichem
11/27/00	<0.040	7.86	Verichem
11/28/00	<0.040	7.84	Verichem
11/29/00	< 0.040	7.8	Verichem
11/30/00	< 0.040	7.86	Verichem
12/1/00	< 0.040	7.82	Verichem
12/4/00	<0.040		Verichem
12/6/00	1.96		Endyne
12/12/00	2.62		Endyne

6. Groundwater Sampling Results

OMYA commenced groundwater monitoring of well #5 (the CDP well), the Pittsford-Florence gravel well, and neighboring residential wells on November 20, 2000. Samples were submitted to Verichem, Inc. for analysis of OPP and pH. Laboratory results are presented in the following summary table.

Sample Date	Location	OPP Result	pН	Laboratory
		(ppm)		
11/20/00	Well #5	<0.040		Verichem
11/21/00	Well #5	<0.040		Verichem
11/21/00	Town	<0.040		Verichem
11/22/00	1261 W. Creek Rd	<0.040	7.44	Verichem
11/22/00	7797 Whipple Hollow Rd	<0.040	7.56	Verichem
11/22/00	1377 W. Creek Rd	<0.040	7.61	Verichem
11/22/00	Well #5	<0.040	7.36	Verichem
11/22/00	Town	<0.040	7.51	Verichem
11/23/00	Town	<0.040	7.6	Verichem
11/24/00	Town	<0.040	8.26	Verichem
11/25/00	Town	< 0.040	7.49	Verichem
11/26/00	Town	<0.040	7.57	Verichem
11/27/00	Well #5	<0.040	7.43	Verichem
11/29/00	Well #5	<0.040	7.38	Verichem
11/30/00	Well #5	<0.040	7.38	Verichem
11/30/00	Town	<0.040	7.55	Verichem
12/1/00	Well #5	< 0.040	7.34	Verichem
	Town	<0.040	7.57	Verichem
12/1/00 12/6/00	Well #5	<0.002		Endyne

Note: The lowest concentration detectable by Verichem, Inc. is 40 ppb.

Groundwater samples collected by H&N from on site wells 96-1 and well #2 were submitted to Endyne, Inc. for laboratory analysis via EPA Methods 8260, 8270C and 642.

Groundwater collected from well 96-1 on December 4, 2000 contained no VOCs, no OPP, and no SVOCs as analyzed by EPA methods 8260, 8270C, and 642 testing. No unidentified peaks were detected by any analysis.

Well 96-1

		***	,,, ,,, ,	
!	Sample Date	Test Method	Compound	Detection (ppm)
	12/4/00	EPA 642	OPP	< 0.001
	12/4/00	EPA 8270C	OPP	<0.005

Groundwater collected from well #2 on December 4, 2000 contained no OPP and no target or non-target compounds by EPA 8260. However, two non-target compounds were detected during the 8270C analysis that were characterized as dimethyl-1,4-dioxane (about 15 ppb) and 2-methyl-3-buten-2-ol (about 5 ppb). These compounds are not believed to be OPP breakdown products, but may have originated in petroleum compounds from the historical industrial activity in the area where the well is located.

Well#2

Sample Date	Test Method	Compound	Detection (ppm)
12/4/00	EPA 642	OPP	<0.001
12/4/00	EPA 8270C	OPP	<0.005
		Dimethyl-1,4-Dioxane*	0.015
		2-Methyl-3-buten-2-oi*	0.005

^{*}Substance was detected as an Unidentified Peak in spectral analysis. Identification and quantitation of substance are based on a best match to library data and are not confirmed.

V. CONCEPTUAL MODEL

OPP is present in all of the lagoons and quarries at the OMYA site. It is degrading rapidly, through aerobic and anaerobic processes. The end products of degradation are nonhazardous, although intermediate products may be formed, including some hazardous organic chemicals. Such intermediate chemicals would be very short-lived, however, and exist only in very low concentrations. Migration of OPP and breakdown products into groundwater and surface waters is unlikely, except possibly in low levels close to the quarries.

A. Contaminant Distribution

The OPP spill entered the Dogleg Quarry, and high levels of 500 to 650 ppm were initially observed. Because OPP is poorly soluble in water with a pH of less than 12, a portion of the chemical partitioned onto the sediment covering the quarry bottom. Levels detected in the sediment core sample indicate that the concentrations decreased quickly with depth into the sediment layer. In the core sample at SS-1, taken from near the sediment surface (15 feet below the water surface), the OPP concentration of 42.3 ppm contrasts with the 97 to 168 ppm concentration in quarry water at the same location.

When it became necessary to discharge flotation water to the Dolomite Quarry, OPP was introduced to it. OPP levels were first detected there at 10.0 ppm, but declined due to degradation and dilution with inflowing water after the discharge ended, to under 5

ppm. This water discharges through a culvert/penstock to the Pittsford Italian Quarry (PIQ). Following the cessation of discharge to the Dolomite Quarry, OPP was present in lower levels near 2 ppm in the PIQ.

Groundwater sampling at wells 96-1, #2, and #5 indicates that the bedrock aquifer in these regions contains no OPP or breakdown products. The well #5 data, collected from a point downgradient from both the Dogleg and Dolomite Quarries, shows no OPP, including results from an analysis performed by Endyne. It is unlikely that the bedrock aquifer downgradient from the Dogleg, Dolomite or Pittsford-Italian Quarries would be contaminated, because all of these quarries contain significant depths of relatively impermeable sediment. Significant exfiltration into the aquifer could only occur if portions of the quarry sidewalls are above the sediment layer.

B. Contaminant Transport

Contaminants are being transported from the Dogleg Quarry to the Dolomite Quarry to the Pittsford-Italian Quarry and back to the Dogleg Quarry via OMYA's surface water circulation system.

The permeability test conducted on the sediment core from the Dogleg Quarry indicates a low permeability, equivalent to the low end of the range associated with silts. This indicates that pure water could migrate through the sediment at a rate of 0.023 ft/day, assuming a porosity of 0.35 and a gradient of 10% as reflected by surface topography. At this rate, it would take over two years for pure water to travel through the sediment layer, with an estimated thickness of 20 feet.

Travel of dissolved contaminants would be retarded because of OPP's poor solubility and greater affinity for carbon-based material. The factors needed to quantify retardation are unknown, but would likely increase the travel time for OPP and breakdown products to well over two years. With a half-life ranging from 1 to 28 days, concentrations would be reduced by a factor of at least 2×10^9 before leaving the system. Similar conditions are believed to exist in the Pittsford-Italian and Dolomite Quarries. Depth soundings from the Dolomite Quarry show a 20 to 30 foot depth of sediment (Appendix 1, page 8), which is the same material as found in the Dogleg.

Contaminated water from the Dogleg Quarry enters the Dolomite Quarry after it is first pumped into the flotation system, and then flotation wastewater is sent to the Dolomite Quarry. This only occurs when the Dogleg Quarry becomes full, and wastewater must be directed elsewhere. From the Dolomite Quarry, water overflows at a weir into a

culvert/penstock which leads to the Pittsford-Italian Quarry. After mixing with the clean water in the PIQ, levels of 1-2.6 ppm have resulted. This water is then pumped back to the plant for use in process water.

Whether flow out of the quarries occurs is uncertain, and will be verified via further groundwater sampling at well 96-2 and a proposed well. The ability of water to enter the aquifer depends in part on the water balances in the quarries, piezometric conditions of the rock aquifer, the extent to which flotation waste sediment covers the sidewalls, and whether OMYA pumps more out than flows in.

Insufficient data currently exists to understand the flow regimes between the quarries and the aquifers. Whether flow is from a quarry to the adjacent aquifer, or vice-versa can be determined from the water elevations. The table below summarizes the available information.

Flevation*	ws. Depth, Ft	ws Elevation, Ft
	47.75	495.25
	6.38	513.62
l		511
		501
471		471
	543 520 511 501	520 6.38 511 501

^{*} Elevations are from Vermont Center for Geographic Information, Digital Elevation Model, and are not field verified or surveyed.

From this elevation data, well 96-1 can be considered to be downgradient of the Dolomite and Dogleg Quarries, and well #2 is higher in elevation than either. Site observations suggest that the Digital Elevation data may not be accurate, as the Dogleg Quarry appeared to be higher in elevation than both well #2 and well 96-1. The Digital Elevation data suggest that groundwater discharges from the Dogleg, to the aquifer near 96-1, and that the Dolomite receives water from the aquifer near well #2. More monitoring points, and surveyed elevations, around the three quarries are needed to make a conclusion regarding the water balances. Such wells are proposed (see section VI below).

In the unlikely event that the bedrock aquifer were to be contaminated, it is possible for the gravel aquifers to become contaminated. Because the OMYA gravel wells #1 and #2, and the Pittsford-Florence Water System gravel well are consistently pumped at high rates, they could induce the flow of water from the bedrock aquifer to the confined gravel layer. These wells are quite distant from the OMYA site, so the possibility of contamination is remote.

The town well is adjacent to the Otter Creek, which is likely to be a significant source of recharge water to the gravel aquifer. Therefore, pumping the town well is unlikely to influence flow from the bedrock aquifer.

C. Contaminant Degradation

Laboratory results from the Dogleg Quarry and the Dolomite Quarry indicate steadily declining levels of OPP. Low concentrations of possible breakdown products indicate that intermediate products may be forming, but they are not persistent. Research indicates that OPP probably degrades into numerous intermediate products – oxygenated hydrocarbons – which are ultimately broken down to carbon dioxide and water.

In the north end of the Dogleg Quarry, OPP levels declined from 950 ppm to 51 ppm in 10 days before water was transferred to the Dolomite Quarry. Dilution with fresh flotation wastewater accounted for only a small fraction of this reduction. Natural degradation resulted in the majority of the decrease in OPP levels. With high levels of aerobic bacteria in the quarry, and available electron acceptors in the form of oxygen dissolving from the atmosphere, biodegradation could occur rapidly. Further degradation has occurred, so that less than a month after the release, OPP levels in the Dogleg Quarry have decayed to less than 0.3 ppm.

Similarly, in the Dolomite Quarry, OPP levels declined from a high of 10 ppm following the first discharge of contaminated water to that quarry, to 3.57 ppm six days later.

Potential breakdown products are present in only very low levels, indicating their instability. The most prevalent likely breakdown product is acetone, which has been tentatively detected at levels ranging from 0.23 ppm to 0.95 ppm. Compared to the OPP levels, the acetone concentrations clearly show that breakdown products are not accumulating. Other possible degradation products: phenols, ketones, and catechols, have been present at concentrations of less than 0.05 ppm, despite the presence of the parent OPP at over 900 ppm.

If OPP and breakdown products were to migrate into the groundwater, the degradation reactions would continue. Groundwater contains moderate levels of dissolved oxygen which could support aerobic degradation. Other electron acceptors should also be readily available, including iron and manganese which are probably present in high concentrations, given the description of the bedrock water as ochre-colored. In the

event that anoxic conditions exist in groundwater or in the quarry sediments, anaerobic degradation would occur, with the abundant calcium carbonate behaving as the electron donor required for such a reaction.

D. Conceptual Model Summary

OPP is distributed in the lagoons and quarries at the site, and travels primarily through OMYA's water circulation system. It is degrading rapidly, through aerobic and anaerobic processes. Intermediate breakdown products are short-lived, and may exist only in very low concentrations. Water and carbon dioxide are the ultimate end products. Migration of OPP and breakdown products into groundwater and surface waters is minimized by the impermeable sediments in the Dogleg, Dolomite and Pittsford-Italian Quarries. Further testing and well installation is needed to verify flow patterns between the quarries and the aquifer.

VI. MONITORING PROGRAM

It is necessary to further characterize the degree and extent of contamination, the direction of contaminant flow, and the chemical breakdown processes which are occurring, including the formation of daughter products. The proposed monitoring program involves sampling of surface water from the open-water quarries and groundwater from bedrock wells, private wells, and the town well. A test will be performed to determine if the bedrock aquifer could communicate with the gravel aquifer, and sampling of gravel wells will occur if a communication exists. Monitoring will be conducted periodically until it can be assured that the contamination has attenuated.

A. Groundwater Monitoring

A groundwater monitoring program is necessary to investigate whether the bedrock and gravel aquifers have been, or will be, contaminated with OPP or its breakdown products. Bedrock aquifer contamination could occur if contaminated water were to exfiltrate from the settling ponds, the Dogleg Quarry, the Dolomite Quarry, or the Pittsford-Italian Quarry. To date, wells 96-1, #2, and #5 have been sampled with no detections of OPP or breakdown products. For gravel aquifer contamination to occur, the bedrock aquifer would first need to become contaminated. Then, pumping of gravel wells would need to occur at a high enough rate to depressurize the gravel aquifer, allowing water to flow from the bedrock aquifer.

There are six wells at the OMYA property that have not been filled in or abandoned. Bedrock wells 96-1, #2, #5, and 96-2 are useful for groundwater sampling. However, they do not provide a complete picture of the aquifer. Additional wells are needed to determine whether certain areas downgradient from the quarries have been contaminated, and to ascertain the direction of flow between the aquifer and the three quarries.

1. Current Well Network

Well 96-1, with a depth of 500 feet and a 5 gpm drillers yield, is unused. It can be sampled with a portable pump, with sufficient purging to acquire a valid sample. The well drillers log indicates that a single water-bearing fracture was encountered, at the depth of 420 feet. The well is located near to a H&N fracture trace which runs to the Dogleg Quarry. Because of the low yield, it is doubtful that the well intersected the fracture. This well is located laterally downgradient from the Dogleg Quarry. Therefore, water from the well could manifest effects to the aquifer from the Dogleg Quarry. However, because well 96-1 is only laterally downgradient, and because the sole fracture is 420 feet deep, it is probable that aquifer contamination could exist, but not be detectable at this well.

Well #2 (the southern of the two wells #2) is also unused. Like well 96-1, it may be sampled with a field pump. No drillers log exists; however the Geomapping report describes a 150 foot total depth and a 50 gpm yield. The well is reported to tap into a solution cavity which produced silty water, hence it is now silted in to a depth of only 22 feet. The well is topographically downgradient from the Dogleg Quarry, but with a higher water elevation, and adjacent to the Dolomite Quarry, and thus could manifest aquifer contamination from either of these sources.

Well #5 probably provides the most useful data of the existing wells. As it is used to supply a contractors trailer-office, the well is pumped regularly and thus would be likely to draw in contamination. Sampling is conducted at the tap. The well is located topographically downgradient from both the Dogleg and Dolomite Quarries; it is aligned in the dominant fracture direction with the Dolomite Quarry.

Well 96-2, currently unused, could also be sampled with a field pump. This well would indicate the effects of the Pittsford-Italian Quarry, since it is located topographically downgradient from that quarry, along a fracture trace which represents the contact between the Shelburne Limestone and the Clarendon Springs Dolomite. Measurement of water levels in this well will help to determine the water balance at the PIQ.

Gravel wells #1 and #2, located in the valley west of the plant, could be sampled at the taps if evidence indicates that the gravel aquifer is at risk of contamination. They will be used to test for communication with the bedrock aquifer, by being pumped, while water levels in the unused bedrock wells are continuously recorded.

2. Gaps in the Current Well Network and Proposed Wells

The current well network does not test the bedrock aquifer at locations directly downgradient from the Dogleg Quarry, nor in the direction of the PCWS wells. No wells exist on the east side of the Dolomite Quarry. Insufficient wells exist surrounding the three quarries to determine whether they are gaining or losing water to the aquifer.

Three bedrock wells are proposed. The first, located downhill of the Dogleg Quarry, is situated on a strong fracture trace leading to the Dogleg. A well at this site would best show the water quality and flow effects of the Dogleg Quarry on the aquifer.

A second well, east of the Dolomite Quarry, would check for exfiltration into the aquifer on that side of the quarry, where no monitoring points currently exist. This will provide key data on groundwater flow into or out of that quarry.

A third well site is shown south of the Pittsford-Italian Quarry, on the east-west trending fracture traces. This well would serve as an early warning system for the Public Community Wells, as it is located between them and the spill area. Based on the general west to east regional groundwater flow, this well site is probably downgradient from the quarries. Depth measurements at this well would indicate the flow relationship between the PIQ and the aquifer. Proposed bedrock monitor well locations are shown on the Orthophoto Map in Appendix 1, page 5.

3. Monitoring Program

We propose to install the three monitoring wells using an air-rotary drill rig to a depth of 300 feet. We believe that deeper fractures could not be affected by the contamination, due to the shallow bedding angle of the bedrock and the depths of the quarries. Lithology will be carefully logged during drilling. Elevations of all existing and proposed onsite well casings will be surveyed, as will measurement points at each quarry.

Once the proposed wells are drilled, water level monitoring will indicate whether the gravel wells communicate with the bedrock aquifer. The monitoring would compare

water level fluctuations in the bedrock wells with pumping cycles in the OMYA gravel wells #1 and #2, and the Pittsford-Florence Water System well. Water quality sampling at the OMYA gravel wells (nonpotable water) would occur if a hydraulic connection with the bedrock aquifer were observed. The town well will continue to be sampled periodically.

We propose to conduct rounds of well sampling monthly, at first. The sampling frequency will be phased, as shown in the monitoring summary below. The following wells would be tested: 96-1, #2, #5, 96-2, the three proposed wells, the three private wells, and the active PCWS well. The OMYA gravel wells would be included in the program if the testing indicates a hydraulic connection. The wells in use would be sampled at the taps. Other wells, not in use, would be purged with a field pump until groundwater conductivity and temperature readings stabilized. Samples would be analyzed by Endyne for OPP and potential breakdown products by EPA methods 642, 8270C, and 8260.

The three private wells and the Pittsford-Florence water system well have been resampled for analysis by Sci-Test, and results will be reported when available. These wells will be included in the sampling routine.

Additional data on the flow between the Dogleg, Dolomite, and PIQ Quarries and the bedrock aquifer will be collected by installing multi-level piezometers into the sediment. A measurement of gradient will indicate whether the quarry is discharging or receiving water.

B. Surface Water Monitoring

Water quality monitoring in the three onsite quarries will continue. Water will be sampled and tested by Endyne for OPP and potential breakdown products by EPA methods 642, 8270C, and 8260. Samples will be acquired every other week from the quarries until such time as it becomes apparent that concentrations have declined to non-detectable levels.

Currently, we do not plan to sample nearby streams and wetlands since there is no likely risk of contamination.

C. Monitoring Summary

Monitoring will be conducted monthly at first, and will be scaled back through time. We propose a two-year long monitoring program, which will encompass two full hydrologic cycles.

The table below summarizes the monitoring proposal.

Activity	Location	Schedule
Install bedrock monitor wells	3 locations (See maps)	Early 2001
Test for communication between gravel aquifer and bedrock aquifer	3 proposed bedrock wells, OMYA gravel wells, PCWS town well	Following well installation
Install piezometers in Dogleg and Dolomite Quarries	Sediment at edge of the quarries	December 2000
Sample water quality for OPP and potential breakdown products	Dogleg, Dolomite, PIQ	Every other week until contaminants are below detectable levels, then follow schedule for wells, below.
	Wells 96-1, 96-2, #2, #5; 3 proposed wells, 3 private wells, PCWS well.	Monthly for 4 months; quarterly for balance of 1 year; semi-annually for 2 nd year.
Sample water quality for OPP and potential breakdown products	Gravel wells #1 and #2	Monthly for 4 months; quarterly for balance of 1 year; semi-annually for 2 nd year, <u>if</u> communication with bedrock is detected
Measure water levels / gradient	Bedrock wells, 3 quarries, quarry piezometers	Monthly for 4 months; quarterly for balance of 1 year; semi-annually for 2 nd year.

VII. REMEDIATION PROGRAM

The design and feasibility of an active remediation system is currently being evaluated. A remediation plan will be implemented in the event that a plume of OPP is detected in on-site and/or off-site groundwater monitor wells. The results of groundwater monitoring to date do not indicate the need for active remediation.

Potassium permanganate was identified as a potential chemical oxidant for OPP. H&N has utilized potassium permanganate to chemically oxidize other types of organic compound plumes. A laboratory bench treatment test was conducted to determine the effectiveness of potassium permanganate as a chemical oxidant for OPP. The details of the test are presented below.

A. Potassium Permanganate Bench Treatment Test

A sediment core sample from the Dogleg Quarry was split to conduct a laboratory bench test of a potential treatment scheme for chemically oxidizing. OPP into inert breakdown products. Core sample SS#1(0-4") from (15-18') was split into two aliquots. One 20-gram aliquot was reacted with 2.5 milliliters of a 5% potassium permanganate solution (KMnO₄) for 90 minutes. The other 20-gram aliquot was untreated. Each aliquot (about 56% solids) was extracted with methylene chloride as a soil with an ultrasonic disruptor to break up any solid clumps.

Sample ID	Resulting [OPP] (mg/kg)	% Reduction
Control: SS#1 (0-4")	10.6	0
Test: SS#1 (0-4") + KMnO ₄	<0.10	>99.06

OPP in the untreated portion was detected at 10.6 mg/kg, while less than 0.1 mg/kg was detected in the permanganate treated aliquot. The treatment resulted in greater than 99% reduction of OPP concentration in the sludge layer tested, in 90 minutes. Evaluation of the chemical mass balance indicates that significantly more potassium permanganate was added than necessary to oxidize the OPP. Additional testing will be performed to fine tune the ratio of potassium permanganate needed to oxidize the OPP with minimal excess permanganate remaining.

B. Additional Investigation

On December 12, 2000, additional surface water samples were collected from the Dolomite Quarry and submitted to Endyne, Inc. for analysis of OPP and total organic carbon (TOC). The water samples will also be utilized for further testing of the proposed potassium permanganate remediation plan.

The additional investigation will focus on determining potassium permanganate consumption by degradation of TOC and OPP. The presence of varying concentrations of TOC can significantly impact the mass of potassium permanganate utilized to oxidize OPP. Therefore, strict stoichiometric calculations will underestimate the concentration of permanganate necessary to oxidize the OPP. Testing of the waters potentially requiring treatment will provide the most accurate permanganate concentration estimates.

A series of empirical tests utilizing various concentrations of potassium permanganate will be performed to determine the optimal concentration necessary to adequately treat

the OPP. The data generated from the additional testing will be utilized to develop a full-scale remediation plan if it becomes necessary.

REFERENCES CITED

Freeze, R.A., and Cherry, J.A., 1979. Groundwater Prentice-Hall:Englewood Cliffs, NJ.

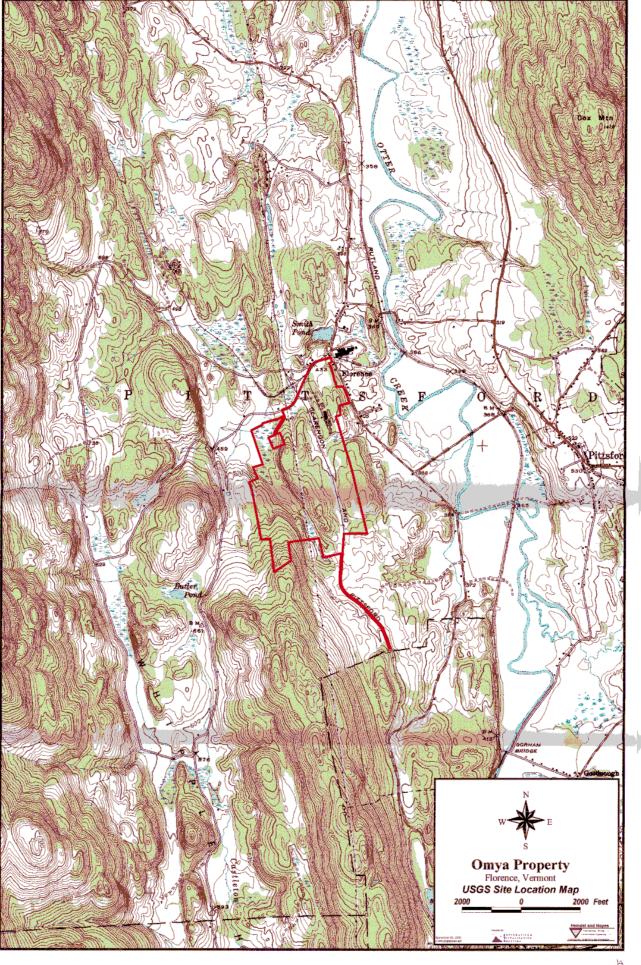
Howard, P.H., Boethling, R.S., Jarvis, W.F., Meylan, W.M., Michalenko, E.M., 1991. Handbook of Environmental Degradation Rates Lewis Publishers: Chelsea, MI.

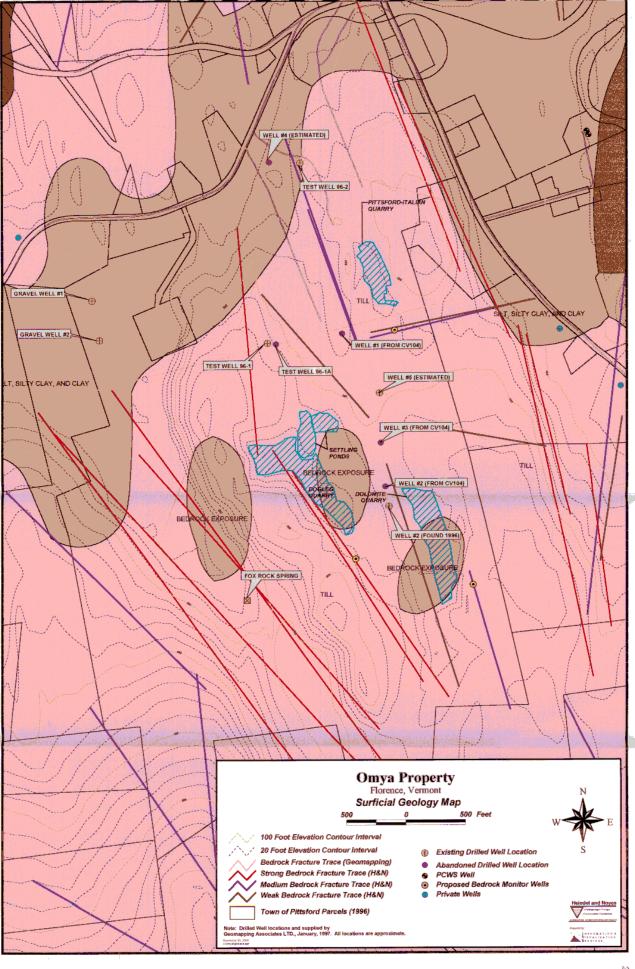
Howard, P.H., Meylan, W.M., 1997. <u>Handbook of Physical Properties of Organic</u> Chemicals CRC Press: Boca Raton, FL.

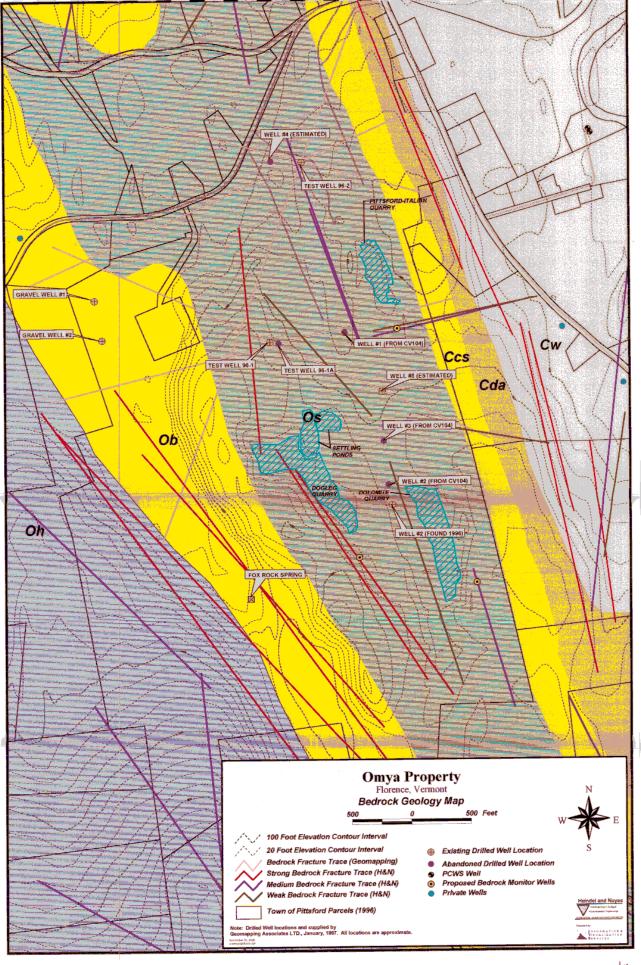
Sutherland, J.B., Fatemeh, R., Kahn, A.A., Cerniglia, C.E., 1995. "Mechanisms of Polycyclic Aromatic Hydrocarbon Degradation". In <u>Microbial Transformation and Degradation of Toxic Organic Chemicals</u>, 1995, Wiley-Liss: New York. Lily Y. Young and Carl E. Cerniglia, Eds.

Winholz, M., Budavari, S., Blumetti, R.F., and Otterbein, E.S., 1983. <u>The Merck Index:</u> An Encyclopedia of Chemicals, Drugs, and Biologicals Merck & Co., Inc.:Rahway, NJ.

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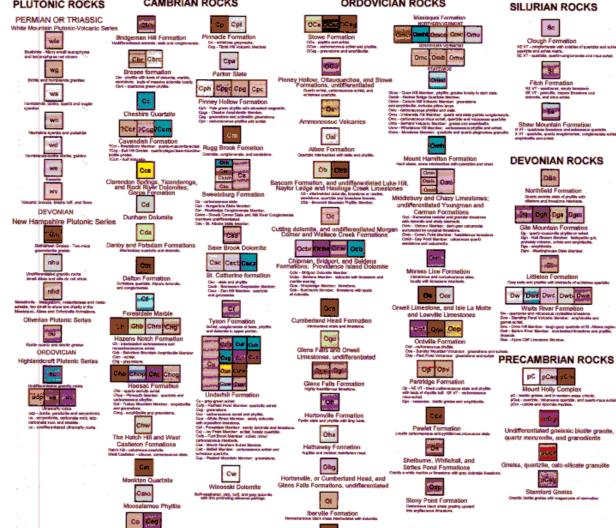






BEDROCK GEOLOGY LEGEND





SOURCE NOTES

Bodrock Geology was digitized and scenned, by Wagner, Heindel, and Noyee, into a PC ARCKNFO database from 1:52500 original State of Vermont bedrock geology base maps (1956-1966) These maps were made available by Dr. Barry Docien, Geology Department Chalipperson, University of Vermont. Bedrock data for most of the state is available, in 15 minute quade, from IVS (802) 985-0437.

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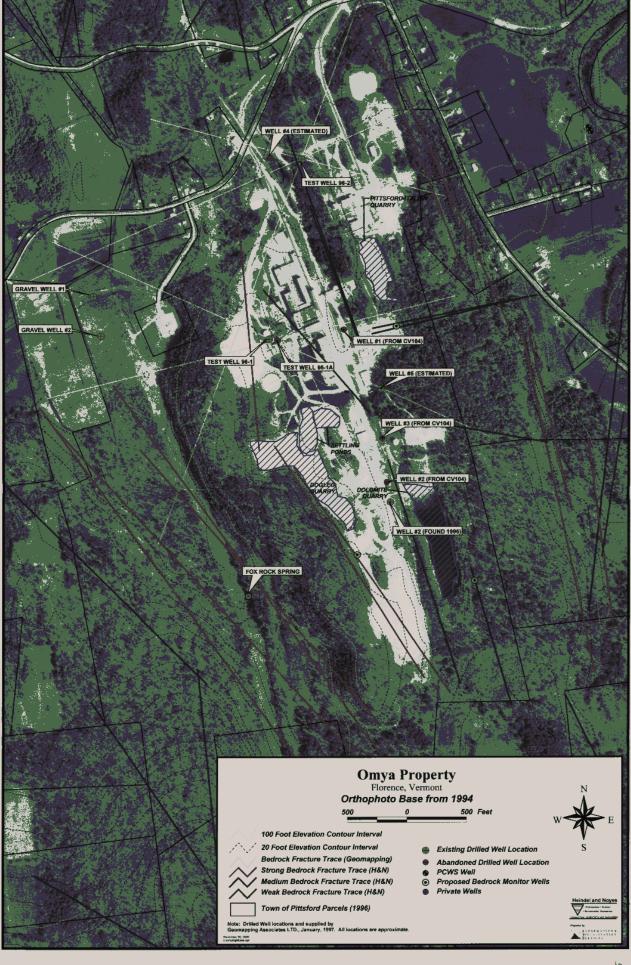
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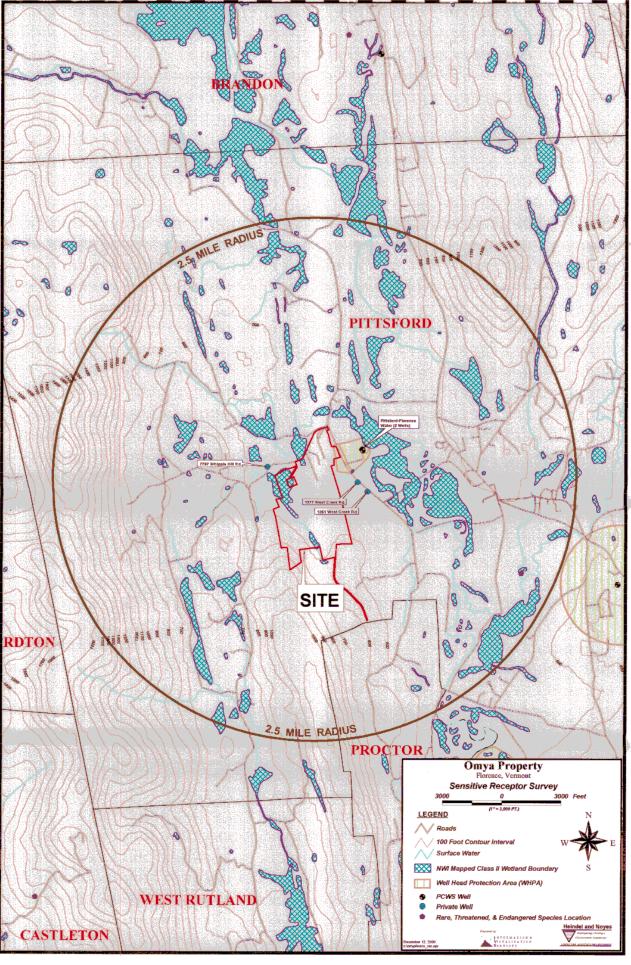
Read Centrafines were generated from pre-1990 1:5000 orthophotos (or better). Road data (RDSnn) is available from the Vermont Center for Geographic Information, VCGI (802) 656-4277.

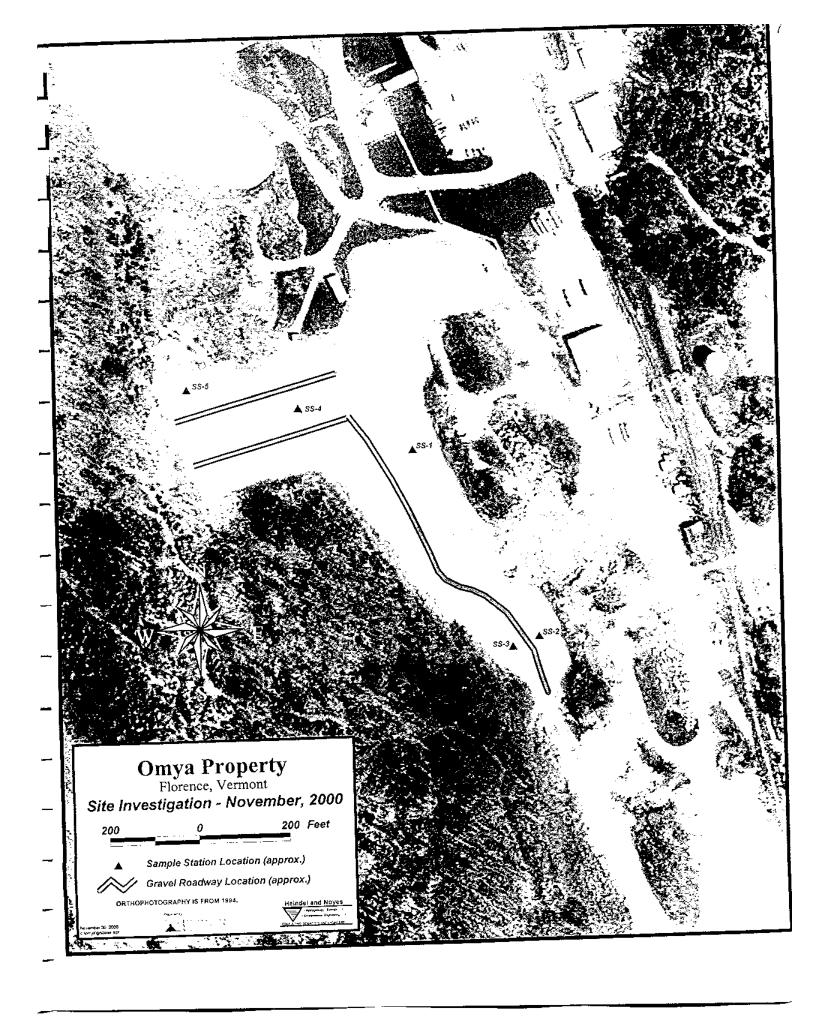
Linear Surface Waters are Digital Line Graph Data, generated from 1:24,000 USGS topographic maps. This data is available from VGIS.

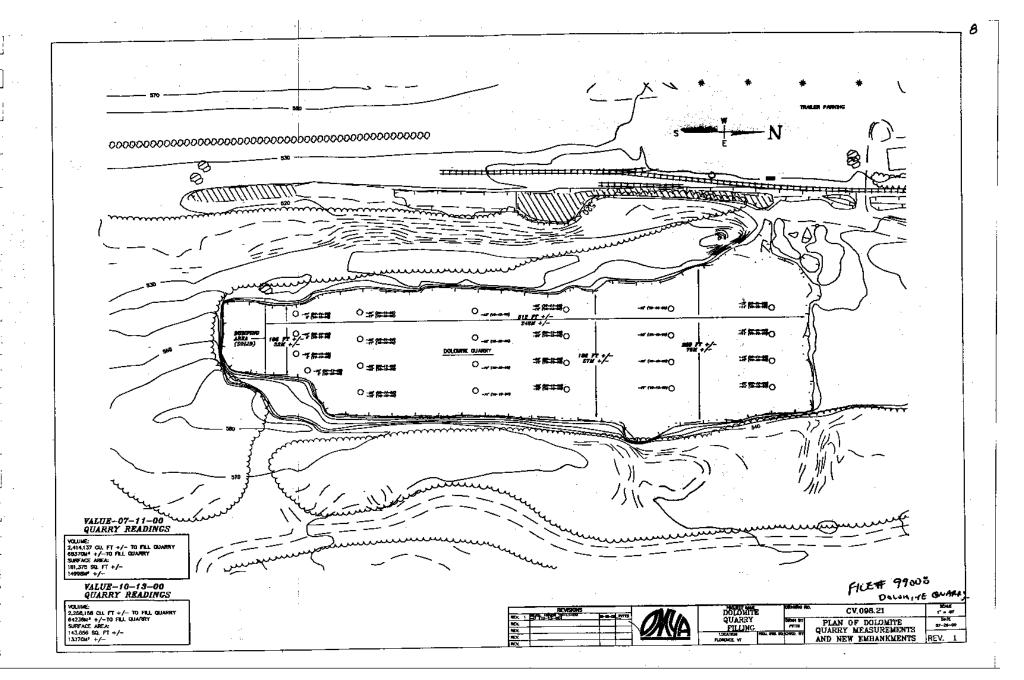
Town Boundaries were digitated from pre-1990 1:2000 USGS topographic maps. This coverage was created by the EPA and is available through VGIS.











Map lorbel: Well #4

Form WR-59 Rev. 7-22

State of Vermont DEPARTMENT OF WATER RESOURCES

WELL COMPLETION REPORT (This report must be completed and DO NOT FILL IN submitted to the Department of Water Resources, State Office Building, Montpeller, Vermont 05602, no later than 60 days after completion of well. Complete or line out all blanks.) Mailing Address (Please locate well on a large TOWN IN WHICH WELL IS LOCATED: scale map to accompany this report.: Maps are available on DATE WELL WAS COMPLETED: request.) 🏻 Domestic PROPOSED USE OF WELL: **Industrial** Other (Specify) ☐ Municipal TWAK ☐ Air Percussion Rotary Cable Tool DRILLING EQUIPMENT Other (Specify) STATIC WATER ... TOTAL DEPTH OF WELL: ft. Diameter.....in. Material... CASING DETAILS: Length... Weight.... SCREEN DETAILS: Make ... Diameter in. Slot Size METHOD OF SEALING CASING TO SCREEN OR BEDROCK: TNAL YIELD TEST: Bailed, or Pumped, or Compressed Air Hours at gallons per minute Water level during yield test..... Give description of formations penetrated, such as: peat, silt, sand, gravel, clay, hardpan, shale, limestone, granite, etc. Include size of gravel (diameter): WELL LOG and sand (fine, medium, coarse, color of material, structure (loose, packed, cemented, hard). For example: Surface to 27 ft. fine, packed, yellow sand; Depth From Ground Surface 27 ft. to 134 ft. gray granite. Surface to YIELD TEST DATA IN G.P.M. If yield was tested at different depths during drilling, List Below · ft. G.P.M. @ G.P.M. @ G.P.M. @ ft. G.P.M. @ WATER ANALYSIS: Has water been analyzed? [] Yes Zi No If Yes, Include Analysis DRILLED BY: DOING BUSINESS AS: Company WELL DRILLERS LICENSE NO. DATE OF REPORT: 5M 6-76

Gravel Well #2

PUMPING TEST RECORD



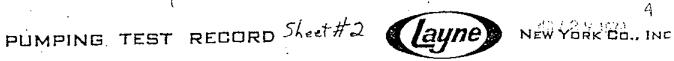
NEW YORK Co., INC

LOCATION: Florence, VT. PUMPING WELL #36" X16" Gravel Packed

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PUMPING TEST RECORD Sheet#3 (Layne)

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	9:00	-	60.00		40.26	,	23.62	26.92	26.56	/	1	\
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QUALITY: TIME FE NN	HARD	PH		DOR.	TASTE	
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PUMP TYPE STAGES			•			
SETTINGFT. TO SUCTION FLANGE	1	ή	<u>'</u>	ij 	·	_
SUCTION FT. OF IN PIPE/HOSE		<u> </u>	<u> </u>	}	<u> </u>	-
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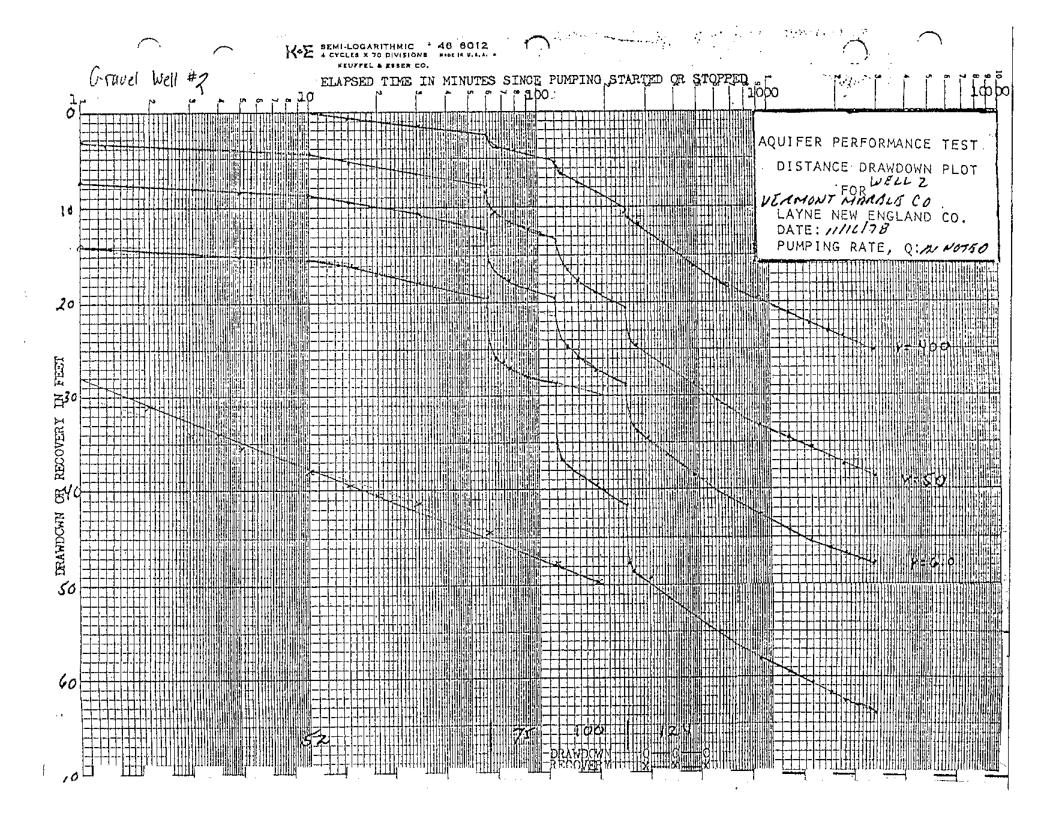
PUMPING TEST RECORD Steet 4 (Layne) NEW YORK Co., INC.

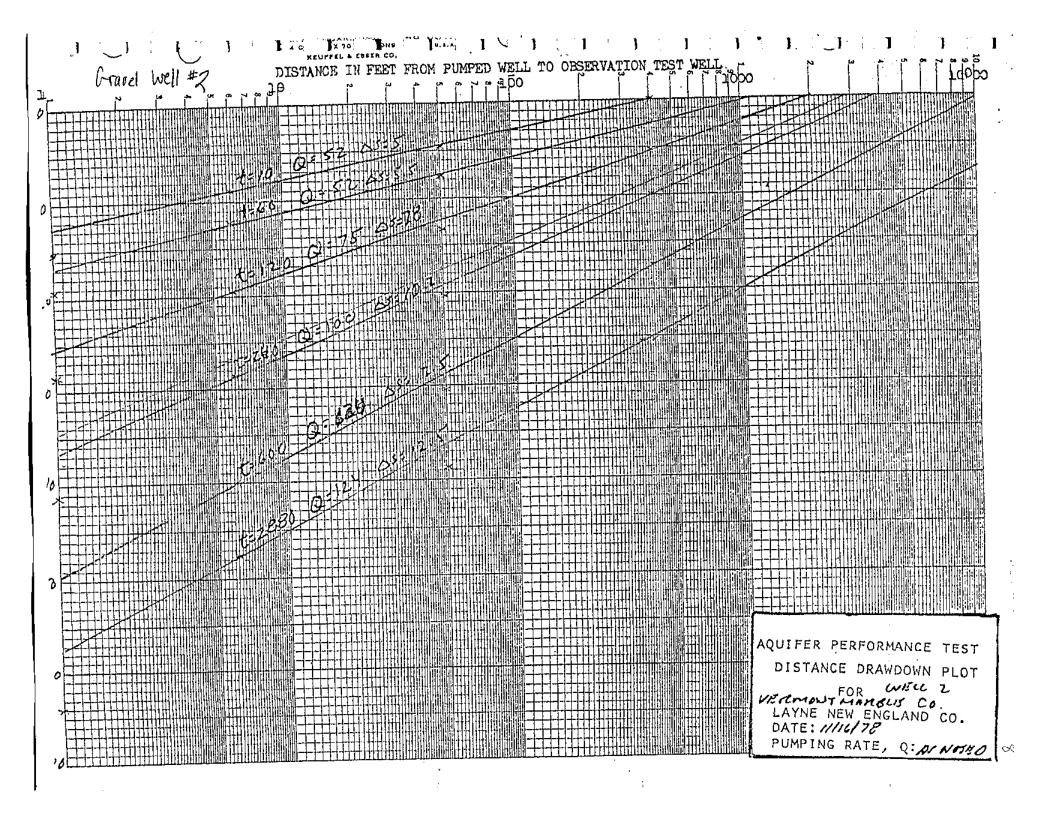
FOR PURCHASER



FOR LAYNE - NEW YORK CO., INC.

		cm cut				DATE	. 261	NOU.	11/		11/5	ivel Paci	امعا
Loc	ATION: _	Flower	ce, V)	-		РОМ	PING	WELL:	9		· Gra	iver faci	N EO
WEI	LL NO	1	····			JOB							
TIME	AIR PRESS.	PUMPING LEVEL	OHS. /	WELL 50	OBS	WELLE .	6	OBS. 7	WELL.	4	темр.	ORIFICE INCHES	GP)
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2:00		60.50'	50.23	40.78'		24.16	, ,	27.45	27.0			24"	12%
3:00	<u></u>	60.50'	50.27	40.88				27,55	27.			24"	
4:00		60.50'	50.31	41.00		24.36		27.651				24"	ì
5:00		60.501	50.43'	1		24.4	2'	27,76'	27,	42'	-	24"	-
6:00		60.751	50.51		<u> </u>			27.85				24	12
7:00		61,00	50,58'	41.28		24.65	y'	27.99	27.5	59		24"	12
7:01		45.00'	39,10'		7.4				<u> </u>			0	10
7:02	!··- 	33,00'	36.05'			<u> </u>						<u> </u>	0
7:03		30.00	34.84'			<u> </u>							0
7:04		30.00	33.82				· · · -		<u> </u>				10
7:05		28.50	,	32.28	1	<u> </u>							0
7:10		26.00	1	1	<u>, </u>	22,5	0					<u> </u>	0
7:15	 	24.50	28.98'	29,05	<u> </u>	21.6	6'		ļ				0
7:20	1	23.50	J .	27.89		20,9	15'	24.20		<u>/3/</u>		<u> </u>	· ·
7:30	- I	22.45	7	1		19.5	71		-	1391	<u></u>		5
8:00	1	19.50	.1	22.85		17.3	41	20.67			<u> </u>		0
9:00	i	16.191	1 , .	19.55		19.6	<u> 2'</u>	18.03		<u>63'</u>			10
10:00		14.08		17.46		12.9	13 ¹	16.31	1 15,	93'	<u> </u>		10
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OF Pr	RIFICE	IN. C	SIZE	ST.	AGES		WE	EL PUM	-, WELL				
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su	JCTION		FT. OF		וא פו	PE/HOSE	 -				<u> </u>		1





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-			· · · · · · · · · · · · · · · · · ·	PIPE 55' × 16" × 3/8" STEEL Grave Well #7
Ť	ORIGINAL GROUND		ELEV.	55' x6" x 3/6" STEEL Grave Well #1
-	LEVEL	 _		The state of the s
- }	Z FILL		1.	
ļ	}			SCREEN: 10' X6" STAINLESS STEEL JENNYOT
- 1			7 -7	Screening 20 1
		4	المحمد أن من المرا الم	PLUG: STAINGELT /TEEL PERTE
- 1	3/ /			1
		1 1		LEGAL PARCHETO 1/1
- {				
-	1/1	- 1	Jord .	SIZE NUMBER
ļ	/ /	<u> </u>	1	STAGES TYPE
				SETTING COLUMN SHAFTING
Ì	GREY /	1		Ingine
ı			. Super 3	BOME
	CLAY			IMP. SHAFI
				STRAINER HEAD AIR LINE
	1 / 1	1 1		MOTOR
				MAKE TYPE
				VOLTS CYCLE
			1/ //	PHASE AMP.
				H. P. RPM
			38	FRAME NON REV.
				MODEL .
				UPPER BRG. GEAR DRIVE
	- 1/1			Mrg. Model
	\ /\$			RATIO SERIAL NO.
		ļ	1 1/.	HVY. THRUST NON-REV.
			1 1/1	ENGINE
	1 1			MFG. MODEL
				RPM CONTINUOUS H. P.
				FUEL SERIAL NO. W.E.L.L.
				STARTED 10/1/78 CLEAR DEPTH 7/ FIRST TEST 10/21/78 METHOD CORLE TOOL
			61'	FINAL TEST GUAR. CAP.
	162			Contract Towns
	SMVD			- 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1
	GNAVEL	-		DISK TO G.W. DRILLER & CARPEN F.
	DND			DATE 10/18/3
	BOULDER	1		STATIC LEV. 73
	7,	_ !	→ -, 7/,	PRODUCTION (&
	MANDPAN			Pump. LEVEL 97
				WATER TEMP. 50
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	1	£_142.		Delli veneri
		Det	120/20/25	Market Control of the
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			•	
				LAYNENEW YORK CO., INC LINDER, N. J.
				WATER SUPPLY CONTRACTORS
			መሮል ለ፣ 4050	VERMONT MARKECO
	1		DEC 04 1978	DRAWN BY LAYNE WELL # / CUSTOMER WELL #
	1		<u>.</u>	LAYNE WELL # / CUSTOMER WELL # /

PUMPING TEST RECORD Grand Well *1



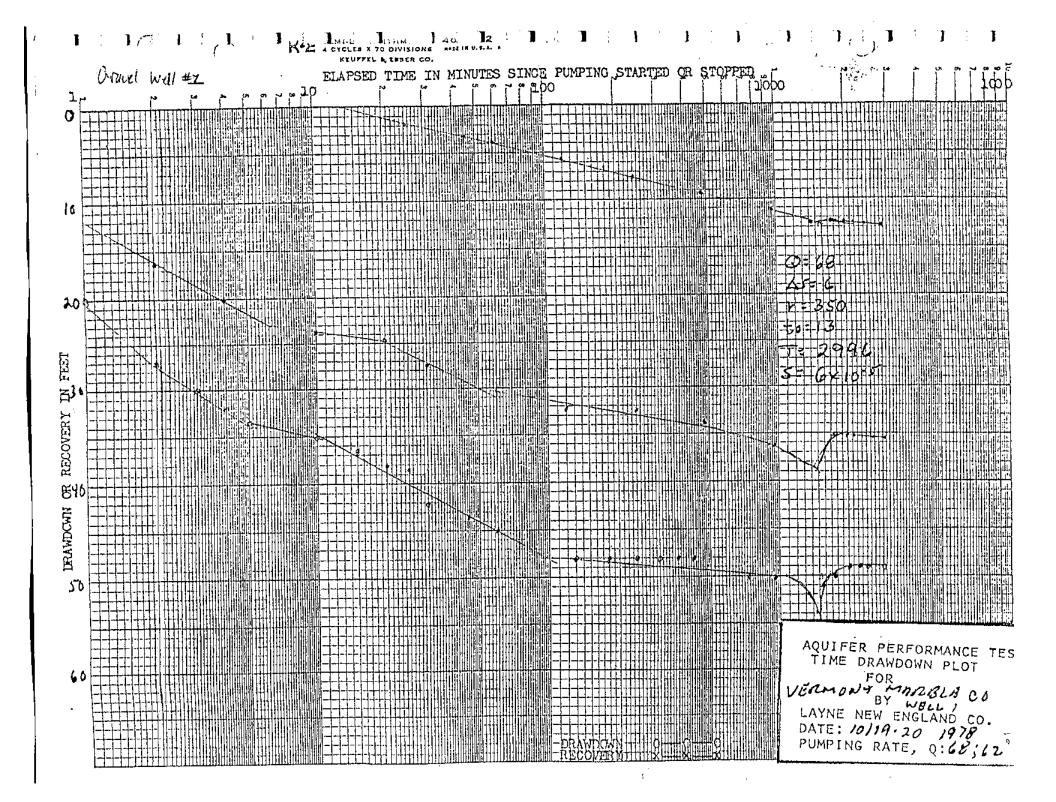
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<u>.</u>	FOR: N	on: [Floren	ce, Ver	most_		PUMP	ing WELL:	#1	G"XC)		
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· 	SETT	ING		FT. OF		IN PIP		<u>a</u>	400'		<u> </u>		
•			PIPE_10		_2	IN PIP	E,HOSE	3	350'				
_								_	FARTIN	NE . NEW	YORK CO.	. INC.	
-			FOR PURC	HASER					FOR LAY			-	

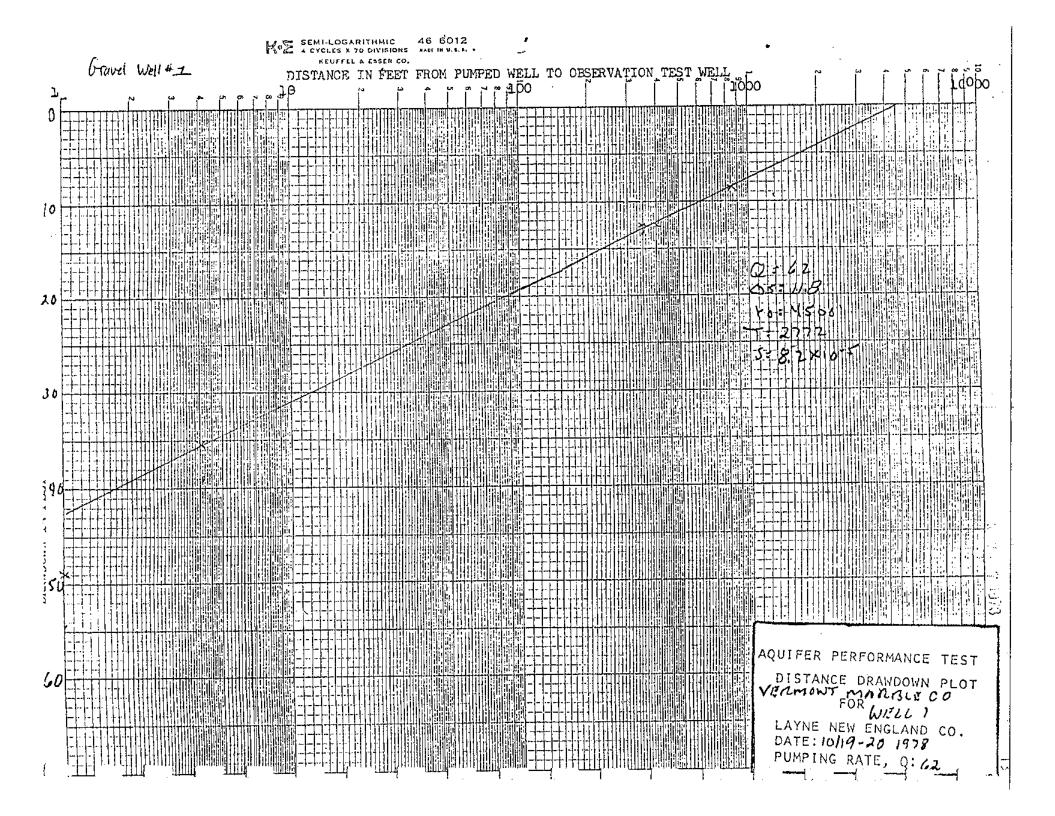
PUMPING TEST RECORD Gravel Well #1



ROY 2.7 . 9 New York Co., Inc

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	TIME	AIR PRESS.	PUMPING LEVEL	OBS.	WELL_3	OBS.	086.		OBS.	OE WELI	ıs.	TEMP.	ORIFICE	
	STATIC	LEVEL	Flasing	1									0	
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		11.00		38.10	15,55	14.60							<u> </u>	
	2:00	l		38,18	15,60	14.66						ļ. <u></u>	1	⊥ '
	3:00	11.00		3810				_						17
1760	4:00	11.00		38,30	16.60	14,70					 .	<u> </u>		<u> </u>
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			FOR PURCHA	SER	, , , , , , , , , , , , , , , , , , , 				FOR	LAYNE	- NEW	YORK CO.,	INC.	





GEOMAPPING ASSOCIATES LID.

PROPERTY: OMYA, INC., VERPOL PLANT

maj label:

Well 96 i

WELL NO. 96-1 BOREHOLE DIAMETER: COORDINATES (SCALED) N TOTAL DEPTH: 500 CASING: 8" I.D., STEEL. 29 LBS. / FT. GROUND ELEV: +/- 534 DRILLING METHOD: AIR ROTARY 8/14/96 - 8/15/96 DATE DRILLED: STATIC LEVEL: NOT MEASURED DRILLER: GREEN MT. DRILLING CO., INC., CENTER RUTLAND, VT SURFACE DESCRIPTION: BLASTED BENCH IN BEDROCK, LAWN AREA

SITE LOCATION: ADJACENT TO SOOK WATER STORAGE TANK

GEOLOG	SIST:	JIM f	PURDY	LOG	DATE: 8/14/96				RC.	SAMPLE
FOOTAGE	BFLA LC	PH#C	GEOLOGIC DESCRIPTION		REMARKS	FRACTURES	ARLET MACHARGE	BET.	LOSS	MT.
25			O - I TOPSOL FILL I - 50' OFF WHITE - LT. GRAY FINE GRANED MARBLE, BLAST FRACTURED NEAR SURFACE	1	NSTALLED 9' OF 8' STEEL CASING W/ DRIVE SHOE SEY B' IN COMPETENT ROCK					
50		-	50 - 500' WHITE - GRAY FINE GRANED MARBLE, OXDE STAINED, THIN HEAVILY STAINED SCHISTOSE MARBLE	ox0E	COARSE DRLL CUTTINGS & - 27 RETURNED AT SURFACE 50 - 450°				<u> </u>	
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175 200		<u>+</u>					ļ	00.1%		
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450	13	<u>+</u>	-							
475		+ -				!		ļ		
500) 	_+	500' END OF HOLE						V	

maj label: well 96-1

Water Supply Division Well Report
Well Statistics

Printed: 11/13/97

Well Tag Number: 1	996-35	Date Report Received:	8/27/96
Well Report Number:	3,168	Map Cell: 16C4	
•	'A, INC.		
Purchaser's Name:	Λ, πτο.		
		Date Well Was Completed:	8/16/96
Town Name: Pittsford	6 L	Date Well Was Completed.	0. (0.00
Purposed Use of Well:	Abandoned		
Reason for Drilling Well:	Test or explo	ration	
Drilling Equipment:			
	☐ Well Ha	s Screen	
Total Depth of Well (in feet):	300.00		
Casing Finish	: Above groun	d, unfinished	
Total Casing Length (in feet):	10.00		
Casing Diameter	; 8.00		
		el Casing	
Mathed of Cooling Cooling		-	
Method of Sealing Casing:		•	
Yield Test Method	•		
Yield Tested At (Gallons pe		5.00	
Static Water Level (in feet):			
	☐ Well is	OverFlowing	
	Has Wa	iter been Analyzed	
	ON: TEST OR E	EXPLORATION & PROVIDE ADDI'NDONED.	TIONAL
Reason for Well Dev	elopment:		
Well Drille	r: Green Mour	ntain Drilling Company, Inc.	
Tax Ma	ıp:		
	Gravel	Well	
Overburden Thickne	ss (in feet):	0	

Water Supply Division Well Report

Printed: 11/13/97

Well Lithology

Lithology

Clay

Town: Pittsford

Well Report Number:

3,169

Starting Depth

Ending Depth

Driller's Description:

1

CLAY

500

Rock, bedrock, ledge, etc.

Driller's Description:

MARBLE (MIXED COLORS)-WATER

WELL NO. / TAG NO. 1996-34 (140 Ordina in Use)

This report must be completed and submitted to

State of Vermont Dept. of Environmental Conservation 103 S. Main St., The Old Pantry Waterbury, VT 05671-0403

WELL COMPLETION REPORT

DEPARTMENT USE ONLY - E.C. _____U.S.G.S. ____ Field Location O Map area____ Latitude Elay.

Location may days after completion of the well. WELL OWNER DAYA, Jour. OR WELL PURCHASER DOTE WELL TOWN PITTSFOAD SUBDIVISION LOT NO. DATE WELL WAS COMPLETED B-15 9 0 PROPOSED USE OF WELL DOMAIN. Delan Englan Comment will be found to Capturellian. Proposed USE OF WELL DOMAIN. Delan town. Delan Comment will be found to Capturellian. Delan Editional States, Domain Comment will be found to Capturellian. Delan Editional States, Domain Comment will be found to Capturellian. Delan Editional States, Domain Comment will be found to Capturellian. Delan Britanian States, Domain Comment will be found to Capturellian. Delan Britanian States, Domain Comment will be found to Capturellian. Delan Britanian Comment will be found to capturellian. Dela	
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LOCATION OF WELL. TOWN PITTSFOOD SUBDIVISION LOT NO. DATE WELL WAS COMPLETED 8-15-96 PROPOSED USE OF WELL. Downstie. Delies. Commeta Cleaning Subdiving Well. Delies. Delies. Commeta Cleaning Subdiving Well. Delies. Delie	-1) 6-1
DATE WELL WAS COMPLETED 8-15-96 PROPOSED USE OF WELL Domestic, Deplet Commetted Comme	-1) 6-1
DATE WELL WAS COMPLETED	-1) 6-1
REASON FOR ORILLING WELL. O New Supply. O Replace Creating Supply. O Cooper Calenting Well. & Treat or Engineering. DRILLING EQUIPMENT. O Cashe Tool. Apolory with A-P. O Oliver TYPE OF WELL. Open Hole in Bearing. O Open End Casing, O Screened or Stating; O Oliver TOTAL DEPTH OF WELL: 500 1441 Supply Supply Cases of Stating; O Oliver CASING FINISH: O above provide, Foreign grands, Unfalliabled, O Barring, O in Pil., O Reasoned, O Name used, O Oliver CASING DETAILS: fairstangly 10/2 is Langill believe 2 9 is Olive. Stating of STOTE wit. 79 is./fil LINER OR INNER CASING DETAILS: Cangill believe 2 11 Diameter in Material Weight in Material in Material Oliver. SCREEN DETAILS: Male and Type. It offers Share, O Grant - Type	-1) 6-1
DRILLING EQUIPMENT. © Cable Tool. Apolory with A-P, © Olare TYPE OF WELL. Apolor Hole in Bestrois. © Open End Casing, © Screened or Stating, © Older TOTAL DEPTH OF WELL: 500 Jack Soling and Casing, © Screened or Stating, © Older CASING FINISH: © above promod, Formbook, Apolor promod, United Soling, © In Pil., © Removed, © Name used, © Older CASING DETAILS' total toughts 10/2 11 Longits below 1.3 9 11 Old. 8 14 Material STOCK 12 Mil. 13 /11 LINER OR INNER CASING DETAILS: Congressed 12 Olders Soling 13 Olders 14 Material 15 /11 METHOD OF SEALING CASING TO BEDROCK, Apolice Soling, © Great - 1728 Dollard 15 /11 16 Bedrack 11 m Bedrack	16-1
DRILLING EQUIPMENT. © Casio Toss. Apology with A-P. © Oliver TYPE OF WELL. & Open Hole in Baseroca. © Open for Casing, © Screened or Stating; © Oliver TOTAL DEPTH OF WELL: 500 Jeel basew and surface CASING FINISH: © Above proved, Formore, & Space proved, Unfilliated, © Buring, © In Pill, © Removale, © Name used, © Oliver CASING DETAILS: fairstample 10/2 11 Langib below 1.3 9 11 Ola. 8 14 Material STEEL 12 79 19./11 LINER OR INNER CASING DETAILS: Langib below 1.3 11 Olambiar 12 Olambiar 13 Material 14 Material 14 Material 15 Material	
TYPE OF WELL. Some rest in Secreta. O one feet Casing, O Secretarial Colors TOTAL DEPTH OF WELL: 500 1441 Secretarial Colors CASING FINISH: O above provide fractional Colors young, Unduring O Secretarial Colors of Casing Details: Tarastery of 10/2 is Leagth below 12 9 is 014. 8 is Material STOCK of 19./11 LINER OR INNER CASING DETAILS: Cangin below 12 9 is 014. 8 is Material STOCK of 19./11 LINER OR INNER CASING TO BEDROCK. South Store, O Great - 17.04 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000 0.000	
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METHOD OF SEALING CASING TO BEDROCK Delice State Delical - 1794	
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3 SCREEN DETAILS, Make and From	
3. SCREEN DETAILS, Hear and Free	
	
Sigt Size	
4. YIELD TEST Decises. Compressed Air, for	
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STATIC WATER LEVEL	
S WATER ANALYSIS me the only from endered T O Tot SK No. 11 Ton. Where	
7 SPECIAL NOTES:	
8. WELL LOG	A11, 4-0/b
Out from Land Surface water formation Description "Search badget placed afford name and quiddington let number	14 184 11911
Free Free Original Clay	
1 500 V MARBLE (MIXED COLORS)	
STAR: PU	1 -
TALK H	ou se
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ि दिस्पन् मास्य र	ांत्र म्र
20. TESTED YIELD WELL DRILLED BY: RAY & JEFF LEOWARD	
If the grand and formed or detroined marine during distance, and desired	tic
DOING BUSINESS AS: GREEN MT. DRIlling Co.	<u></u>
420 5 Day 04 7	Uh,
REPORT FILED BY: Armented Signature	
1	U

PROPERTY: OMYA. INC., VERPOL PLANT **WELL NO. 96-1A** (SCALED) N BOREHOLE DIAMETER: COORDINATES TOTAL DEPTH: 300 CASING: 8" LD., STEEL, 29 LBS, / FT. +/- 500 GROUND ELEV: DRILLING METHOD: AIR ROTARY 8/16/96 DATE DRILLED: DRILLER: GREEN MT. DRILLING CO., INC., CENTER RUTLAND, VT STATIC LEVEL: NOT MEASURED SURFACE DESCRIPTION: BLASTED RAMP IN BEDROCK, ASPHALT DRIVE SITE LOCATION: NORTH SIDE OF CASCADE MILL LOWER LEVEL ACCESS ROAD

map label: weil 96-1 A

GEOLOGIST: JM PURDY			LOG DATE: 8/16/96							
FOOTAGE	C GRA		BEOLOGIC DESCRIPTION		REMARKS	FRACTURES	ARLET Nacharge	RET.	C. LOSS	BAMPLE INT.
25		Ŧ	O - I' FILL, I - 45' HEAVILY WEATHERED, FRACTURE OXDE STANED, WHITE, FINE GRANED MY		NSTALLED 9' OF 8' STEEL CASING W/ DRIVE SHOEL SET 8' N COMPETENT ROCK	45'	5 GPM			30') 33'
50		-	45' MUD SEAM, HEAVLY WEATHERED, O STANED MARBLE, CLIVE GRAY CLAY FIL SOLUTION FRACTURE	LEU	COARSE DRLL CUTTINGS (1 - 21) RETURNED AT SURFACE 0 - 300					22.
<i>7</i> 5			55 - 80' HEAVLY WEATHERED, FRACTUOXDE STANED, WHITE FINE GRANED MASS - 90' OFF WHITE - LT. GRAY FINE	RED. \R8LE	Sola neer o			 ရ		80°
100	E		GRANED MARBLE 90 - 300' WHITE FNE GRANED MARBLE 0XDE STANED, THIN HEAVLY OXDE STANED SCHISTOSE MARBLE	5.				ING FLUID		10°
125		t 	STANED SCHISTOSE MARBLE		OCHRE COLOR AND FINE SEDMENT IN ARLIFT DISCHARGE DOES NOT CLEAR WITH DEVELOPMENT	:		AIR / DRILLING		HG.
150		E						R		***
175		- +						RETURN		80"
200		+						100%		300,
225	E	+								120'
250		± £,								240°
275		<u>_</u>								280"
300		+	300' END OF HOLE					V		
325										
350										
375										
400	1=									
425	1=									
450	13									
475										
500)]									
			<u> </u>		l					Щ.

WELL NO. / TAG NO.

State of Vermont

Dept. of Environmental Conservation

DEPARTMENT USE ONLY

1996-35	103 S. Main St., The Old Pantry Waterbury, VT 05671-0403	E.C. U.S.G.S
which could biviting III South Miles Street	WELL COMPLETION REPORT	LatitudeElev Longitude
The Old Pantry, Waterbury, VT 05671-0403 to later than 90 days after completion of the well.	Location map attached to WCR	Date in Town Files D
WELL OWNER OMYA	Tuc P.O. Box 10	Florence, UT. 05744
	Promi	or Halland Adelace
Roses	PITTSFORD SUBDIVISION	LOT NO.
	· 8-14-49	P.o. # 82501
ASSESSED HER OF WELL'S	Quantity of glass	(w-z.)
THE PART OF THE LINE WEL	.L Sveply, Asplace Estating Supply, Overson Estating	Well, X test or Exploreties. W96-1A
•	Prazies Additional Supply. O Other	
DRILLING EQUIPMENT O	ible Tool Saldry with A-F. C CIAN	
TVOE OF WELL X	edrack, O Open End Casing, C Screened or Sinited; C Other	
TOTAL OF OTH OF WELL	300 feel below had owlded	
· · · · · · · · · · · · · · · · · · ·	Manager of Samuel Control of the Pill Assessed	C Mone wood, C Other
CASING DETAILS THE TOTAL	3/2 11 Lingto believe 3 4 11 Dia Bun Maris	161 37 CCC 40 40 CC (18./1)
		of Watch!
2. METHOD OF SEALING CASIN	TAILS, Leagle and	Oritina m. Agia 11 in Bodrecs
	(I) Other	
4. YIELD TEST! 🛭 منابع , 🔾 وشنيع	Compressed Air, for	_ Gallans per minute _ Permanant Airline installed
	face below land surface . Data or Elma managerad	
6 WATER ANALYSIS	Ell to be Abandared ber	RUCK OF MUDSER
17 SPECIAL NOTES:	D GUANTITY OF WATER.	
IB. WELL LOG	2 3 4 4	19. SITE MAP Shows purchased structure such as buildings, as pric tasks, and/or a structure fund may be defined and the self-
Gapin from Land Surface marer feet feet dearing	Farmetren Genérication	
Grown / F'	-//	
	ARBLE (MIXED GloRS)	MARBLE CLIFF ?
1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1		2. 1. 8 0. 80
1	40 SEAM (DIRTY WATER)	DRIVE
45 V M	40 SEAM (DIRTY WATER)	
		CRUSHING PLANT
מח דבידה עובי ה	WELL DRILLED BY: 1974	STEFF LEONARD
20. TESTED YIELD	,	
4401 641	DOING BUSINESS AS: GA	EER MT. Unilling Log IN
45	5	immend L. Teon and &
	REPORT FILED BY:	Aumorited Signalure
	9.23	-96 WELL POHLEDS IN NO 5/

Water Supply Division Well Report Well Statistics

Printed: 11/13/97

Well Tag Number:	1996-34	Date Report R	
Well Report Number:	3,169	Map Cell:	1604
Owner's Name:	OMYA, INC.		
Purchaser's Name:			0.45/06
Town Name: Pittsford		Date Well Was Con	npleted: 8/15/96
Purposed Use of \	Vell: OTHER		
Reason for Drilling V	Vell: Test or ex	ploration	
Drilling Equipr	nent: Rotary (A	P)	
ı	☐ Well	Has Screen	
Total Depth of Well (in t	eet): 500.00		
Casing F		ound, unfinished	
Total Casing Length (in t			
Casing Diar			
		Steel Casing	
Method of Sealing Ca	sing: Drive sho	e only	
Yield Test Me		sed air	
Yield Tested At (Gallo		5.00	
Static Water Level (in			
Otatio Mater Estat (☐ Well	is OverFlowing	
	Has	Water been Analyzed	
Comments:			
Reason for We	li Development:		
		ountain Drilling Company, Ir	ıc.
	ах Мар:	•	
	•	vei Well	
Overburden Th	ickness (in feet):	1	
Overpuiden in	ionitodo (m. 1044).		

Water Supply Division Well Report

Printed: 11/13/97

Well Lithology

Town: Pittsford

Well Report Number:

3,168

Starting Depth

Ending Depth

Lithology

,

*

300

Dirt, soil, topsoil, loam

Driller's Description:

FILL

Rock, bedrock, ledge, etc.

Driller's Description:

1

MARBLE (MIXED COLORS)-WATER

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OTTEN I __ OT

PROPERTY: OMYA, INC., VERPOL PLANT WELL NO. 96-2 BOREHOLE DIAMETER: (SCALED) COORDINATES TOTAL DEPTH: 590" CASING: 81 I.D., STEEL, 29 LBS. / FT. +/- 484 GROUND ELEV: AIR ROTARY DRILLING METHOD: 8/20/96 DATE DRILLED: STATIC LEVEL: NOT MEASURED DRILLER: GREEN MT. DRILLING CO., INC., CENTER RUTLAND, VT SURFACE DESCRIPTION: EDGE OF WOODS, SIDE OF GRAVEL ROAD, LOW RIDGE CROSS-OVER ROAD. EAST + WEST PLANT AT INTERSECTION SITE LOCATION:

LOG DATE: 8/20/96 JIM PURDY GEOLOGIST: ARLET CRG. SAMPLE DISCHARGE RET. LOSS MT. FRACTURES GRAPHIC REMARKS GEOLOGIC DESCRIPTION POOTAGE LOG NSTALLED 9' OF 8' STEEL CASING W/ DRIVE SHOE, SET 5' IN COMPETENT ROCK 0 - 3' FINE SANDY LOAM TILL 25 3 - 100' LT. GRAY - BUFF DOLOMITE 50 75 COARSE DRILL
CUTTINGS RETURNED
SURFACE 1-2 SIZE)
100-180', OCHRE RED
STAIN 100 - 180' LT. GRAY - BUFF DOLOMTE WITH HEAVY OXDE STAIN FRACTURED 100 125 150 175 3 180 - 380' LT. GRAY - OFF WHITE DOLOMITE OXDE STANED, INTERVALS 220 - 260' WATER BEARING INTERVAL NOT WEATHERED ю сем 196 200 / DRILLING 200 22.73 225 ₹ î. 250 200 RETURN 275 20 GPM WATER SEARING INTERVAL NOT WEATHERED 290 300 325 350 375 380 - 400' REDDISH BROWN - GRAY - BUFF ALTERED DOLOMITE ALTERATION ZONE AT DIKE CONTACT 400 400 - 460' DK. GRAY DIABASE DXE. FINE GRANED DIKE ROCK FRESH NO OXDE STAIN PROBABLE HIGH ANGLE DIXE 70 - 90 DEG. FROM VERTICAL. 425 450 ALTERATION ZONE AT OKE CONTACT 460 - 480' REDDISH BROWN - GRAY - BUFF ALTERED DOLOMITE 475 480 - 590' LT. GRAY - GRAY DOLOMITE NO OXDE STAINING 490° 400°5 500 590' END OF HOLE

Map label: well 46-2

State of Vermont Dept. of Environmental Conservation 103 S. Main St., The Old Pantry Waterbury, VT 05671-0403

DEPARTMENT	USE	ONLY
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E.C		. U.S.6	G. S	
Field Locati				
Latitude	<u> </u>	•	-Elev	
Longitude Scale: 62,59	•		Торо	 r
Scale: 62,59	00 □,2	5,00	00,24	0000

This report must be completed and submitted to the Department of Environmental Conservation, WELL COMPLETION REPORT Water Supply Division, 103 South Main Street,

The Old P	miry. W	aterbury, \ after compi	/T 0567 ccion of	I-0103 no the well.	Location map attached to WC	R	Scale: 62,500 [], 25,000 [], 24,000 [] Data in Town Files []
WEL	L OY OR	NER	00	IYA, INC.	P.O. Box 10	Permonest M	Florence, by, 05744
· WEL	_	IRCHAS	ER		<u></u>		Kalibig Address
	. –			OWN PITTSE	at D sugaryising		
Loc.	AHUN	Or WE	LC. 1	OWNN	30001VISION		P. 4 8 2 5 0 1
DAT	E WE	LL WAS	COM	PLETED 8-7			
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					ry 415 ±-P, □ 01641		
TYP	E OF	WELL!	ו••	Halle in Badress , () Open (ne Ceeing, 🖸 Screened or Slottad) 🗇 O	l e 41	
					feel below land switece		
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O. CAS	ING D	ETAILS:	1419114	414 10/2 12 Leads	18 ballou L 5 <u> </u>		STEEL W. 29 18/11
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		•			O 01N44		
3. SCF	REEN	DETAILS	3; 444	and Type			stangthtt.Diameterm,
\$161 3	i-14		. 01414 14	lag af sçidan in lant þelaw l	and surfacef1 , Gravel s	actifused: G	ravel Sire or Type
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5 STA	TIC W	ATER	FVFI				, Overtions at
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				water they work to a			
	נו ונ						19. SITE MAP
		4~4 5+/1444				Υ	Show permanent afracture such as buildings, supric tanks, and/or permanent afracture and leading and leading and leading and the series
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1111		fael		Orders the two are			
-				1 O	DOING BUSINESS AS:	<u> </u>	EN Mr. DRIlling Co. Inc
		00		15	REPORT FILED BY:	lain	in l. Tenaul fr.
	_	<u>, </u>			HEFORT FILED BY.	10	Authoritied Signature
}					1 25 25 25 27: 8	- 23~	96 5/

Well 96-2

Water Supply Division Well Report
Well Statistics

Printed: 11/13/97

Well Tag Number: 199	6-36 Date Report Received: 8/27/96					
Well Report Number:	3,167 Map Cell: 16C4					
Owner's Name: OMYA	, INC.					
Purchaser's Name:						
Town Name: Pittsford	Date Well Was Completed: 8/20/96					
Purposed Use of Well:	Domestic					
Reason for Drilling Well:	Test or exploration					
Drilling Equipment:	Rotary (AP)					
1	Well Has Screen					
Total Depth of Well (in feet):	590.00					
Casing Finish:	Above ground, unfinished					
Total Casing Length (in feet):	10.00					
Casing Diameter:	8.00					
	☐ Not Steel Casing					
Method of Sealing Casing:	Drive shoe only					
Yield Test Method:	Compressed air					
Yield Tested At (Gallons per	Minute): 15.00					
Static Water Level (in feet):	0.00					
•						
	Has Water been Analyzed					
Comments: WELL REASON SUPPLY.	N: TEST OR EXPLORATION & PROVIDE ADDITIONAL					
Reason for Well Deve						
Well Driller:						
Tax Map	•					
	Gravel Well					
Overburden Thickness	s (in feet): 3					

Well 96-2

Water Supply Division Well Report

Printed: 11/13/97

Well Lithology

Town: Pittsford

alian ar Danasala

Well Report Number:

3,167

Starting Depth

Ending Depth

Lithology

- 0

3

3 Clay

Driller's Description:

CLAY

Rock, bedrock, ledge, etc.

Driller's Description:

ILDEE (MINER C

MARBLE (MIXED COLORS)-WATER

390

490

390

Rock, bedrock, ledge, etc.

Driller's Description:

DARK MARBLE

490

590 Rock, bedrock, ledge, etc.

Driller's Description:

MARBLE (MIXED COLORS)

PRECAUTIONARY STATEMENTS

HAZARDS TO HUMANS AND DOMESTIC ANIMALS.....

DANGER: Do not breathe dust. Corrosive. Causes eye and skin initalion. Do not get in eyes, skin or on clothing. Wear goggles or face shield and rubber gloves when handling. Harmful and can be fatal if swallowed or inhaled.

ENVIRONMENTAL HAZARDS

This pesticide is toxic to fish: Do not discharge effluent containing this product into takes, streams, ponds, estuatios, oceans or other waters unless in accordance with the requirements of a National Pollutant Discharge Elimination System (NPDES) permit and the permitting authority has been notified in writing prior to discharge. Do not discharge effluent containing this product to sewer systems without previously notifying the local sewage treatment plant authority. For guidance, contact your local State Water Board or Regional Office of the EPA.

STATEMENT OF PRACTICAL TREATMENT

If in eyes, flush with water for at least 15 minutes. Call a physician or get medical attention.

If on skin, wash immediately with soap and water. Wash thoroughly after handling.

If inhaled, immediately remove person to fresh air. Then get immediate medical attention.

If swallowed, immediately drink large amounts of water. Do not induce vomiting. Call a physician immediately.

DIRECTIONS FOR USE

It is a violation of Federal Law to use this product in a manner inconsistent with its labeling.

Detailed directions for use can be found in the accompanying technical brochure entitled "P-1® Preservative Directions for Use".

Information on the chemical and physical properties and efficacy can be found in the accompanying brochure entitled "P-1® Preservative - Product Information".

Obtain and read both the "P-1® Preservative - Directions for Use" and the "P-1® Preservative - Product information" before undertaking the use of P-1® Preservative in order to avoid process mistakes.

NOTE TO USER: For worker protection during mixing, loading and during application, wear a hat, long sleeve shirt, and long legged trausers or overalls. In addition, during mixing and loading, wear rubber or neoprone gloves and protective goggles. Protective dothing should be laundered separately following application.

P-1® PRESERVATIVE

FOR THE PRESERVATION OF POLYMER EMULSIONS, (COATING, PVA SYSTEMS AND RUBBER), THICKENERS (STARCH, CELLULOSE, GUAR, ETC.), PAPER, STARCRES, CLAY SLURRIES AND SLURRIES.

Active Ingredient: o-Phenylphenol Inert Ingredients	99.9% 0.1%
High migrations in the same of	100%

STOP • READ LABEL BEFORE USE

DANGER

• KEEP OUT OF REACH OF CHILDREN

LOT NO.
EPA Reg. No. 39967-11-67869
EPA Est. No.: 39967-DEU-01
kg
Made In U.S.A.

SOLD BY:



VERICHEM INC.

3499 Grand Avenue Plttsburgh, PA 15225

REV. 11/98

PESTICIDE STORAGE AND DISPOSAL

PESTICIDE STORAGE: Keep containers tightly and securely closed. Do not contaminate water, tood or feed by storage. Store "Defow 122°F (50°C), protect from exposure to light and moisture. The product has unlimited shelf life if protected from light and moisture. The flakes agglomerate if the storage temperature is too high. See Verlchem's Material Safety Data Sheet "P-1® Preservative" for additional instructions.

PESTICIDE DISPOSAL: Do not contaminate water, food or feed by storage and disposal. Pesticide wastes are acutely hazardous. Improper disposal of excess pesticide spray mixture, or rinsate is a violation of Federal Law. It these wastes cannot be disposed of by use according to label instructions, contact your State Pesticide or Environmental Control Agency, or the Hazardous Waste Representative at the nearest EPA Regional Office for guidance.

CONTAINER DISPOSAL: Triple rinse (or equivalent). Then offer for recycling or reconditioning, or puncture and dispose of in a sanitary landfill, or by incineration, or, if allowed by state and local authorities, by burning. If burned, stay out of smoke.

CHEMICAL OR PHYSICAL HAZARDS

Do not store near heat or open flame. Store in a cool, dry place away from open flame and extreme heat. Store in original container and out of reach of children preferably in a locked storage area. Handle and open container in a manner as to prevent spillage. If container is feaking, invert to prevent leakage. If the container is leaking or meterial is spilled for any reason or cause, carefully dam up spilled material to prevent runoff. Refer to Precautionary Statements on label for hazards associated with the handling of this material. Do not walk through spilled material. Absorb spilled material with absorbing type compounds, and dispose of as directed for pesticides above. In spill or leak incidents, keep unauthorized people away. In case of emergency, call CHEMTREC at 800-424-9300.

SEE SIDE PANEL FOR STATEMENTS OF PRACTICAL TREATMENT AND PRECAUTIONARY STATEMENTS.

NONWARRANTY: Seller (whether manufacturer, distributor or other) makes no warranty express or implied, including the implied warranty of merchantability, regarding this product except the composition as set forth in the ingredient statement heroin, and shall not be liable for special or consequential damages, the exclusive remedy being replacement of the product. Buyer or user assumes all rick of possession, handling or use of this material when such use and/or handling is contrary to label instructions.

P-1@ Preservative is a registered TM of Verichem Inc., Pgh., PA.

MATERIAL SAFETY DATA SHEET

CAS NO: 90-43-7

DATE: 1/01/98

SUPERSEDES: Original

CHEMICAL PRODUCT AND COMPANY IDENTIFICATION 1.

PRODUCT NAME:

P-1 PRESERVATIVE

SYNONYMS:

Ortho phenyl phenol, ortho hydroxy biphenyl, 2-phenyl phenol

1900,11-20

CHRMICAL FAMILY:

· Phenol

MOLECULAR FORMULA: C12H10O

MOLECULAR WEIGHT:

VERICHEM, 3499 Grand Avenue, Pittsburgh, PA, USA, 15225 (412+331-7299, 8:30 am to 5:00 pm; 412-521-6393 24 hours).

EMERGENCY PHONE: For any emergency involving spill, leak, fire, exposure, or

accident call CHEMTREC: 1-800-424-9300. Outside the USA and Canada call: 202-483-7616.

COMPOSITION/INFORMATION ON INGREDIENTS 2.

REGULATED COMPONENTS

COMPONENT	CAS NUMBER	PERCENT	Worker Exposure	REFERENCE
O-phenylphenol	90-43-7	99.9%	Noue Established	,

HAZARDS IDENTIFICATION 3.

EMERGENCY OVERVIEW

APPEARANCE AND ODOR: White flakes with phenolic odor.

EFFECTS OF OVEREXPOSURE: May cause skin and respiratory irritation. Causes eye

burns, Harmful if swallowed.

PHYSICAL/CHEMICAL HAZARDS:

4. FIRST AID MEASURES (IN CASE OF CONTACT)

EYE: Immediately flush with plenty of water for at least 15 minute and get medical attention.

SKIN: Immediately flush with plenty of water for at least 15 minutes, wash following rinse with soap and water. If irritation persists, consult a physician.

ORAL: If swallowed, do not induce vomiting. Give victim glass of water. Call a physician immediately. Never give anything by mouth to an unconscious person. **DERMAL**:

INHALATION: Remove to fresh air, administer oxygen, and artificial respiration. Call a physician immediately. Never give anything by mouth to an unconscious person.

5. FIRE FIGHTING MEASURES

FLAMMABLE PROPERTIES

FLASH POINT:

FLAMMABLE LIMITS (% BY VOLUME):

280°F/138°C.

Lower explosive limit 1.4% Upper explosive limit 9.5%

968°F/>520°C.

AUTOIGNITION TEMPERATURE: DECOMPOSITION TEMPERATURE:

EXTINGUISHING MEDIA AND FIRE FIGHTING INSTRUCTION

In case of fire use water spray, alcohol foam, carbon dioxide, or dry chemicals. Exercise caution when fighting fires involving chemical substances. Respiratory protection is required. Burning will produce toxic fumes. Containers exposed to heat will result in pressure increase in containers. Cool hot containers with cool water.

6. ACCIDENTAL RELEASE MEASURES

STEPS TO BE TAKEN IN CASE MATERIAL IS RELEASED OR SPILLED

Immediately dike material to keep away from water sources. Absorb with vermiculite or other suitable absorbent. Discard in accordance with Federal, State, and local regulations. Residual amounts left on surface maybe cleaned with soap and water. Rinse residuals to municipal treatment plant.

7. HANDLING AND STORAGE

FROM : VERICHEM

EXPOSURE GUIDELINE(S): Avoid breathing dirt. Avoid contact with eyes and skin. Wash thoroughly after handling. Reseal containers after use.

VENTILATION: General recommended.

SHELF LIFE: Greater 2 years.

PERSONAL PROTECTION 8.

EYE PROTECTION REQUIREMENTS: Chemical goggles or face shield. SKIN PROTECTION REQUIREMENTS: Rubber gloves with long sleeved shirt to minimize contact.

VENTILATION REQUIREMENTS: Use local and general exhaust ventilation to control levels of exposurc.

RESPIRATOR REQUIREMENTS: An air-purifying particulate/organic vapor cartridge respirator is required any time inhalation of dust or vapor is possible Observe OSHA regulations for respirator use (29 CFR 1910.134).

9. PHYSICAL AND CHEMICAL PROPERTIES

White flakes. APPEARANCE AND ODOR:

547°F/286°C. BOILING POINT: 134°F/56-58°C. MELTING POINT:

0.0007 @ 68°F/20°C; **VAPOR PRESSURE:**

212°F/100°C. 1.0

1,26 at 68°F/20°C. SPECIFIC GRAVITY:

Unknown. VAPOR DENSITY:

PERCENT VOLATILE (BY WEIGHT): Negative.

6.1 of a saturated solution. pH:

SATURATION IN AIR

(PERCENT BY VOLUME): Unknown.

Unknown. **EVAPORATION RATE:** 2,000 mg/l @ 68°F/20°C. SOLUBILITY IN WATER:



TECHNICAL
DATA SHEETS

11:51

P - 1[®] PRESERVATIVE

DESCRIPTION

P-1 PRESERVATIVE is effective against a wide variety of mold fungi and bacteria for the preservation of glues and adhesives, polymer emulsions (coatings, PVA systems and rubber), thickeners (starch, cellulose, guar, etc.), paper, textiles, dyes, temporary sapstain control for fresh cut lumber, concrete additives / masonry, metalworking fluids, air filter oils, starches, clay slurries, slurries, printing inks, graphite, and hard surface treatment products (polishes, waxes and ceramic glazes).

The following guidance is given as an approximation for each use pattern, but field testing is recommended to achieve optimum effectiveness.

RECOMMENDED USES AND APPLICATIONS

GLUES AND ADHESIVES

P-1 PRESERVATIVE can be dissolved in suitable systems such as ethanol or 1,2-propanediol, or an alkaline homogeneous concentrate can be prepared and added to the substrate for the effective preservation of glues and adhesives. P-1 PRESERVATIVE can be directly mixed until homogeneous into to dry glues before they are concentrated.

Recommended Dosages:

7	
Bone Glues	0.10-0.15%
Skin Glues	0.15-0.40%
Leather Glues	0.15-0.40%
Fish Glues	0.15-0.40%
Plant Glues	0.10-0.25%
Gelatin-Based Blues	0.15-0.30%
Albumin-Containing Glues	0.20-0.40%
Starch-Based Glues (Liquid)	0.10-0.25%
Starch-Based Glues (Solid)	0.15-1.00%
Polymer Dispersion-Based Adhesives	0.10-0.30%
Dextrin-Based Adhesives	0.10-0.30%
Cellulose-Based Adhesives	0.10-0.30%
Casein-Containing Adhesives	0.20-0.40%

STAINS AND PAINTS

P-1 PRESERVATIVE is generally incorporated into pigments and fillers, but it can also be first dissolved in a suitable solvent or added directly to the pre-heated binder systems. For best results, the preservative should be homogeneously incorporated into the stain or paint. Discoloration may occur, and should be evaluated.

11:51

RECOMMENDED USES AND APPLICATIONS (CONT)

STAINS AND PAINTS (cont)

TO

Recommended Dosages:

Latex-Based Systems	0.15-0.25%
Casein-Systems	0.20-0.40%
Oil-Containing Systems	
Alkyd Resin-Based Systems	
Synthetic Resin Dispersions	0.10-0.25%

TEMPORARY SAPSTAIN CONTROL FOR FRESH CUT LUMBER AND OTHER LUMBER

P-1 PRESERVATIVE is converted to alkaline liquid and water dilutable formulations by adding sodium hydroxide or other bases or by preparing emulsion concentrates of P-1 PRESERVATIVE. Auxiliaries may also be added. For the temporary protection of freshly sawn lumber against staining and molding, P-1 PRESERVATIVE solutions as prepared above are applied to the freshly sawn lumber by either dipping or spraying.

Recommended Dosages:

Fresh Cut Lumber	1.00-4.00%
Fruit and Vegetable Containers	1.00-3,00%
Pallets	
Construction Woods	

DYES, PIGMENTS AND FILLER SUSPENSIONS

P-1 PRESERVATIVE should be evenly mixed with dispersing agents for processes where the temperature does not exceed 100°C. For all other processes, P-1 PRESERVATIVE is metered directly into the material to be preserved during the cooling step after the temperature falls below 50°C and homogeneously distributed by stirring.

Recommended Dosages:

Clay Slurries	0.05-0.20%
Kaolin Slurries	
Starch Slurries	
Other Filler Suspensions	
Calcium Carbonate Slurries	
Iron Oxide Pigment Slurries	0.05-0.20%
Carbon Black Pigment Slurries	
Titanium Dioxide Pigment Slurries	
Organic Dyes and Pigment Slurries	

POLYMER DISPERSIONS AND EMULSIONS

P-1 PRESERVATIVE is added immediately after the preparation of the polymer dispersion or emulsion during the cooling process. Losses of active ingredient caused by elevated temperatures should be avoided. For homogeneous distribution, P-1 PRESERVATIVE should be dissolved in suitable solvents such as ethanol, 1,2-propanediol, water dilutable alkaline solutions then added to emulsion concentrates or polymer dispersions.

RECOMMENDED USES AND APPLICATIONS (CONT)

POLYMER DISPERSIONS AND EMULSIONS (cont)

TO

LEATHER

P-1 PRESERVATIVE may be dissolved in five times the amount of alcohol and then stirred into pickle solutions. In the case of chrome leather, an effective protection against mold attack may be obtained without additional labor by stirring P-1 PRESERVATIVE together with soda into the basifying solution, since P-1 PRESERVATIVE dissolves freely in the alkaline solution.

The oil-soluble P-1 PRESERVATIVE can also be dissolved in the fat liquor and applied in the drum simultaneously. Finished vegetable-tanned leathers and chrome leathers may also be protected against mold attack by a treatment on both sides of the material with an 0.2-0.4% solution of P-1 PRESERVATIVE.

Recommended Dosages:

Pickle Solutions/Pickled Hides	0.15-0.25%
Chrome Leather (osiculated on pelt wt)	
Leather Pasting Adhesives	
Leather Pigment Finishes	

TEXTILES

P-1 PRESERVATIVE should be dissolved in suitable solvents such as ethanol, 1,2-propanediol, mineral spirits, or paraffins or can be converted to a water dilutable alkaline concentrate and applied directly to the textile normally by dipping or spraying. The preservative is affixed to the fiber with a suitable acid like acetic acid. For the preservation of auxiliaries, P-1 PRESERVATIVE is diluted in solvents as mentioned above or is homogeneously distributed with the aid of alkaline solutions or emulsion concentrates. Print thickeners are preserved by adding P-1 PRESERVATIVE directly to the emulsion or homogeneously metered directly into the dried product. Milling may be necessary in order to obtain consistent particulate sizes.

Recommended Dosages:

0.50-5.00%
0.50-5.00%
0.50-2.00%
0.50-2.50%
0.50-2.50%
0.50-2.00%
0.50-2.00%

RECOMMENDED USES AND APPLICATIONS (CONT)

TEXTILES (cont)

Recommended Dosages:

ΤO

Cotton Production	0.50-1.50%
Wood Production	
Fire Hoses	1.00~ 3,00%
Auxiliaries (Sizing Agents, Finishing Agents, Spin	nning Preparations.
Wetting Agents)	
- - ·	

LUBRICOOLANTS AND MINERAL OIL BASED PRODUCTS

P-1 PRESERVATIVE is dissolved in the necessary amount of mineral oil, auxiliary (for example dispersing agent), solvents such as Ethanol, Glycols or water dilutable alkaline concentrates, and added to form a homogeneous mixture in the substrates (metalworking fluids, etc.,) by stirring.

Recommended Dosages:

Mineral Oil-Based Pro	oduct	S
-----------------------	-------	---

Cooling Fluids - Concentrate	1.00-5.00%
Ready for Use	0.05-0.25%
Boring and Cutting Oils	
Fuel Oils	0.02-0.15%
Hydraulic Oils	

Non-Mineral Oil-Based Products

Cutting Oils	0.05-0.30%
Cooling Fluids	.0.05-0.30%

OTHER APPLICATIONS

Recommended Dosages:

Concrete and Masonry Additives	0.10-0.30%
Biopolymers (e.g. Xanthan)	0.05-0.20%
Fire Extinguishing Medium	
Photographic Gelatins	
Tints	
Cleaning Solutions	0.05-0.40%
Wax Emplsions	
Chemical Cleaning Solutions	
Concrete	
Caulking Materials	
Rubber Systems	
Ceramic Glazes	
Polishes	
Paper Auxiliaries	0.05-0.30%
Paper Siurcies	

ACTIVITY

MINIMUM INHIBITION CONCENTRATIONS OF O-PHENYLPHENOL, DETERMINED ON SPECIAL NUTRIENT MEDIA

;	Minimum Inhibition Concentration
Microorganisms	mg/liter
Bacillus punctatus	200
Bacillus subtilis	100
Bacterium vulgare	100
Escherichia coli	200
Leuconostoc mesenterioides	100
Pseudomonas aeruginosa	1500
Pseudomonas fluorescens	
Staphylococcu aureus	100
Candida albicans	100
Saccharomyces cerevisiae	200
Saccharomyces acidifaciens	200
Torula rubra	100
Aspergillus flavus Link	25
Asperillus niger	30
Chaetomium globosum	. 60
Cladosporium herbarum link and fries	κΛ
Cladosporium sphaerospennum	40
Coniophora cerebella	24
Lentinus tigrinus	
Penicillium citrinum	***************************************
Penicillium digitatum	
Penicillum fimicolosum QECD F 10	33
Penicillum italicum	·············/D
Polyporus versicolor	
Pullularia pullulans Fusey	
Trichiphyton pedis	۵۵
Trichoderma viride	75

TYPICAL. **PROPERTIES**

Physical FormFlakes, colorle	ess to light yellow with slight phenolic odor.
Composition2-Phe	mylphenol, active ingredient content 99.9%
Density	
Rulk Density	
Malanda 317. '-7.	0.60 – 0.65
TATOTECOTAT AA GIGUL	770
TATOLETING T OHILL	56 8000
Boiling Point	
Vanor Pressure	
Glach Doint	0.5 torr at 100°C
TIASH FULLE	13897

TO

Methanol	800
Acetone	
Ethanol	
Ether	
Benzene	
Pine Oil	
Propylene Glycol	300
Ethylene Glycol	
Carbon Tetrachloride	
NaOH, 10%	50
White Spirit	
Water	

PRECAUTIONS AND HANDLING

PESTICIDE STORAGE AND DISPOSAL

Pesticide Storage

Keep containers tightly and securely closed. Do not contaminate water, food or feed by storage and disposal. Store below 122°F (50°C). Protect from exposure to light and moisture. The product has a shelf life of at least two (2) years if protected from light and moisture. The flakes agglomerate if the storage temperature is too high. See Material Safety Data Sheet for additional instructions.

Pesticide Disposal

Pesticide wastes are acutely hazardous. Improper disposal of excess pesticide, spray mixture, or rinsate is a violation of Federal Law. If these wastes cannot be disposed of by use according to label instructions, contact your State Pesticide or Environmental Control Agency, or the Hazardous Waste Representative at the nearest EPA Regional Office for guidance.

Container Disposal

Triple rinse (or equivalent). Then offer for recycling or reconditioning, or puncture and dispose of in a sanitary landfill, or incineration, or, if allowed by state and local authorities by burning. If burned, stay out of smoke.

HEALTH AND SAFETY INFORMATION

Appropriate literature has been assembled which provides information pertaining to health and safety concerns that must be observed when handling Verichem products, appropriate industrial hygiene and other safety precautions recommended by their manufacturer should be followed. Before working with any product mentioned in this publication, you must read and become familiar with available information concerning its hazards, proper use and handling. This cannot be overemphasized. Information is available in several forms, such as Material Safety Data Sheets and Product Labels. Consult your Verichem representative or contact the Product Safety Manager.

F&DA

TO

175.105 (Adhesives) for use as a preservative in the manufacture of complying adhesives

176.210 (Defoaming agents used in the manufacture of paper and paperboard)

177.1632 (Poly(phenyleneterephthalamide)resins), for use as a fungicide, not to exceed 0.01 wt% based on the base polymer

177.2600 (Rubber articles intended for repeated use) for use as an antioxidant and antiozonant, total antioxidants and antiozonants not to exceed 5 wt% of rubber product

178.1010 (Sanitizing solutions) for use in sanitizing solutions specifically listed in 178.1010

AVAILABILITY

P-1 PRESERVATIVE is available in commercial quantities.

Shipping Weight55.9 pounds

NOTICE: This product is for industrial use only by persons having technical skill in formulating coatings, adhesives, elastomers, paints and the like. VERICHEM MAKES NO WARRANTY OF ANY KIND, EXPRESS OR IMPLIED, WITH RESPECT TO THE GOODS WHEN NOT USED IN STRICT ACCORDANCE WITH DIRECTIONS AND RECOMMENDATIONS GIVEN HEREWITH. Recommendations for the use of this product are based on laboratory tests that Verichem considers to be reliable. The performance of this product varies from application to application. Verichem is not responsible for results when the product is used with incompatible or contaminated raw materials; when the product is used under unsanitary conditions; or when manufacturing equipment is constructed of such raw materials as to inactivate this product. Under no circumstances will Verichem be liable for consequential damages or damages to anyone in excess of the purchase price of the product. User assumes all risks of use, storage and handling, not in strict accordance with the precautionary statements and storage recommendations of the labeling.

EFFECTIVE DATE: 1/1/98

REPLACES: ORIGINAL







MATERIAL SAFETY DATA TEMPEZ

MATERIAL SAFETY DATA SHEET

CAS NO: 6152-33-6

DATE: 7/21/99

SUPERSEDES: Original

1. CHEMICAL PRODUCT AND COMPANY IDENTIFICATION

PRODUCT NAME:

P-J PRESERVATIVE 20% Solution

SYNONYMS:

Sodium O-phenylphenate

CHEMICAL FAMILY:

Phenol

MOLECULAR FORMULA:

C12H oONa

MOLECULAR WEIGHT:

192

VERICHEM, 3499 Grand Avenue, Pittsburgh, PA, USA, 15225 (412-331-7299, 8:30 am to 5:00 pm; 412-521-6393 24 hours).

EMERGENCY PHONE: For any emergency involving spill, leak, fire, exposure, or

accident call CHEMTREC: 1-800-424-9300. Outside the USA and Canada call: 202-483-7616.

2. COMPOSITION/INFORMATION ON INGREDIENTS REGULATED COMPONENTS

	COMPONENT	CAS NUMBER	PERCENT	WORKER EXPOSURE	REFERENCE
. [Sodium Orthophenylphenate	132-27-4	20%	None Established	

3. HAZARDS IDENTIFICATION

EMERGENCY OVERVIEW

APPEARANCE AND ODOR: Clear Pale to Dark Yellow Solution

EFFECTS OF OVEREXPOSURE: Na OPP has been shown to cause bladder

tumors in mice fed high doses.

PHYSICAL/CHEMICAL HAZARDS: Not regulated for transportation.

FIRST AID MEASURES (IN CASE OF CONTACT) 4.

EYE: Flush with plenty of water for at least 15 minutes. Consult a physician.

SKIN: Flush with plenty of water for at least 15 minutes, wash with soap and water. If

irritation persists, consult a physician.

INGESTION: Do not induce vomiting. Call a physician immediately. Never give

anything by mouth to an unconscious person.

INHALATION: No toxic effects expected.

5. FIRE FIGHTING MEASURES

FLAMMABLE PROPERTIES

FLASH POINT:

FLAMMABLE LIMITS (% BY VOLUME): **AUTOIGNITION TEMPERATURE:**

DECOMPOSITION TEMPERATURE:

Not flammable.

Not applicable. Not applicable.

Not applicable

EXTINGUISHING MEDIA AND FIRE FIGHTING INSTRUCTION

In case of fire use water spray, alcohol foam, carbon dioxide, or dry chemicals. Exercise caution when fighting fires involving chemical substances. Respiratory protection is required. Burning will produce toxic fumes. Containers exposed to heat will result in pressure increase in containers. Cool hot containers with cool water.

ACCIDENTAL RELEASE MEASURES б.

STEPS TO BE TAKEN IN CASE MATERIAL IS RELEASED OR SPILLED

Immediately dike material to keep away from water sources. Absorb with vermiculite or other suitable absorbent. Discard in accordance with Federal, State, and local regulations. Residual amounts left on surface maybe cleaned with soap and water. Rinse residuals to municipal treatment plant.

. 7. HANDLING AND STORAGE

HANDLING: Avoid breathing vapors. Avoid getting in eyes or on skin. Wash thoroughly after handling,

STORAGE: Store in a cool place.

8. PERSONAL PROTECTION

EVE PROTECTION REQUIREMENTS: Chemical safety glasses or face shield. SKIN PROTECTION REQUIREMENTS: Rubber gloves, long sleeve shirts and pants. VENTILATION REQUIREMENTS: Local and general exhaust ventilation to control levels of

RESPIRATOR REQUIREMENTS: Air purifying particulate / organic vapor cartridge respirator.

. 9. PHYSICAL AND CHEMICAL PROPERTIES

APPEARANCE AND ODOR:

BOILING POINT:

MELTING POINT: VAPOR PRESSURE:

SPECIFIC GRAVITY: VAPOR DENSITY:

PERCENT VOLATILE (BY WEIGHT): рĦ:

SATURATION IN AIR

(PERCENT BY VOLUME):

EVAPORATION RATE: SOLUBILITY IN WATER: Clear light yellow to dark brown.

Similar to water. Not applicable.

Unknown. 1.068 ± 0.010 Unknown.

80% 12 - 12.5

Unknown. Unknown,

Complete.

STABILITY AND REACTIVITY 10.

STABILITY (CONDITIONS TO AVOID):

POLYMERIZATION (CONDITIONS TO AVOID):

INCOMPATIBLE MATERIALS:

HAZARDOUS DECOMPOSITION PRODUCTS:

produce CO, CO2 and other toxic furnes.

Stable.

Not applicable. Oxidizing material.

Thermal decomposition may

PRECAUTIONARY STATEMENTS

HAZARDS TO HUMANS AND DOMESTIC ANIMALS

DANGER

Corrosive, May be falal If inhaled.

Harmful if swaltowed or absorbed through skin.

Causes skin burns and irreversible eve damage. Do not get in eyes, on skin or on clothing. Do not breathe vapor or mist. Weer protective eyewear (goggles, faceshield or safety glasses). Wear protective clothing and gloves. Wear a mask or pesticide respirator approved by the National Institute for Occupational Safety and Health. Wash thoroughly with soap and water after handling. Remove contaminated clothing and wash clothing before rouse.

ENVIRONMENTAL HAZARDS

This pesticide is toxic to fish. Do not discharge effluent containing this product into takes, streams, ponds, estuaries, oceans or other waters unless in accordance with the requirements of a National Pollutant Discharge Elimination System (NPDES) permit and the permitting authority has been notified in writing prior to discharge. Do not discharge effluent containing this product to sever systems without previously notifying the local sewage treatment plant authority. For guidance contact your local State Water Board or Regional Office of the EPA.

PHYSICAL OR CHEMICAL HAZARDS

Corrosive to skin and eyes.

STORAGE AND DISPOSAL

PESTICIDE STORAGE: Keep containers tightly and securely closed. Do not contaminate water, food or feed by storage. Store below 122°F (50°C), protect from exposure to light and moisture. The product has unlimited shell life if protected from light and moisture. See Verichem's Material Salety Data Sheet for additional instructions.

PESTICIDE DISPOSAL: Do not contaminate water, food or feed by storage and disposal. Pesticide wastes are acutely hazardous. Improper disposal of excess pesticide spray mixture, or rinsate is a violation of Federal Law, If these wastes cannot be disposed of by use according to label instructions, contact your State Pesticide or Environmental Control Agency, or the Hazardous Waste Representative at the nearest EPA Regional Office for guidance.

CONTAINER DISPOSAL: Triple rinse (or equivalent). Then offer for recycling or reconditioning, or puncture and dispose of in a sanitary landill, or by incineration, or, if allowed by state and local authorities, by burning. If burned, stay out of smoke.

NONWARRANTY: Seller (whether manufacturer, distributor or other) makes no warranty express or implied, including the implied warranty of merchantability, regarding this product except the composition as set forth in the ingredient statement herein, and shall not be liable for special or consequential damages, the exclusive remedy being replacement of the product. Buyer or user assumes all risk of possession, handling or use of this material when such use and/or handling is contrary to label instructions.

P-1® PRESERVATIVE SOLUTION

FOR THE PRESERVATION OF GLUES AND ADHESIVES, POLYMER EMULSIONS, (COATING, PVA SYSTEMS AND RUBBER), THICKENERS (STARCH, CELLULOSE, GUAR, ETC.), LEATHER, PAPER, TEXTILES, DYES, TEMPORARY SAPSTAIN CONTROL FOR FRESH CUT LUMBER, CONCRETE ADDITIVES/MASONRY, METALWORKING FLUIDS, AIR FILTER OILS, STARCHES, CLAY SLURRIES, SLURRIES, PRINTING INKS, GRAPHITE, AND HARD SURFACE TREATMENT PRODUCTS (POLISHES, WAXES AND CERAMIC GLAZES).

> Active Ingredient: TOTAL 100.0%

• KEEP OUT OF REACH OF CHILDREN • DANGER

If Swallowed: Call a doctor or get medical attention. Do not induce vomiting or give anything by mouth to an unconcious person. Drink promptly a large quantity of milk, egg whites, gelatin solution, or it not available, drink large quantities of water. Avoid alcohol.

If Inhaled: Remove victim to fresh air. If not breathing, give artificial respiration, preferably mouth to mouth. Get medical attention.

If on Skin: Wash with plenty of soap and water. Get medical attention.

If in Eves: Hold evelids open and flush with a steady, gentle stream of water for 15 minutes. Get medical attention.

See side panel for additional precautionary statements.

LOT NO.	<u></u>	
EPA Registration No. 67869-24	EPA Est. No.:	67869-PA-0
LB, NET WT.		ko
Made in U.S	S.A.	

SOLD BY:



VERICHEM INC.

3499 Grand Avenue Pittsburgh, PA 15225

Final Approved 5-5/99

DIRECTIONS FOR USE

It is a violation of federal law to use P-1© Preservative Solution is effectipreservation of glues and adhesive: thickeners (starch, cellulose, guar, fresh out lumber, concrete additives slurries, slumies, printing inks, grept and ceramic glazes).

The following quidance is given as recommended to achieve optimum : ADHESIVES AND GLUES

P-1® Preservative Solutions can be 1,2-propanadol, or an alkaline hom substrate for the effective preservat: be directly mixed until homogeneous Recommended Dosages:

Albumin-Containing Glues... Bosa Gluss..... Casein-Containing Adhesive Collinose-Based Adhesivas Dextrin-Based Adhesives ... Fish Gives Gelatin-Based Glues Leather Gines Plant Glues..... Polymer Despersion-Based Skin Glues Starch-Based Glues (Liquit Starch-Based Glues (Solid)

HOUSEHOLD PRODUCTS & (

Recommended Dosages: Biopolymers (e.g. Xanthan) Cautking Materials..... Ceramic Olazos..... Chemical Cleaning Solution Cleaning Solutions..... Concrete Concrete and Masonry Add Fire Extinguishing Modium. Photographic Goldins...... Plasters..... Polishes..... Rubber Systems Tints...... Wax Ernulsions..... Paper Auxiliaries Paper Sturries.....

LEATHER TANNING

P-1© Preservative Solution may be a into picke solutions in the case of may be obtained without additional soda into the basilying solution. Fin also be protected against mold affact - 2.26% solution as P-1® Preservat

Recommended Dosages:

Chrome leafter (calculated Loather pasing achesives. Leather pigment finishes Pickle solutions and pickled

METAL WORKING FLUIDS, LUB! Recommended Dosages:

Mineral Oil-Based Products Boring and Cutting Olls Cooling Fluids - Concentrat Ready for Jae -Fuel Olls HydrauAic Oirs Non-Mineral Of-Based Produ

Cooling Fluids..... Culling O's

FOR USE

of lederal law to use this product in a manner inconsistent with its labeling. ive Solution is effective against a wide variety of model tengt and bacteria for the glues and adhesives, polymer emulsions (coallings, PVA systems and rubber), ich, cellulose, guar, otc.), paper, loxilles, oyes, temporery expession control for ar, concrete additives/mesonry, molalizations fixities, air litter of s, starches, clay s, printing links, graphile, and hard surface treatment products (polishes, waxes a288).

zuidance is given as an approximation for each uso patiem, but hold testing is to achieve optimum effectiveness.

AND GLUES

the Solutions can be dissolved in suitable solvent systems such as elizabli or of an alkaline homograpous concentrate can be propared and added to the na offective preservation of globs and anhesives. P-1® Preservative Solution can ed until homogeneous into to dry gives before they are concentrated.

nded Dosages:	
n-Coarsining Glues	1.1 - 2.3%
Nes	0.6 - 0.9%
Costining Adhesives	1.1 - 2.3%
Containing Adhesives	0.6 - 1.7%
20-D89-60 ART 10-214-0-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1	0.6 - 1.7%
·Based Adhesives	0.2 - 2.3%
ves	
-Based Gives	0.9 - 1.7%
r GCues	9.2 - 2.3%
iles	0.6 - 1.4%
if Dispersion-Based Aultrasives	0.6 - 1.7%
UBS	0.9 - 2.3%
Based Glues (Liquid)	0.6 - 1.4%
	0.9 - 5.7%
Based Glues (Solid)	G.3 - 3.1 /P
D PRODUCTS & CONSTRUCTION PRODUCTS	
nded Dosages:	
mers (e.g. Xanthan)	0.3 - 1.1%
ig Moterials	0.8 - 2.8%
ig rades as a second se	0.6 - 2.8%
c G'azos al Cleaning Solutions	0.1 - 0.8%
AI CEANING SOURCES	0.3 - 2.3%
ig Solutions	0.3 - 2.3%

le______0.6 - 2.8%

0.6 - 2.6%

8.3-1.1%

Sturries 0.4 - 0.5% ANNING

tive Solution may be dissolved in five times the amount of alcohol and then stirred nions. In the case of chrome leather, an effective protection against mold attack ed without additional labor by stirring P-1® Preservative Schrich together with pasitying solution. Finished vegetable-fanted leathers and chicme leathers may ted against mold affack by a trentment on both sides of the material with an 1.13 in as P-18 Preservative Solution.

nded Dosages:

a leafrer (calculated on pelt weight)	0.6 - 1.4%
r nasting achiesives	0.3 - 0.6%
r corment finishes	1.2 - 2.3%
solutions and pickled fildes	0.9 - 1.4%

KING FLUIDS, LUBRICOOLANTS AND MINERAL OIL BASED PRODUCTS aded Dosages:

I-Based Products

and Cutting Oils	0.3 - 1.1%
Fluids - Concentrate	5.7 - 28.3%
ty for .120	U.3 + 1.4%
5	0.1 5:8%
To Ois	03 - 1.4%
at OR-Based Products	
	0.2.17%

PAINTS, COATINGS & STAINS

P-10 Preservatives Solution is generally incorporated directly into pigments and filters, but it can also be first dissolved in a sulfable solvent or added directly to the pre-hasted binder systems. For basi results, the preservative should be homogeneously incorporated into the stain or paint. Discoloration may occur, and should be evaluated.

Recommended Ocsages:

Alkyd Resin-Based Systems	1.4 - 2.8%
Casein-Systems	1.1 - 2.3%
Latex-Based Systems	0.9 - 1.4%
Oil-Containing Systems	1.4 - 2.8%
Synthelic Resin Dispersions	0.6 - 1.4%

PIGMENTS, DYES AND FILLER SUSPENSIONS

P-16 Preservative Solution should be everify imbout with departing agents for processes where the temperature does not exceed 100°C. For all other processes, P-10 Preservative Solution Is metered directly into the material to be preserved during the cooting step after the temperature falls below 50°C and homogeneously distributed by stirring.

Recommended Dosages:

Calcium Carbonate Sturries	0.3 - 1.1%
Carbon Black Figment Sluries	0.3 - 1.1%
Clay Signies	0.3 - 1.1%
Iron Oxide Pigment Sturies	0.3 - 1.1%
Kaolin Sturiles	0,3 - 1,1%
Organic Dyes and Plyment Sturies	0.3 - 1.1%
Other Filter Suspensions	0.3 - 1.7%
Start Studes	0.3 - 1.7%

POLYMER DISPERSIONS AND EMULSIONS

P-1@ Preservative Solution is added immediately after the preparation of the polymer dispersion or enablishin during the ocoling process. Losses of active ingredient caused by elevated temperatures should be avoided.

Recommended Dosages (Polymer Dispersions and Emulsions):

Acrylio-Systems	0.3 - 1.7%
Other Polymer Dispersions	0.3 - 1.7%
Polyester-Based Systems	0.9 - 1.7%
Polyurethane-Based Systems	0.3 - 1.7%
Polyvinylacetate (PVA) Systems	0.3 - 1.7%
Styreno Butadiene Systems	0.3 - 1.7%
Manufic and in Record Systems	0.3 - 1.7%

TEXTILES

P-1@ Preservative Solution should be dissolved in suitable scheda such as albertal, 12proponedial, mineral spirits, or paraffirs or can be convented to a water dilutable alkaline concentrate and applied directly to the textile normally by dipping or spraying. The preservative is affixed to the fiber with a suitable acid like acetic acid. For the preservation of auxiliaries, P. 100 Preservative Solution is diluted in solvents added directly to textile. Print thickenars are preserved by ailding P-K9 Preservative Solution directly to the smulsion or homogeneously metered directly into the direct product. Milling may be necessary in order to obtain consistent particulate sizes.

Recommended Dosages:

Auxillation (String Agents, Finishing Agents, Spinring Preparations, Wetting Agents) 0.1 - 0.3% 28.263%

AMDITOS EDU IZEDEURIOS	L.O LLW
Caroeis	2.8 - 11.3%
Cotton Production	2.8 - 6.5%
Fe15	2.8 - 14.1%
Rio Hoses	2,8 - 17.0%
Geotaxiles	2:8 - 28:3%
Jule, Sis	2.8 - 14.1%
Shower Curiains	2.8 - 11.3%
Uphostery	2.8 - 11.3%
Vicol Production	2.8 - 8.5%

WOOD PRESERVATION

For the temporary protection of freshly sawn lumber against starting and motiting, P-10 Preservative Solutions are applied to the treetry saven leaster by either display or spraying.

Recommended Desages: . .

Maintifferines resulter.	
	5.7 - 22.6%
Fresh Cut Lumber	5.7 - 22.6%
Fruit and Vegetable Containers	67 - 17 0%
Full six Asiastos commitais	E 7 17 000
Palints	2V 1 - 1 (TA 39

NTP CHEMICAL REPOSITORY (RADIAN CORPORATION, AUGUST 29, 1991) O-PHENYLPHENOL

```
-IDENTIFIERS
========
*CATALOG ID NUMBER: 000575
*CAS NUMBER: 90-43-7
*BASE CHEMICAL NAME: PHENYLPHENOL, O-
*PRIMARY NAME: O-PHENYLPHENOL
*CHEMICAL FORMULA: C12H100
*STRUCTURAL FORMULA:
*WLN: QR BR
*SYNONYMS:
 (1,1'-BIPHENYL)-2-OL
 2-BIPHENYLOL
 O-BIPHENYLOL
 2-DIPHENYLOL
 2-HYDROXYBIPHENYL
 O-DIPHENYLOL
 DOWICIDE 1
 DOWCIDE 1
 O-HYDROXYBIPHENYL
 O-HYDROXYDIPHENYL
 NCI-C50351
 2-PHENYLPHENOL
 ORTHOHYDROXYDIPBENYL
 ORTHOPHENYLPHENOL
 ORTHOXENOL
 TORSITE
 USAF EK-2219
 O-XENOL
  2-HYDROXYDIPHENYL
 -PHYSICAL CHEMICAL DATA
  Literature: Light purple colored crystals.
 *PHYSICAL DESCRIPTIONS:
                          Repository: Light lavender crystalline solid.
 *MOLECULAR WEIGHT: 170.21
```

*SPECIFIC GRAVITY: 1.213 @ 25/4 C

*DENSITY: Not available

*MP (DEG C): 55.5 - 57.5 C

*BP (DEG C): 275 C

*SOLUBILITIES:

WATER: <0.1 mg/mL @ 20.5 C (RAD)

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DMSO : >=100 mg/mL @ 22 C (RAD)

95% ETHANOL : >=100 mg/mL @ 22 C (RAD)

METHANOL : Not available

ACETONE : >=100 mg/mL @ 22 C (RAD)

TOLUENE : Not available

OTHER SOLVENTS:

Ligroin: Soluble

Fixed Alkali Hydroxide Solution: Soluble

ETHER : Soluble BENZENE: Soluble

*VOLATILITY : Vapor pressure: 1 mm Hg @ 100 C

*FLAMMABILITY(FLASH POINT):

This chemical has a flash point of 124 C (255 F). It is combustible. Fires involving this chemical can be controlled with a dry chemical, carbon dioxide, foam or Halon extinguisher.

*UEL: Not available LEL: Not available

*REACTIVITY: This compound can react with oxidizing materials.

*STABILITY: This chemical is stable under normal laboratory conditions.

*OTHER PHYSICAL DATA:

Boiling point: 145 C @ 14 mm Hg

Volatile in steam

-TOXICITY

========

*NIOSH REGISTRY NUMBER: DV5775000

*TOXICITY:

typ.	dose	mode	specie	amount	unit	other
	LD50	orl	rat	2000	mg/kg	
	LD50	unk	rat	2700	mg/kg	
	LD50	orl	mus	900	mg/kg	
	LD50	ipr	mus	50	mg/kg	
	LD50	orl	rat	2.48	g/kg	

*AQTX/TLM96: Not available

*SAX TOXICITY EVALUATION: THR: MODERATE via oral route.

*CARCINOGENICITY:

Review: IARC Cancer Review: Animal Inadequate Evidence

IARC: Not classifiable as a human carcinogen (Group 3) [610]

Status: NCI Carcinogenesis Bioassay (Dermal); No Evidence: Male and Female

Mouse [620]

*MUTAGENICITY:

Mutagenic Data: mmo-sat 100 ug/plate dnr-esc 1 mg/disc

cyt-hmn:fbr 200 ug/L

*TERATOGENICITY:

Reproductive Effects Data:

6 gm/kg (6-15D preq) orl-rat TDLo 7 gm/kg (6-15D preg) orl-rat TDLo orl-mus 13050 mg/kg (7-15D preg) TDLo orl-mus 13050 mg/kg (7-15D preg) TDLo

*STANDARDS, REGULATIONS & RECOMMENDATIONS:

osha: None ACGIH: None

NIOSH Criteria Document: None NFPA Hazard Rating: Health (H): 1 Flammability (F): 1 Reactivity (R): 0

H1: Materials only slightly hazardous to health (see NFPA for details). F1: Materials that must be preheated before ignition can occur (see NFPA for details).

RO: Materials which are normally stable even under fire exposure conditions and which are not reactive with water (see NFPA for details).

*OTHER TOXICITY DATA:

Skin and Eye Irritation Data: skn-rbt 25 mg/24H SEV 50 ug/24H SEV eye-rbt

Status: Reported in EPA TSCA Inventory, 1980

Meets criteria for proposed OSHA Medical Records Rule

-OTHER DATA (Regulatory) -----

*PROPER SHIPPING NAME (IATA): Other regulated substances

*UN/ID NUMBER: ID8027

SUBSIDIARY RISK: None PACKING GROUP: None *HAZARD CLASS: 9

*LABELS REQUIRED: Miscellaneous

MAXIMUM QUANTITY: No limit *PACKAGING: PASSENGER: PKG. INSTR.: 906 MAXIMUM QUANTITY: No limit CARGO : PKG. INSTR.: 906

*SPECIAL PROVISIONS: None

Intermediate for dyes and resins, rubber chemicals, fungicide, germicide, preservative, food packaging. Disinfectant and fungicide for impregnation of fruit wrappers and disinfection of seed boxes. Applied during dormant period to control apple canker. Reagent for the determination of trioses. Household disinfectant; dishwashing formulations.

*COMMENTS: Not available

-HANDLING PROCEDURES _____

*ACUTE/CHRONIC HAZARDS: This chemical is a skin and eye irritant. Corneal damage is possible upon eye contact.

First check the victim for contact lenses and remove if present. Flush victim's eyes with water or normal saline solution for 20 to 30 minutes while simultaneously calling a hospital or poison control center.

Do not put any ointments, oils, or medication in the victim's eyes without specific instructions from a physician.

IMMEDIATELY transport the victim after flushing eyes to a hospital even if no symptoms (such as redness or irritation) develop.

*MINIMUM PROTECTIVE CLOTHING:

If Tyvek-type disposable protective clothing is not worn during handling of this chemical, wear disposable Tyvek-type sleeves taped to your gloves.

*RECOMMENDED GLOVE MATERIALS: Not available

*RECOMMENDED RESPIRATOR:

Where the neat test chemical is weighed and diluted, wear a NIOSH-approved half face respirator equipped with an organic vapor/acid gas cartridge (specific for organic vapors, HCl, acid gas and SO2) with a dust/mist filter.

*OTHER: Not available

*STORAGE PRECAUTIONS:

You should store this chemical under refrigerated temperatures, and keep it away from oxidizing materials.

*SPILLS AND LEAKAGE:

Should a spill occur while you are handling this chemical, you should dampen the solid spill material with alcohol, then transfer the dampened material to a suitable container. Use absorbent paper dampened with alcohol to pick up any remaining material. Seal the absorbent paper, and any of your clothes, which may be contaminated, in a vapor-tight plastic bag for eventual disposal. Solvent wash all contaminated surfaces with alcohol followed by washing with a strong soap and water solution. Do not reenter the contaminate area until the Safety Officer (or other responsible person) has verified that the area has been properly cleaned.

*DISPOSAL AND WASTE TREATMENT:

You should dispose of all waste and contaminated materials associated with this chemical as specified by existing local, state and federal regulations concerning hazardous waste disposal. It is suggested that your contaminated materials should be destroyed by incineration in a special, high temperature (>2000 degrees F), chemical incinerator facility.

-EMERGENCY PROCEDURES

*SKIN CONTACT:

IMMEDIATELY flood affected skin with water while removing and isolating all contaminated clothing. Gently wash all affected skin areas thoroughly with soap and water.

IMMEDIATELY call a hospital or poison control center even if no symptoms (such as redness or irritation) develop.

IMMEDIATELY transport the victim to a hospital for treatment after washing the affected areas.

*INHALATION:

IMMEDIATELY leave the contaminated area; take deep breaths of fresh air. If symptoms (such as wheezing, coughing, shortness of breath, or burning in

the mouth, throat, or chest) develop, call a physician and be prepared to transport the victim to a hospital.

Provide proper respiratory protection to rescuers entering an unknown atmosphere. Whenever possible, Self-Contained Breathing Apparatus (SCBA) should be used; if not available, use a level of protection greater than or equal to that advised under Respirator Recommendation,

*EYE CONTACT:

Check for contact lenses and remove them at once if present. You should then immediately flush eyes with water from any source for 15 minutes. Do not use oil or ointment in eyes. Arrange immediate transportation to a medical facility.

*INGESTION:

DO NOT INDUCE VOMITING. Phenols are very toxic poisons AND corrosive and irritating, so that inducing vomiting may make medical problems worse. IMMEDIATELY call a hospital or poison control center and locate activated charcoal, egg whites, or milk in case the medical advisor recommends administering one of them.

If advice from a physician is not readily available and the victim is conscious and not convulsing, give the victim a glass of activated charcoal slurry in water or, if this is not available, a glass of milk, or beaten egg whites and IMMEDIATELY transport victim to a hospital.

If the victim is convulsing or unconscious, do not give anything by mouth, assure that the victim's airway is open and lay the victim on his/her side with the head lower than the body. DO NOT INDUCE VOMITING. IMMEDIATELY transport the victim to a hospital.

*SYMPTOMS:

Symptoms of exposure to this compound include eye irritation with possible corneal injury (necrosis); paleness, cyanosis, weakness, sweating, headache, diarrhea, nausea, vomiting, diarrhea, dizziness, fainting, dark urine, central nervous system depression, and deoxyribonuclease inhibition. Chronic exposure may include irritation and lesions of the respiratory system.

*FIREFIGHTING: Not available

-SOURCES ======

*SOURCES:

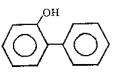
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- [610] Clansky, Kenneth B., Ed. Suspect Chemicals Sourcebook: A Guide to Industrial Chemicals Covered Under Major Federal Regulatory and Advisory Programs. Roytech Publications, Inc. Burlingame, CA. 1990. Update, p. xxiii.
- [620] United States National Toxicology Program. Chemical Status Report. NTP Chemtrack System. Research Triangle Park, NC. November 6, 1990. Listed.

CAS Registry Number: 90-43-7

Molecular Formula: C12H10O



90

2-Phenylphenol consists of white, flaky, crystals which have a mild, distinct odor. It is soluble in fixed alkali hydroxide solutions and most organic solvents, and is insoluble in water (Merck, 1989). 2-Phenylphenol emits acrid smoke and irritating fumes when it is heated to decomposition (Sax, 1989).

Physical Properties of 2-Phenylphenol

Synonyms: 2-biphenylol; o-biphenylol; o-phenylphenol; o-hydroxydiphenyl; orthoxenol; 2-hydroxydiphenyl

170.20 Molecular Weight: 280 - 284 °C **Boiling Point:**

55.5 - 57.5 °C Melting Point:

1.213 at 25/4 °C (water = 1) Density/Specific Gravity: 1 mm Hg at 100 °C

Vapor Pressure: 1 ppm = 6.96 mg/m^3 Conversion Factor:

(HSDB, 1993; Merck, 1989)

SOURCES AND EMISSIONS

A. Sources

- 2-Phenylphenol is used as a preservative in water-oil emulsions, in vegetable wax in dishwashing formulations, and in paper used for food storage. It is also used in the rubber industry and as a laboratory reagent (HSDB, 1993).
- 2-Phenylphenol (ortho-phenylphenol) is registered as a disinfectant, bactericide, and virucide. It is registered for agricultural use in disinfecting fruits, vegetables, and eggs. 2-Phenylphenol is registered for use as a general surface disinfectant in hospitals, nursing homes, veterinary hospitals, poultry farms, dairy farms, commercial laundries, barbershops, and food processing plants. It is also used to sterilize hospital and veterinary equipment (DPR, 1996).

Toxic Air Contaminant Identification List Summaries - ARB/SSD/SES September 1997

The licensing and regulation of pesticides for sale and use in California are the responsibility of the Department of Pesticide Regulation (DPR). Information presented in this fact sheet regarding the permitted pesticidal uses of 2-phenylphenol has been collected from pesticide labels registered for use in California and from DPR's pesticide databases. This information reflects pesticide use and permitted uses in California as of October 15, 1996. For further information regarding the pesticidal uses of this compound, please contact the Pesticide Registration Branch of DPR (DPR, 1996).

The primary sources of 2-phenylphenol emissions in California reported in the United States Environmental Protection Agency's (U.S. EPA) 1995 Toxics Release Inventory (TRI) Public Data Release Report were the chemical and allied products industries (U.S. EPA, 1996b).

B. Emissions

In California, approximately 10 pounds of 2-phenylphenol emissions were reported in the U.S. EPA 1995 TRI Public Data Release Report (U.S. EPA, 1996b).

- C. Natural Occurrence
- 2-Phenylphenol is not known to occur as a natural product (HSDB, 1993).

AMBIENT CONCENTRATIONS

No Air Resources Board data exist for ambient measurements of 2-phenylphenol.

INDOOR SOURCES AND CONCENTRATIONS

2-Phenylphenol is a common ingredient used in certain disinfectant products. In a U.S. EPA non-occupational pesticide exposure study, 32 household pesticides were measured in homes in three cities over several seasons from 1986-1988 (Immerman and Schaum, 1990). The indoor concentrations of 2-phenylphenol were moderately elevated. The mean indoor concentrations of 2-phenylphenol for homes in Springfield/Chicopee, Massachusetts for spring and winter were 22.8 and 44.5 nanograms per cubic meter (ng/m³), respectively. The mean indoor concentrations for homes in Jacksonville, Florida over three seasons ranged from 59.0 to 96.0 ng/m³. In contrast, outdoor concentrations of 2-phenylphenol were much lower and ranged from below 0.05 to 1.6 ng/m³.

ATMOSPHERIC PERSISTENCE

Based on its vapor pressure, 2-phenylphenol can be expected to exist primarily in the gas phase in the ambient atmosphere, although a small percentage may be associated with the particulate phase. The calculated half-life and lifetime of 2-phenylphenol due to reaction with the hydroxyl radical are 10 hours and 14 hours, respectively (Atkinson, 1995). It may also degrade

Toxic Air Contaminant Identification List Summaries - ARB/SSD/SES September 1997

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2-Phenylphenol

very rapidly in nighttime air by the reaction with nitrate radicals. The particulate phase of 2-phenylphenol may be physically removed from air via wet and dry deposition (HSDB, 1993).

AB 2588 RISK ASSESSMENT INFORMATION

2-phenylphenol emissions are not reported from stationary sources in California under the AB 2588 program. It is also not listed in the California Air Pollution Control Officers Association Air Toxics "Hot Spots" Program Revised 1992 Risk Assessment Guidelines as having health values (cancer or non-cancer) for use in risk assessments (CAPCOA, 1993).

HEALTH EFFECTS

Probable routes of human exposure to 2-phenylphenol are inhalation, and dermal contact.

Non-Cancer: 2-Phenylphenol or the sodium salt is irritating to the eyes, skin, and respiratory tract (Clayton and Clayton, 1981; HSDB, 1995). The U.S. EPA has not established a Reference Concentration (RfC) or an oral Reference Dose (RfD) for 2-phenylphenol (U.S. EPA, 1995a). It is fetotoxic in rodents, but it is not teratogenic in mice or rats (HSDB, 1995; Reprotox, 1995).

Cancer: The sodium salt of 2-phenylphenol causes bladder cancer in rats (HSDB, 1995). The U.S. EPA has not evaluated 2-phenylphenol for its carcinogenicity (U.S. EPA, 1995a). The International Agency for Research on Cancer has classified 2-phenylphenol (ortho-phenylphenol) in Group 3: Unclassifiable as to its carcinogenicity, based on no data in humans and inadequate data in experimental animals (IARC, 1987a).

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O-PHENYLPHENOL

CASNO: 90-43-7

FORMULA: C12-H10-O

SYNONYMS/COMMON NAMES

• (1,1'-BIPHENYL)-2-OL (9CI)

KNOWN USES:

• FUNGICIDE. GERMICIDE. HOUSEHOLD DISINFECTANT. PRESERVATIVE IN WATER-OIL EMULSIONS. DISH-WASHING FORMULATIONS.RUBBER INDUSTRY. PRODUCED FROM COAL TARS (TDB)

CHEMICAL HEALTH AND SAFETY INFORMATION:

SHORT-TERM TOXICITY

14-DAY (Topical) (C50351), COMPLETED
 MICE:SWISS

LONG-TERM CARCINOGENICITY

- 2-YEAR (Topical) (C50351)
 - o TR-301 (NTIS # PB86-217239) (PEER REVIEW 03/85)
 - o MICE:SWISS
 - o CARCINOGENESIS RESULTS (see RESULTS definitions)
 - MALE MICE = NO EVIDENCE
 - FEMALE MICE = NO EVIDENCE
 - DOSE: 55.5 MG/0.1 ML/50 PER GROUP

GENETIC TOXICOLOGY

- IN VITRO CYTOGENETICS
 - NEGATIVE (CHROMOSOME ABERRATIONS)
 - WEAKLY POSITIVE (SISTER CHROMATID EXCHANGES)
- DROSOPHILA (SEX-LINKED RECESSIVE LETHAL/RECIPROCAL TRANSLOCATION)
 - o NEGATIVE
 - CITATION Woodruff, R.C., Mason, J.M., Valencia, R., and Zimmering, S. Chemical mutagenesis testing in Drosophila: V. Results of 53 coded compounds tested for the National Toxicology Program. Environ. Mutagen. 7 (1985): 677-702.
- MOUSE LYMPHOMA
 - o POSITIVE
- MOUSE LYMPHOMA
 - o POSITIVE
- SALMONELLA

- NEGATIVE
- SALMONELLA
 - WEAKLY POSITIVE
 - o CITATION Haworth, S., Lawlor, T., Mortelmans, K., Speck, W., and Zeiger, E. Salmonella mutagenicity test results for 250 chemicals. Environ. Mutagen. 5(Suppl 1) (1983): 3-142.

Last Updated 11/13/00

2-Hydroxybiphenyl



• Formula: C₁₂H₁₀O

• MW: 170,21

SMILES String: Oc1ccccc1c2cccc2

Synonyms: o-Hydroxybiphenyl; o-Hydroxydiphenyl; 2-Hydroxydiphenyl; 2-Phenylphenol; Torsite; Orthoxenol;
 2-Biphenylol; o-Xonal; 1,1'-Biphenyl-2-ol; (1,1-Biphenyl)-2-ol; orthohydroxydipbenyl; biphenyl-2-ol;
 Biphenylol; Dowicide 1; Hydroxdiphenyl

CAS Reg. 90-43-7

ChemFinder Entry

• UM-BBD reactions whose product is 2-Hydroxybiphenyl

o 2'-Hydroxybiphenyl-2-sulfinate ----> 2-Hydroxybiphenyl (reacID# r0236)

[Dibenzothiophene Desulfurization] [BBD Main Menu]

Page Author(s): Margie Romine
March 15, 1998 6:14:11 PM CST BBDMaster@email.labmed.umn.edu

This is the UM-BBD compound page for compID# c0269. It was generated on December 6, 2000 3:15:39 PM CST.

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The UM-BBD Biphenyl Pathway

- Reactions in the Biphenyl Pathway
 - o Biphenyl ----> cis-2,3-Dihydro-2,3-dihydroxybiphenyl (reacID# r0388)
 - o cis-2,3-Dihydro-2,3-dihydroxybiphenyl ----> 2,3-Dihydroxybiphenyl (reacID# r0389)
 - o 2,3-Dihydroxybiphenyl ----> 2-Hydroxy-6-oxo-6-phenylhexa-2,4-dienoate (reacID# r0390)
 - o 2-Hydroxy-6-oxo-6-phenylhexa-2,4-dienoate ----> Benzoate + cis-2-Hydroxypenta-2,4-dienoate (reacID# r0391)
- · Compounds in the Biphenyl Pathway
 - o Benzoate (compID# c0121)
 - o Biphenyl (compID# c0371)
 - o cis-2,3-Dihydro-2,3-dihydroxybiphenyl (compID# c0372)
 - o cis-2-Hydroxypenta-2,4-dienoate (compID# c0102)
 - o 2,3-Dihydroxybiphenyl (compID# c0373)
 - o 2-Hydroxy-6-oxo-6-phenylhexa-2,4-dienoate (compID# c0374)
- Enzymes in the Biphenyl Pathway
 - o EC 1.3.1.56 2,3-dihydro-2,3-dihydroxybiphenyl dehydrogenase (enzymeID# e0134)
 - o EC 1.13.11.39 2,3-dihydroxybiphenyl 1,2-dioxygenase (enzymeID# e0127)
 - o EC 1.14.99 biphenyl dioxygenase (enzymeID# e0089)
 - o EC 3.7.1.8 2-hydroxy-6-oxo-6-phenylhexa-2,4-dienoate hydrolase (enzymeID# e0033)

[Biphenyl Pathway Map] [BBD Main Menu]

Page Author(s): David Linden and Zhifu Sun
April 21, 1998 2:51:46 PM CDT BBDMaster@email.labmed.umn.edu

This is the UM-BBD compounds and reactions page for the Biphenyl Pathway. It was generated on December 6, 2000 3:19:53 PM CST.

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Wagner, Heindel, and Noyes, Inc.

Consulting Geologists

434 Shelburne Rd. Burlington, VT 05401

PAGE	_ OF _	Page No.
PROJECT:		

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000090-	43-7	Ί	_	2-PHENYLPHENOL	•	CAS *: 000067-	-66-3	7	CHLOROFORM			
C ₁₂	H ₁₀ C			. OH	.	Mol Weight	HCI ₃		· · · · · · · · · · · · · · · · · · ·	Çı		
170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170.21 170			ick			MP (deg C): +63.52 BP (deg C):		F2 100	÷ 5.:	CI CI		
BP pressure (mm Hg):		_	Data			8P pressure (mm Hg):	:		1 0200			
Property/ Value	Units	Terro	Type	Reference		Property/ Value	Units	Tems	Cara Type	Reference		
ws 7.00E+002	mgl	25	EXP	IARC (1983)	_ :	7.95E+003	mg1.	25	EXP	MACKAY,D ET AL. (1980)		
3.09			EXP	HANSCH,C ET AL. (1995)		1.97			EXP	HANSCH,C & LEO,AJ (1985)		
2.00E-003	пип На	25	EXT	KUNDEL,H ET AL (1975)		VP 1.97E+002	пит. Не	25	EXP	BOUBLIK,T ET AL. (1984)		
9,92	ρCı	25	EST	LIPNICK, RL ET AL. (1985)		ōc	рКа	Ì				
1.058-006	atm m3/mol	25	EST	VP/WSOL		HL 3.67E-003	a:m m3 mor	24	EXP	GOSSETT,JM (1987)		
сн 2.90≅-011	em3/mole sec	25	EST	MEYLAN,WM & HOWARD,PH (1993)	╝	04 1.03E-013	ст3 maid sec			ATKINSON,R (1989)		

AS #: 000108-93-0				CYCLOHEXANOL	<u>:</u>
Formula C ₆	H ₁₂ O			ОН	
Mor Wegot.	00.16				
MP (deg C): 23-25		75 .50	; 5 i		
32 (5eg Č):	51				
3.P pressure (mm Hg)					
Property/ Value	Units	Temp	Cara Type	Reference	
WS 3.75E+004	mat	25	EXP	RIDDICK, JA ET AL. (1986)	
1.23			EXP	HANSCH,C & LEO,AJ (1985)	:
8.00E-001	mm Hg	25	EXP	DAUBERT,TE & DANNER,RP (1989)	
oc	рКа				
2.81E-006	asm m3:mei	25	EST	VP/WSQL	
0H 1.74E-011	cm3/moid	25	EŞT	ATKINSON,R (1988)	

°°° 000065-	85-0	T		BENZOIC ACID			
C ₇	C ₇ H ₆ O ₂						
,							
122.12 M* dwg C; F* dwg Ck 122.4			2 €k				
→ → +			-				
249.2 Procure ,745 ng.				HO 0			
Property	Units	Terrs	Data Type	Releience			
3.40E+003	mgl	Ŧ "	i —	YALKOWSKY,SH & DANNENFELSER,RM (1992)			
1.87	<u> </u>		εхр	HANSCH,C & LEO,AJ (1985)			
1.20E-003	лл Нд	25	EXT	MALASPINA,L ET AL. (1973)			
4.21	экт	25	EXP	P SERJEANT, EP & DEMPSEY, B (1979)			
1.08E-007	am mlumos	25	EST	T MEYLAN, WM & HOWARD, PH (1991)			
1.15E-012	Sec Sec	25	EST	MEYLAN,WM & HOWARD,PH (1993)			

CAS #: 000108-	95-2		PHENOL					
Formula: C _d	H ₆ 0			ОН				
Moi Weight.	4.11			1				
MP (ceg C) 40.90		FP (SMG C)						
	1,839							
&P pressure (mm Hg)		_	Data					
Property/ Varue	Unns	Temp	Type	Reference				
ws 8.28E+004	mg/L	25	EXP	SOUTHWORTH, GR & KELLER, JL (1986)				
1.46			EXP	HANSCH,C & LEO,AJ (1985)				
3.50E-001	nan Hg	25	EXP	JONES,AH (1960)				
9.99	pKa	25	EXP	P SERJEANT, EP & DEMPSEY, B (1979)				
a.33E-007	alm m3 mat	25	EXP	GAFFNEY, JS ET AL. (1987)				
ਰਜ 2.63E-011	cm3 moic sec	25	EAD	ATKINSON,R (1989)				

CAS #: 000067-	64-1			ACETONE			
Formula: Cg	H ₆ O			0-1			
Mot Weight: 5	8.08			l ĭ			
MP (deg C): +94.7		يعن) وت	c.				
92 (04g C): 56.07							
8P pressure (mm Hg)							
Property/ Value	Units	Тетр	Data Type	Reference			
ws 1.00E+006	mal		EXP	RIODICK,JA ET AL. (1986)			
			EXP	HANSCH,C & LEO.AJ (1985)			
-0.24	l						
-0.24 VF 2.31E+002	rism hg	25	EXP	BOUBLIK,T ET AL. (1984)			
-0.24	mm hg pKa	25		BOUBLIK,T ET AL. (1984) SERJEANT,EP & DEMPSEY,B (1979)			
-0.24 VP 2.31E+002	 		EXP				



160 James Brown Orive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

MEMORANDUM

TO: Jon Wilkinson FR: Harry Locker DT: December 4, 2000

RE: Omya Draft

On November 27, 2000 Endyne received samples for chemical analysis from H&N for a project identified as OMYA. The samples were collected on 11/21/00 and consisted of varying concentrations of carbonate sludge in an aqueous matrix. The lab was requested to focus on samples SS#1 (166075), SS #2 (166076), SS-1 (core), 15-18 (166047), SS-1 (core) 0-4" (166051). Specifically, Endyne was tasked to analyze these samples for o-phenyl phenol (OPP) and potential breakdown metabolites including phenol, by EPA method 642.

Samples 166075, 166076, and 166047 each settled into two distinct fractions, carbonate sludge and a floating aqueous layer - when allowed to sit idle for several hours. The water phase of these samples was decanted and processed via EPA method 642. Sample 166051, however, did not separate into 2 phases when allowed to settle; and as such, was processed like a solid sample. This sample 166051 was determined to be 56% solid.

The following table summarizes the EPA 642 results.

Sample ID	Ref#	OPP	Phenol*
SS #1	166075	97.0 mg/L	6.36 mg/L
SS #2	166076	5.32 mg/L	0.6 mg/L
SS-1 (core) 15-18'	166047	168 mg/L	26.4 mg/L
SS-1 (core) 0-4"	166051**	42.3 mg/Kg	5.75 mg/Kg

^{*} Due to limitations inherent in HPLC/UV analyses, values reported for phenol are tentative. The early portion of the HPLC chromatogram is susceptible to potential interferences and as such, with ID by retention time only, we must note the detection as tentative.

^{**} Sample 166051 was extracted using an 18 hr soxlet procedure for this result.



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

The lab was also requested to evaluate the effect of KM_n04 on the contaminants present in the sludge material. Specifically, we were instructed to react sample 166051 with 0.5% by weight KM_n04 .

A 40 ml vial of the sample (166051) was split into 2 aliquots. One approximately 20g aliquot was extracted inside a soxlet sock with approximately 20 g of sodium sulfate. This mixture was extracted with MeCl₂ and the aid of ultrasonic disruption. The other approximate 20 g aliquot was reacted with 2.5 ml of a 5% solution of KM_n04 and allowed to react at room temperature for 90 min. This sample was then extracted as described above for the split sample. The results of this area as follows:

Sample ID	Ref#	OPP	Phenol*
SS-1 (core) 0-4"	166051	10.6 mg/kg	3.78 mg/kg
SS-1 (core) reacted with	166051 K	< 0.100	< 0.500
KM ₀ 04			

The extraction procedure used for these samples may not be as comprehensive as the 18hr soxlet procedure used on the original analysis. This most likely explains the reduced total concentration of OPP on the control sample relative to the previous result obtained using the 18 hr extraction.

Analysis of the extracts by GC/MS is ongoing in an effort to confirm tentative phenol detections.

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160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

To: Jeff, Jon, Dori and Meddie

Fr: H Locker Dt: 12/07/2000 Re: OMYA OPP

Previously, Endyne Inc. has reported tentative detection of Phenol in OMYA samples 166075, 166076, 166047, 166051 by EPA Method 642. Analysis of sample 166075 by SW 846 Method 8270C - where phenol is a target compound - did not detect the presence of phenol in this sample. Based on this finding, we must change our previous tentative identification to non detects for phenol for each of these samples analyzed by Method 642. This method, as we have it implemented, can not reliably be used for the determination of phenol.

As noted in the preliminary report, EPA Method 642 with UV detection has inherent limitations specifically with respect to non-target analytes. UV detection is a one dimensional technique with compound identification by retention time only. In addition to these instrumental limitations, the Method 642 extraction procedure may not be well suited for phenol analysis. Method 642 employs a liquid/liquid extraction of water samples at neutral pH. The reduced aromaticity of phenol with respect to OPP may result in a reduced extraction efficiency of phenol by Method 642. Additionally, the presence of other organics such as emulsifiers or floculants may further impair the extraction efficiency of phenol by this pH neutral extraction procedure.

Therefore, to definitively evaluate these samples for the presence of phenol requires analysis by Method 8270C. This method employs a serial liquid / liquid extraction of the sample at pH <2 and pH > 12. This procedure would be expected to have an improved specificity for the extraction of Phenol. In fact, phenol is a target compound by method 8270C. Method 8270C utilizes GC/MS as the analytical technique. This analytical methodology provides a two dimensional characterization utilizing both retention time and mass spectral fingerprint. This procedure will also better address sample matrix effects because it utilizes an isotopically labeled d5 phenol analog as an extraction surrogate. This extraction surrogate will provide insight into the extraction efficiency of phenol for each individual sample.

We regret any inconvenience this retraction may cause; in our haste to deliver timely results, we were unable to identify all the complexities inherent in the analytical methodology and the unique sample matrix.

Currently the lab is evaluating the addition of OPP as a target compound to Method 8270C. Preliminary results show excellent correlation of results of OPP by Method 8270C with Method 642 results.

Much concern has been given to the potential breakdown pathway of OPP and potential intermediate metabolites. Obviously the plethora of potential bacteria each with its own



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

Currently the lab is evaluating the addition of OPP as a target compound to Method 8270C. Preliminary results show excellent correlation of results of OPP by Method 8270C with Method 642 results.

Much concern has been given to the potential breakdown pathway of OPP and potential intermediate metabolites. Obviously the plethora of potential bacteria each with its own specific enzymatic weaponry, combined with potential chemical and photochemical routes, could result in a myriad of potential pathways and intermediates. In the absence of a known and definitive pathway it would seem prudent to continue with an 8260 and 8270 testing regime. Together, these methods provide a comprehensive screen for potential metabolites. Although this may not be the most scientifically rigorous solution, it does provide a pragmatic approach to detecting the presence of a broad array of potential intermediates. The most likely intermediates would include oxygenated (carboxcylic acids, ethers, esters, ketones etc.) hydrocarbons of 12 carbons or less which are readily detected by these methods. Since each of these methods employs GC/MS detection, UIP characterization can provide an extremely comprehensive screen for potential metabolites.



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

Special Report

Client: Heindel & Noyes

Project: OMYA

Date Sampled: November 21, 2000

Order Id: 10366

Analysis Date: December 4, 2000

The laboratory was requested to evaluate the effect of KMnO₄ addition on the o-Phenyl phenol contaminate present in the sludge material of sample SS-1(core) 0-4"(reference #166051).

A 40 ml vial of the sample (166051) was split into 2 aliquots. One approximately 20g aliquot was extracted inside a soxlet thimble with approximately 20g of sodium sulfate. This mixture was extracted with MeCl₂ and the aid of ultrasonic disruption. The other approximate 20g aliquot was reacted with 2.5ml of 5% solution KMnO₄ and allowed to react at room temperature for 90 min. This sample was then extracted as described above for the split sample. The results of this analysis follows:

Sample ID	Ref#	OPP
SS-1 (core) 0-4"	L	10.6 mg/kg
SS-1 (core) reacted with KMnO4	166051K	< 0.100 mg/kg

Sincerely,

Harry B. Locker, Ph.D. Laboratory Director



LABORATORY REPORT

Laboratory Services

160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

Heindel & Noyes PO Box 64709,

Burlington, VT 05406-4709

Attn: J Noyes

PROJECT: OMYA ORDER ID: 10376

RECEIVE DATE: November 27, 2000

REPORT DATE: December 4, 2000

Enclosed please find the results of the analyses performed for the samples referenced on the attached chain of custody. Different groups of analyses may be reported under separate cover.

All samples were prepared and analyzed by requirements outlined in the referenced methods and within the specified holding times.

All instrumentation was calibrated with the appropriate frequency and verified by the requirements outlined in the referenced methods.

Blank contamination was not observed at levels affecting the analytical results.

Mun

Analytical method precision and accuracy was monitored by laboratory control standards which include matrix spike, duplicate and quality control analyses. These standards were determined to be within established laboratory method acceptance limits, unless otherwise noted.

Reviewed by,

Harry B. Locker, Ph.D. Laboratory Director

enclosures



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT

SW 8260

CLIENT: Heindel & Noyes

PROJECT: OMYA

SITE: SS #1

DATE RECEIVED: November 27, 2000 REPORT DATE: December 4, 2000 ANALYSIS DATE: November 27, 2000 ORDER ID: 10376

REFERENCE NUMBER: 166075

DATE SAMPLED: November 21, 2000

TIME SAMPLED: 11:30 AM

SAMPLER: DB/CA ANALYST: 725

	Result		Result
Parameter	ug/L	<u>Parameter</u>	<u>ug/L</u>
Benzene	< 1.0	1,1-Dichloropropene	< 1.0
Bromobenzene	< 1.0	cis-1,3-Dichloropropene	< 1.0
Bromochloromethane	< 2.0	trans-1,3-Dichloropropene	< 1.0
Bromodichloromethane	< 1.0	Ethylbenzene	< 1.0
Bromoform	< 1.0	Hexachlorobutadiene	< 5.0
Bromomethane	< 5.0	Isopropylbenzene	< 1.0
n-Butylbenzene	< 1.0	p-Isopropyltoluene	< 1.0
sec-Butylbenzene	< 1.0	Methylene Chloride	< 5.0
tert-Butylbenzene	< 1.0	МТВЕ	< 2.0
Carbon Tetrachloride	< 1.0	Naphthalene	< 5.0
	< 1.0	n-Propylbenzene	< 1.0
Chlorobenzene	< 5.0	Styrene	< 1.0
Chloroethane	< 1.0	1,1,1,2-Tetrachloroethane	< 2.0
Chloroform Chloromethane	< 10.0	1,1,2,2-Tetrachloroethane	< 2.0
2-Chlorotoluene	< 1.0	Tetrachloroethene	< 1.0
4-Chlorotoluene	< 1.0	Toluene	< 1.0
Dibromochloromethane	< 1.0	1,2,3-Trichlorobenzene	< 2.0
1,2-Dibromo-3-Chloropropane	< 2.0	1,2,4-Trichlorobenzene	< 2.0
1,2-Dibromoethane	< 2.0	1,1,1-Trichloroethane	0.1>
Dibromomethane	< 2.0	1,1,2-Trichloroethane	< 1.0
1,2-Dichlorobenzene	< 1.0	Trichloroethene	< 1.0
	< 1.0	Trichlorofluoromethane	< 2.0
1,3-Dichlorobenzene	< 1.0	1,2,3-Trichloropropane	< 2.0
1,4-Dichlorobenzene Dichlorodifluoromethane	< 10.0	1,2,4-Trimethylbenzene	< 1.0
	< 1.0	1,3,5-Trimethylbenzene	< 1.0
1,1-Dichloroethane	< 1.0	Vinyl Chloride	< 2.0
1,2-Dichloroethane	< 1.0	Xylenes, Total	< 2.0
1,1-Dichloroethene	< 1.0	Surrogate 1	100.%
cis-1,2-Dichloroethene	< 1.0	Surrogate 2	100.%
trans-1,2-Dichloroethene	< 1.0	Surrogate 3	98.%
1,2-Dichloropropane	< 1.0	UIP's	1.
1,3-Dichloropropane	< 1.0	-	
2,2-Dichloropropanc	~ 1.0		



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT

CHARACTERIZATION OF UNIDENTIFIED PEAKS

CLIENT: Heindel & Noyes PROJECT NAME: OMYA

REPORT DATE: November 28, 2000 DATE RECEIVED: November 27, 2000

ORDER ID: 10376 SAMPLER: DB/CA

Unidentified peak characterization is achieved by direct comparison of sample and library spectral data:

REFERENCE NUMBER: 166,075

SITE: SS #1

NUMBER UIPs: 1

CHARACTERIZATION: Unidentified peak is Acetone at approximately 950 ug/L.

— LIND I IND, IND.

160 James Brown Drive
Williston, Vermont 05495

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5	Nitrate N		10	Alkalinity		15	Conduc		20	8010/8020	25	8270 B/N or Acid	30		
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LABORATORY REPORT

Laboratory Services

160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

Heindel & Noyes

PO Box 64709,

Burlington, VT 05406-4709

Attn: J Noyes

PROJECT: OMYA ORDER ID: 10366

RECEIVE DATE: November 27, 2000

REPORT DATE: December 4, 2000

Enclosed please find the results of the analyses performed for the samples referenced on the attached chain of custody. Different groups of analyses may be reported under separate cover.

All samples were prepared and analyzed by requirements outlined in the referenced methods and within the specified holding times.

All instrumentation was calibrated with the appropriate frequency and verified by the requirements outlined in the referenced methods.

Blank contamination was not observed at levels affecting the analytical results.

Mun

Analytical method precision and accuracy was monitored by laboratory control standards which include matrix spike, duplicate and quality control analyses. These standards were determined to be within established laboratory method acceptance limits, unless otherwise noted.

Reviewed by,

Harry B. Locker, Ph.D. Laboratory Director

enclosures



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT

SW 8260

CLIENT: Heindel & Noyes

PROJECT: OMYA SITE: SS-1(core) 15-18'

DATE RECEIVED: November 27, 2000 REPORT DATE: December 4, 2000 ANALYSIS DATE: November 27, 2000 ORDER ID: 10366

REFERENCE NUMBER: 166047

DATE SAMPLED: November 21, 2000

TIME SAMPLED: NI SAMPLER: DB/CA ANALYST: 725

	Result		Result
Parameter	ug/L	<u>Parameter</u>	ug/L
Benzene	< 1.0	1,1-Dichloropropene	< 1.0
Bromobenzene	< 1.0	cis-1,3-Dichloropropene	< 1.0
Bromochloromethane	< 2.0	trans-1,3-Dichloropropene	< 1.0
Bromodichloromethane	< 1.0	Ethylbenzene	< 1.0
Bromoform	< 1.0	Hexachlorobutadiene	< 5.0
Bromomethane	< 5.0	Isopropylbenzene	< 1.0
n-Butylbenzene	< 1.0	p-Isopropyltoluene	< 1.0
sec-Butylbenzene	< 1.0	Methylene Chloride	< 5.0
tert-Butylbenzene	. < 1.0	MTBE	< 2.0
Carbon Tetrachloride	< 1.0	Naphthalene	< 5.0
Chlorobenzene	< 1.0	n-Propylbenzene	< 1.0
Chloroethane	< 5.0	Styrene	< 1.0
Chloroform	< 1.0	1,1,1,2-Tetrachloroethane	< 2.0
Chloromethane	< 10.0	1,1,2,2-Tetrachloroethane	< 2.0
2-Chlorotoluene	< 1.0	Tetrachloroethene	< 1.0
4-Chlorotoluene	< 1.0	Toluene	2.3
Dibromochloromethane	< 1.0	1,2,3-Trichlorobenzene	< 2.0
1,2-Dibromo-3-Chloropropane	< 2.0	1,2,4-Trichlorobenzene	< 2.0
1,2-Dibromoethane	< 2.0	1,1,1-Trichloroethane	< 1.0
Dibromomethane	< 2.0	1,1,2-Trichloroethane	< 1.0
1,2-Dichlorobenzene	< 1.0	Trichloroethene	< 1.0
1,3-Dichlorobenzene	< 1.0	Trichlorofluoromethane	< 2.0
1,4-Dichlorobenzene	< 1.0	1,2,3-Trichloropropane	< 2.0
Dichlorodifluoromethane	< 10.0	1,2,4-Trimethylbenzene	< 1.0
1.1-Dichloroethane	< 1.0	1,3,5-Trimethylbenzene	< 1.0
1,2-Dichloroethane	< 1.0	Vinyl Chloride	< 2.0
1,1-Dichloroethene	< 1.0	Xylenes, Total	< 2.0
	< 1.0	Surrogate 1	96.%
cis-1,2-Dichloroethene	< 1.0	Surrogate 2	102.%
trans-1,2-Dichloroethene	< 1.0	Surrogate 3	100.%
1,2-Dichloropropane	< 1.0	UIP's	1.
1,3-Dichloropropane	< 1.0	-	
2,2-Dichloropropane	~ 1.0		



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT

SW 8260

CLIENT: Heindel & Noyes

PROJECT: OMYA SITE: SS-1(core) 0-4"

DATE RECEIVED: November 27, 2000 REPORT DATE: December 4, 2000 ANALYSIS DATE: November 27, 2000 ORDER ID: 10366

REFERENCE NUMBER: 166051 DATE SAMPLED: November 21, 2000

TIME SAMPLED: NI

SAMPLER: DB/CA ANALYST: 725

	Result		Result
<u>Parameter</u>	ug/kg, as received	Parameter Parameter	ug/kg, as received
Benzene	< 5.0	1,1-Dichloropropene	< 5.0
Bromobenzene	< 5.0	cis-1,3-Dichloropropene	< 5.0
Bromochloromethane	< 10.0	trans-1,3-Dichloropropene	< 5.0
Bromodichloromethane	< 5.0	Ethylbenzene	< 5.0
Bromoform	< 5.0	Hexachlorobutadiene	< 25.0
Bromomethane	< 25.0	Isopropylbenzene	< 5.0
n-Butylbenzene	< 5.0	p-Isopropyltoluene	< 5.0
sec-Butylbenzene	< 5.0	Methylene Chloride	< 25.0
tert-Butylbenzene	< 5.0	MTBE	< 10.0
Carbon Tetrachloride	< 5.0	Naphthalene	< 25.0
Chlorobenzene	< 5.0	n-Propylbenzene	< 5.0
Chloroethane	< 25.0	Styrene	< 5.0
Chloroform	< 5.0	1,1,1,2-Tetrachloroethane	< 10.0
Chloromethane	< 50.0	1,1,2,2-Tetrachloroethane	< 10.0
4-Chlorotoluene	< 5.0	Tetrachloroethene	< 5.0
2-Chlorotoluene	< 5.0	Toluene	< 5.0
Dibromochloromethane	< 5.0	1,2,3-Trichlorobenzene	< 10.0
1,2-Dibromo-3-Chloropropane	< 10.0	1,2,4-Trichlorobenzene	< 10.0
1,2-Dibromoethane	< 10.0	1,1,1-Trichloroethane	< 5.0
Dibromomethane	< 10.0	1,1,2-Trichloroethane	< 5.0
1,2-Dichlorobenzene	< 5.0	Trichloroethene	< 5.0
1,3-Dichlorobenzene	< 5.0	Trichlorofluoromethane	< 10.0
1.4-Dichlorobenzene	< 5.0	1,2,3-Trichloropropane	< 10.0
Dichlorodifluoromethane	< 50.0	1,2,4-Trimethylbenzene	< 5.0
1,1-Dichloroethane	< 5.0	1,3,5-Trimethylbenzene	< 5.0
1.2-Dichloroethane	< 5.0	Vinyl Chloride	< 10.0
1,1-Dichloroethene	< 5.0	Xylenes, Total	< 10.0
cis-1.2-Dichloroethene	< 5.0	Surrogate 1	106.%
trans-1,2-Dichloroethene	< 5.0	Surrogate 2	104.%
1,2-Dichloropropane	< 5.0	Surrogate 3	100.%
1,3-Dichloropropane	< 5.0	UIP's	t.
2,2-Dichloropropane	< 5.0	Percent Solids	NA





160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT

CHARACTERIZATION OF UNIDENTIFIED PEAKS

CLIENT: Heindel & Noyes PROJECT NAME: OMYA

REPORT DATE: November 28, 2000 DATE RECEIVED: November 27, 2000

ORDER ID: 10366 SAMPLER: DB/CA

Unidentified peak characterization is achieved by direct comparison of sample and library spectral data:

REFERENCE NUMBER: 166,047

SITE: SS - 1 (core) 15-18'

NUMBER UIPs: 1

CHARACTERIZATION: Unidentified peak is Acetone at approximately 550 ug/L.

REFERENCE NUMBER: 166,051

SITE: SS - 1 (corc) 0-4"

NUMBER UIPs: 1

CHARACTERIZATION: Unidentified peak is Acetone at approximately 470 ug/kg.

Date pages 7	From (119.	Co,	Phone v	Fax 4	
7671					
Post-it* Fax Note	7000 m	Co/Dept.	Phone #	Fax #	

Billing Address: Reporting Address: HIN Project Name: OMYA HIN Sampler Name: A+V Company: Endyne Order ID: Contact Name/Phone #: J. Noyes -Į Phone #: 10300 (Lab Use Only) 658 0820 Sample Containers 8 M P Analysis Sample Rush Date Time Field Results/Remarks Ref# Matrix Required Preservation Sample Identification No. Type/Size (Lab Use Only) 8760 + HPLC 2-8760=HCl 8760 + HPLC 3- HPLC = Nowle ZY 40ml VOA 6 15-181 H-70 (ore 1-8760 HCC 2-Hplc=Noc 3 6-9" 166047 SS-3 lone - 14(し vl 166049 SS-4 (come - None 75-281 1-401 3 ન 166050 SS-5 (cone 12-15 - Wove_ 8260 +HPLK+ % Solids None 0-Y" Slugge (0605/155-1 (Cone) 0-Y" (one 0-44 160033 SS-4 (ore 41 B-4" 6602571 SS-5 (ore Date/Time Received by: Received by: Date/Time 1/12760 1/50 Requested Analyses No X New York State Project: Yes 8270 PAH 1664 TPH/FOG 26 Sulfate **Total Solids** 16 TKN 11 pН 8015 GRO 27 PP13 Metals 22 17 Coliform (Specify) 12 TSS Total P Chloride RCRA8 Metals 18 23 8015 DRO COD 13 TDS Total Diss. P Ammonia N 29 24 8260/8260B 19 8021B Turbidity COS Nitrite N 30 25 8270 B/N or Acid 15 | Conductivity 20 8010/8020 10 Alkalinity Nitrate N 31 Metals (As Is, Total, Diss.) Ag, Al, As, B, Ba, Be, Ca, Cd, Co, Cr, Cu, Fe, Hg, K, Mg, Mn, Mo, Na, Ni, Pb, Sb, Se, Si, Sr, Ti, Tl, V, Zn TCLP (Specify: volatiles, semi-volatiles, metals, pesticides, herbicides) 33 34 Other



LABORATORY REPORT

Laboratory Services

160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

Heindel & Noyes

PO Box 64709,

Burlington, VT 05406-4709

Attn: Jeff Noyes

PROJECT: OMYA-Dogleg Spill

ORDER ID: 10502

RECEIVE DATE: December 5, 2000

REPORT DATE: December 6, 2000

Enclosed please find the results of the analyses performed for the samples referenced on the attached chain of custody. Different groups of analyses may be reported under separate cover.

All samples were prepared and analyzed by requirements outlined in the referenced methods and within the specified holding times.

All instrumentation was calibrated with the appropriate frequency and verified by the requirements outlined in the referenced methods.

Blank contamination was not observed at levels affecting the analytical results.

1/6//

Analytical method precision and accuracy was monitored by laboratory control standards which include matrix spike, duplicate and quality control analyses. These standards were determined to be within established laboratory method acceptance limits, unless otherwise noted.

Reviewed by,

Harry B. Locker, Ph.D. Laboratory Director

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160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT

SW 8260

CLIENT: Heindel & Noyes

PROJECT: OMYA-Dogleg Spill

SITE: Well 96-1

DATE RECEIVED: December 5, 2000 REPORT DATE: December 6, 2000

ANALYSIS DATE: December 5, 2000

ORDER ID: 10502

REFERENCE NUMBER: 166509 DATE SAMPLED: December 4, 2000

TIME SAMPLED: 12:45 PM

SAMPLER: AP ANALYST: 725

	Result		Result
Parameter	ug/L	<u>Parameter</u>	ng/L
Benzene	< 1.0	1,1-Dichloropropene	< 1.0
Bromohenzene	< 1.0	cis-1,3-Dichloropropene	< 1.0
Bromochloromethane	< 2.0	trans-1,3-Dichloropropene	< 1.0
Bromodichloromethane	< 1.0	Ethylbenzene	< 1.0
Bromoform	< 1.0	Hexachlorobutadiene	< 5.0
Bromomethane	< 5.0	Isopropylbenzene	< 1.0
n-Butylbenzene	< 1.0	p-Isopropyltoluene	< 1.0
sec-Butylbenzene	< 1.0	Methylene Chloride	< 5.0
tert-Butylbenzene	< 1.0	MTBE	< 2.0
Carbon Tetrachloride	< 1.0	Naphthalene	< 5.0
Chlorobenzene	< 1.0	n-Propylbenzene	< 1.0
Chloroethane	< 5.0	Styrene	< 1.0
Chloroform	< 1.0	1,1,1,2-Tetrachloroethane	< 2.0
Chloromethane	< 10.0	1,1,2,2-Tetrachloroethane	< 2.0
2-Chlorotoluene	< 1.0	Tetrachloroethene	< 1.0
4-Chlorotoluene	< 1.0	Toluene	< 1.0
Dibromochloromethane	< 1.0	1,2,3-Trichlorobenzene	< 2.0
1,2-Dibromo-3-Chloropropane	< 2.0	1,2,4-Trichlorobenzene	,< 2.0
1,2-Dibromoethane	< 2.0	1,1,1-Trichloroethane	< 1.0
Dibromomethane	< 2.0	1,1,2-Trichloroethane	< l.0
1,2-Dichlorobenzene	< 1.0	Trichloroethene	< 1.0
1,3-Dichlorobenzene	< 1.0	Trichlorofluoromethane	< 2.0
1,4-Dichlorobenzene	< 1.0	1,2,3-Trichloropropane	< 2.0
Dichlorodifluoromethane	< 10.0	1,2,4-Trimethylbenzene	< 1.0
1,1-Dichloroethane	< 1.0	1,3,5-Trimethylbenzene	< 1.0
1.2-Dichloroethane	< 1.0	Vinyl Chioride	< 2.0
1,1-Dichloroethene	< 1.0	Xylenes, Total	< 2.0
cis-1,2-Dichloroethene	< 1.0	Surrogate 1	100.%
trans-1,2-Dichloroethene	< 1.0	Surrogate 2	97.%
1,2-Dichloropropane	< 1.0	Surrogate 3	91.%
1,3-Dichloropropane	< i.0	UIP's	0.
2,2-Dichloropropane	< 1.0		



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT SW 8260

CLIENT: Heindel & Noyes

PROJECT: OMYA-Dogleg Spill

SITE: Well #2

DATE RECEIVED: December 5, 2000 REPORT DATE: December 6, 2000 ANALYSIS DATE: December 5, 2000 ORDER ID: 10502

REFERENCE NUMBER: 166510 DATE SAMPLED: December 4, 2000

TIME SAMPLED: 3:45 PM

	Result		Result
Parameter	<u>ug/L</u>	<u>Parameter</u>	<u>ug/L</u>
Benzene	< 1.0	1,1-Dichloropropene	< 1.0
Bromobenzene	< 1.0	cis-1,3-Dichloropropene	< 1.0
Bromochloromethane	< 2.0	trans-1,3-Dichloropropene	< 1.0
Bromodichloromethane	< 1.0	Ethylbenzene	< 1.0
Bromoform	< 1.0	Hexachlorobutadiene	< 5.0
Bromomethane	< 5.0	Isopropylbenzene	< 1.0
n-Butylbenzene	< 1.0	p-Isopropyltoluene	< 1.0
sec-Butylbenzene	< 1.0	Methylene Chloride	< 5.0
tert-Butylbenzene	< 1.0	MTBE	< 2.0
Carbon Tetrachloride	< 1.0	Naphthalene	< 5.0
Chlorobenzene	< 1.0	n-Propylbenzene	< 1.0
Chloroethane	< 5.0	Styrene	< 1.0
Chloroform	< 1.0	1,1,1,2-Tetrachloroethane	< 2.0
Chloromethane	< 10.0	1,1,2,2-Tetrachloroethane	< 2.0
4-Chlorotoluene	< 1.0	Tetrachloroethene	< 1.0
2-Chlorotoluene	< 1.0	Toluene	< 1.0
Dibromochloromethane	< 1.0	1,2,3-Trichlorobenzene	< 2.0
1,2-Dibromo-3-Chloropropane	< 2.0	1,2,4-Trichlorobenzene	< 2.0
1,2-Dibromoethane	< 2.0	1,1,1-Trichloroethane	< 1.0
Dibromomethane	< 2.0	1,1,2-Trichloroethane	< 1.0
1,2-Dichlorobenzene	< 1.0	Trichloroethene	< 1.0
1,3-Dichlorobenzene	< 1.0	Trichlorofluoromethane	< 2.0
1,4-Dichlorobenzene	< 1.0	1,2,3-Trichloropropane	< 2.0
Dichlorodifluoromethane	< 10.0	1,2,4-Trimethylbenzene	< 1.0
1,1-Dichloroethane	< 1.0	1,3,5-Trimethylbenzene	< 1.0
1,2-Dichloroethane	< 1.0	Vinyl Chloride	< 2.0
1,1-Dichloroethene	< 1.0	Xylenes, Total	< 2.0
cis-1,2-Dichloroethene	< 1.0	Surrogate I	101.%
trans-1,2-Dichloroethene	< 1.0	Surrogate 2	97.%
1,2-Dichloropropane	< 1.0	Surrogate 3	92.%
1,3-Dichloropropane	< 1.0	UIP's	0.
2,2-Dichloropropane	< 1.0		



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT SW 8260

CLIENT: Heindel & Noyes

PROJECT: OMYA-Dogleg Spill

SITE: Dolomite Quarry

DATE RECEIVED: December 5, 2000 REPORT DATE: December 6, 2000 ANALYSIS DATE: December 5, 2000 ORDER ID: 10502

REFERENCE NUMBER: 166511

DATE SAMPLED: December 4, 2000

TIME SAMPLED: 3:30 PM

	Result		Result
Parameter	ug/L	<u>Parameter</u>	ug/L
Benzene	< 1.0	1,1-Dichloropropene	< 1.0
Bromobenzene	< 1.0	cis-1,3-Dichloropropene	< 1.0
Bromochloromethane	< 2.0	trans-1,3-Dichloropropene	< 1.0
Bromodichloromethane	< 1.0	Ethylbenzene	< 1.0
Bremoform	< 1.0	Hexachlorobutadiene	< 5.0
Bromomethane	< 5.0	Isopropylbenzene	< 1.0
n-Butylbenzene	< 1.0	p-Isopropyltoluene	< 1.0
sec-Butylbenzene	< 1.0	Methylene Chloride	< 5.0
tert-Butylbenzene	< 1.0	MTBE	< 2.0
Carbon Tetrachloride	< 1.0	Naphthalene	< 5.0
Chlorobenzene	< 1.0	n-Propylbenzene	< 1.0
Chloroethane	< 5.0	Styrene	< 1.0
Chloroform	1.9	1,1,1,2-Tetrachloroethane	< 2.0
Chloromethane	< 10.0	1,1,2,2-Tetrachloroethane	< 2.0
4-Chlorotoluene	< 1.0	Tetrachloroethene	< 1.0
2-Chlorotoluene	< 1.0	Toluene	< 1.0
Dibromochloromethane	< 1.0	1,2,3-Trichlorobenzene	< 2.0
1,2-Dibromo-3-Chloropropane	< 2.0	1,2,4-Trichlorobenzene	< 2.0
1,2-Dibromoethane	< 2.0	1,1,1-Trichloroethane	< 1.0
Dibromomethane	< 2.0	1,1,2-Trichloroethane	< 1.0
1,2-Dichlorobenzene	< 1.0	Trichtoroethene	< 1.0
1,3-Dichlorobenzene	< 1.0	Trichlorofluoromethane	< 2.0
1,4-Dichlorobenzene	< 1.0	1,2,3-Trichloropropane	< 2.0
Dichlorodifluoromethane	< 10.0	1,2,4-Trimethylbenzene	< 1.0
1,1-Dichloroethane	< 1.0	1,3,5-Trimethylbenzene	< 1.0
1,2-Dichloroethane	< 10	Vinyl Chloride	< 2.0
1,1-Dichloroethene	< 1.0	Xylenes, Total	< 2.0
cis-1,2-Dichloroethene	< 1.0	Surrogate 1	103.%
trans-1,2-Dichloroethene	< 1.0	Surrogate 2	97.%
1,2-Dichloropropane	< 1.0	Surrogate 3	91.%
1,3-Dichloropropane	< 1.0	UIP's	2.
2,2-Dichloropropane	< 1.0		



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT

CHARACTERIZATION OF UNIDENTIFIED PEAKS

CLIENT: Heindel & Noyes PROJECT NAME: OMYA

REPORT DATE: December 6, 2000 DATE RECEIVED: December 5, 2000

ORDER ID: 10502 SAMPLER: AP

Unidentified peak characterization is achieved by direct comparison of sample and library spectral data:

REFERENCE NUMBER: 166,511

SITE: Dolomite Quarry NUMBER UIPs: 2

CHARACTERIZATION: Unidentified peaks are Acetone at 230 ug/L and 2-Butanone at 33 ug/L.

Toluene was observed at trace levels, below the quantitation limit (approximately 0.9 ug/L).

160 James Brown Drive Williston, Vermont 05495 (802) 879-4333

41139

Project in Endyne (Lab Us	Order ID:	_	3			ny: i	ress: NJ-N Ny None #: Sel		Nos	yes		Sam		bN eddia -C8	fe	- 1	
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11	nonia N	8	Total Diss. P		13	TDS		18	COL)		23	8015 DRO		28	RCRA8 Metals	
4 Nitr	te N	9	BOD		14	Turbidi	ty	19	8021	1B		24	8260/8260B		29		
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LABORATORY REPORT

Laboratory Services

160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

Heindel & Noyes

PO Box 64709,

Burlington, VT 05406-4709

Attn: J Noyes

PROJECT: OMYA ORDER ID: 10376

RECEIVE DATE: November 27, 2000

REPORT DATE: December 11, 2000

Enclosed please find the results of the analyses performed for the samples referenced on the attached chain of custody. Different groups of analyses may be reported under separate cover.

All samples were prepared and analyzed by requirements outlined in the referenced methods and within the specified holding times.

All instrumentation was calibrated with the appropriate frequency and verified by the requirements outlined in the referenced methods.

Blank contamination was not observed at levels affecting the analytical results.

1/1/1

Analytical method precision and accuracy was monitored by laboratory control standards which include matrix spike, duplicate and quality control analyses. These standards were determined to be within established laboratory method acceptance limits, unless otherwise noted.

Reviewed by,

Harry B. Locker, Ph.D. Laboratory Director

enclosures



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT SW 8270C

CLIENT: Heindel & Noyes

PROJECT: OMYA

SITE: SS #1

Page 2 of 2

DATE RECEIVED: November 27, 2000 REPORT DATE: December 11, 2000 ANALYSIS DATE: December 6, 2000 ORDER ID: 10376

REFERENCE NUMBER: 166075

DATE SAMPLED: November 21, 2000

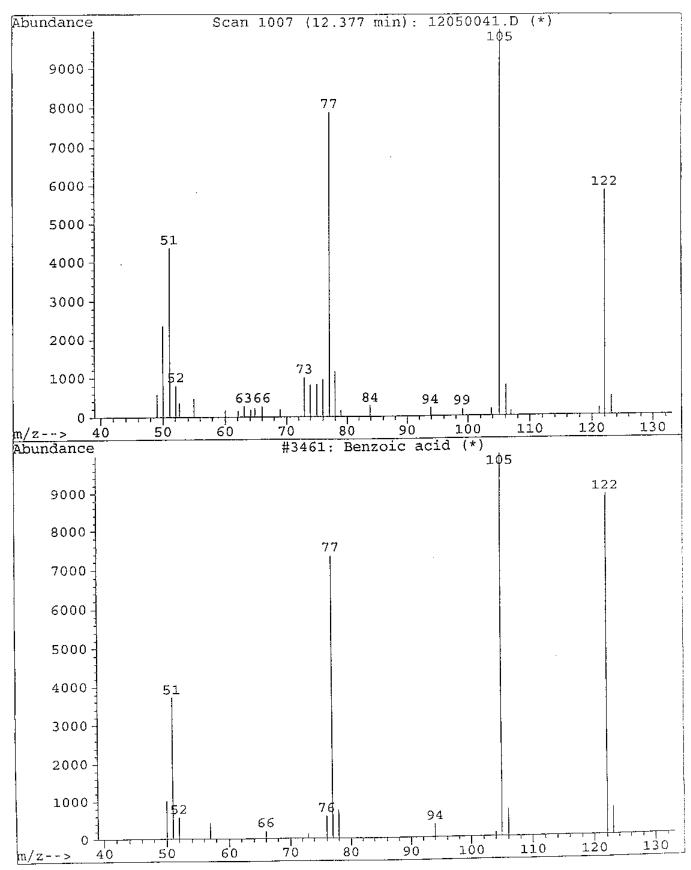
TIME SAMPLED: 11:30 AM

SAMPLER: DB/CA ANALYST: 917

	Result		Result
Parameter	ug/L	<u>Parameter</u>	<u>ug/L</u>
Acenaphthene	< 20.0	1-Methylnaphthalene	< 20.0
Acenaphthylene	< 20.0	2-Methylnaphthalene	< 20.0
Aniline	< 100.	Naphthalene	< 20.0
Anthracene	< 20.0	I-Naphthylamine	< 100.
Azobenzene	< 50.0	2-Naphthylamine	< 100.
Benzidine	< 100.	2-Nitroaniline	< 200.
Benzo(a)anthracene	< 20.0	3-Nitroaniline	< 200.
Benzo(b&k)fluoranthene	< 20.0	4-Nitroaniline	< 200.
Benzo(a)pyrene	< 20.0	Nitrobenzene	< 50.0
Benzo(g,h,i)perylene	< 20.0	N-Nitroso-di-n-butylamine	< 50.0
Bis(2-chloroethyl)ether	< 50.0	N-Nitrosodiphenylamine	< 50.0
Bis(2-chloroethoxy)methane	< 50.0	N-Nitrosodimethylamine	< 100.
Bis(2-ethylhexyl)phthalate	< 100.	N-Nitrosodi-n-propylamine	< 100.
Bis(2-chloroisopropyl)ether	< 100.	N-Nitrosopiperidine	< 100.
4-Bromophenyl phenyl ether	< 20.0	Phenanthrene	< 20.0
Butyl benzyl phthalate	< 100.	Pyrene	< 20.0
Carbazole	< 100.	Pyridine	< 100.
4-Chloroaniline	< 50.0	1,2,4-Trichlorobenzene	< 20.0
1-Chloronaphthalene	< 20.0	Benzyl alcohol	< 100.
2-Chloronaphthalene	< 20.0	4-Chloro-3-methylphenol	< 100.
4-Chlorophenyl phenyl ether	< 20.0	2-Chlorophenol	< 50.0
Chrysene	< 20.0	2,4-Dichlorophenol	< 50.0
Dibenzofuran	< 20.0	2,6-Dichlorophenol	< 50.0
Dibenzo(a,h)anthracene	< 20.0	2,4-Dimethylphenol	< 50.0
Di-n-butylphthalate	< 100.	4,6-Dinitro-2-methylphenol	< 500.
1,2-Dichlorobenzene	< 20.0	2,4-Dinitrophenol	< 100.
1,3-Dichlorobenzene	< 20.0	2-Methylphenol (o-cresol)	< 50.0
1,4-Dichlorobenzene	< 20.0	3&4-Methylphenol (m&p-cresol)	< 50.0
3,3'-Dichlorobenzidine	< 50.0	2-Nitrophenol	< 100.
Diethyl phthalate	< 100.	4-Nitrophenol	< 100.
Dimethyl phthalate	< 100.	Pentachlorophenol	< 150.
2,4-Dinitrotoluene	< 50.0	Phenol	< 50.0
2,6-Dinîtrotoluene	< 50.0	2,4,5-Trichlorophenol	< 100.
Di-π-octylphthalate	< 100.	2,4,6-Trichlorophenol	< 100.
Fluoranthene	< 20.0	Acid Surrogate 1	51.%
Fluorene	< 20.0	Acid Surrogate 2	62.%
Hexachlorobenzene	< 50.0	Acid Surrogate 3	95.%
Hexachlorobutadiene	< 50.0	Base/Neutral Surrogate 1	119.%
Hexachlorocyclopentadiene	< 200.	Base/Neutral Surrogate 2	127.%
Hexachloroethane	< 50.0	Base/Neutral Surrogate 3	149.%
Indeno(1,2,3-cd)pyrene	< 20.0	UIP's	4.
Isophorone	< 20.0		
- 4 44			

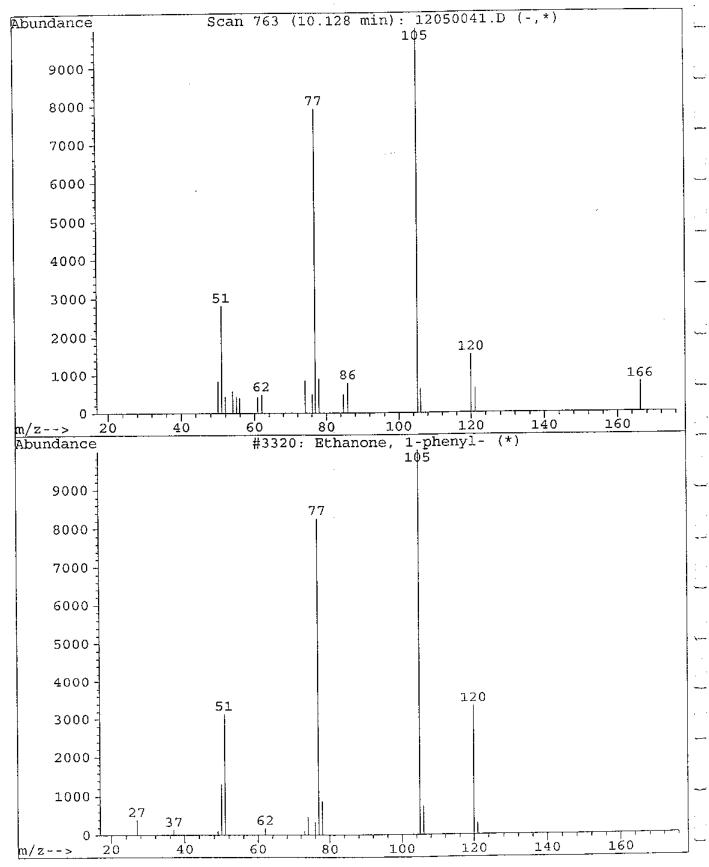
Quality : 90

ID : Benzoic acid



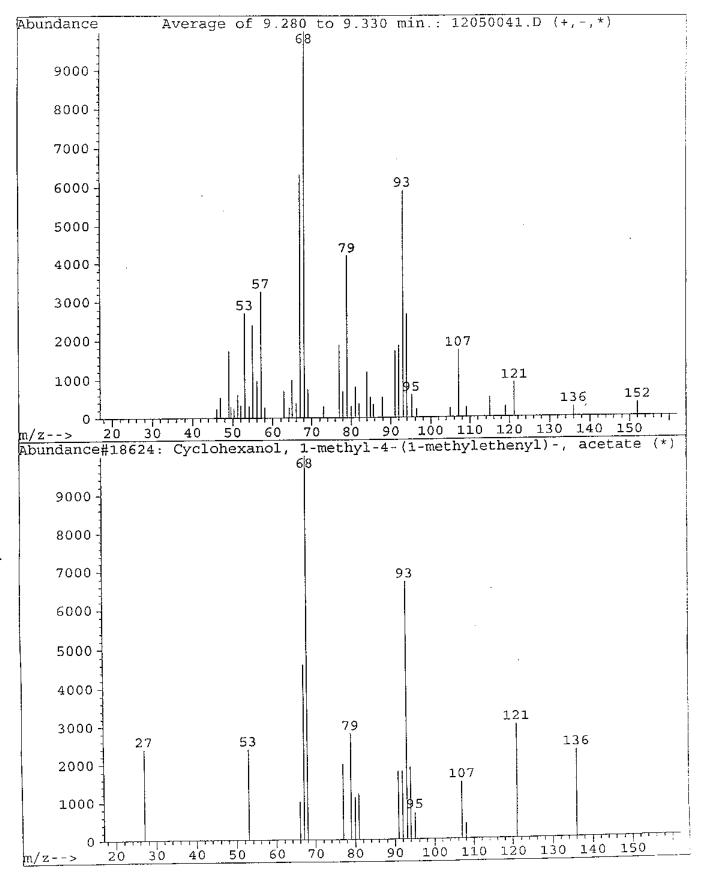
Quality : 45

ID : Ethanone, 1-phenyl-



Quality : 50

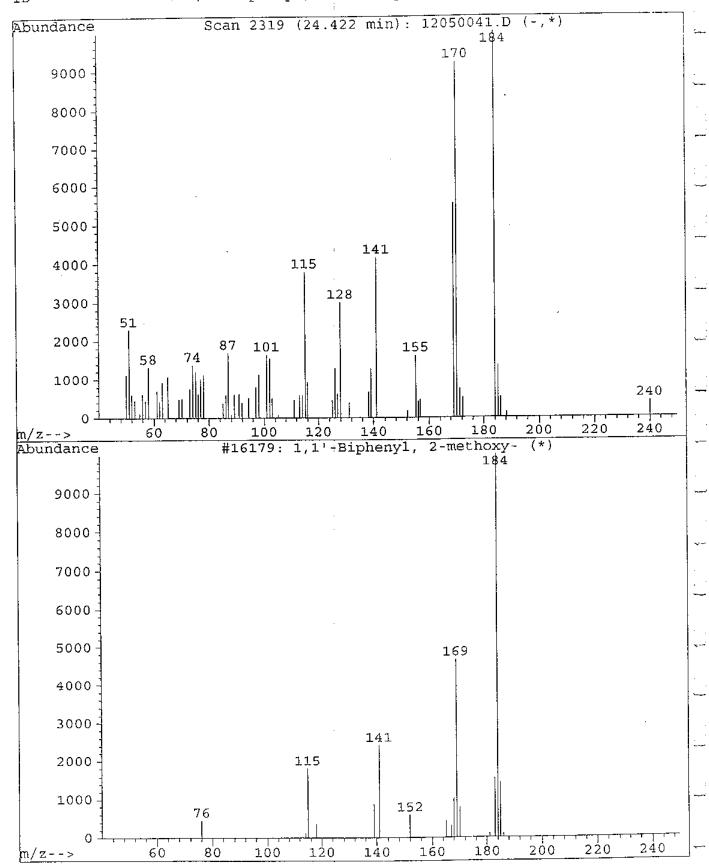
: Cyclohexanol, 1-methyl-4-(1-methylethenyl)-, acetate



Quality

: 47 ID

: 1,1'-Biphenyl, 2+methoxy-





160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT

CHARACTERIZATION OF UNIDENTIFIED PEAKS

CLIENT: Heindel & Noyes PROJECT NAME: OMYA

REPORT DATE: December 10, 2000 DATE RECEIVED: November 27, 2000

ORDER ID: 10376 SAMPLER: DB/CA

Unidentified peak characterization is achieved by direct comparison of sample and library spectral data:

REFERENCE NUMBER: 166,075

SITE: Dolomite Quarry NUMBER UIPs: 4

CHARACTERIZATION:

	Best Library match	Estimated Concentration ug/L
· 1	Benzoic Acid	15
2	1-Phenyl ethanone	5.
3	2-Methoxy-1,1-Biphenyl	2.
4	1-Methyl-4-(1-methylethenyl) cyclohexanol	2.

160 James Brown Drive

160 James Brown Drive Williston, Vermont 05495

Project Name: Reporting Address:					ess:	۶/	J		ng Address:	H;'N		
1	Endyne Order ID: (Lab Usc Only) Company: H Contact Name/Phone #						Sampler Name: DB 1 C 14 Phone #: 658 - 0820					
Ref# (Lab Use Only)	Sample Identifica	tion Matr	eix G	CO M P	Date/Time	No.	Type/Size	Field	Results/Remarks	Analysis Required	Sample Preservation	Rush 24h
166075	- SS #1	b	OV		11.30	2	JLamen Jone 1/00		47C opp 5260	Horg	none	297)
166076	, SS#2	Ha	o V		12:10	2 2	1602 Plas 1602 4006	Y.C.	(-PLC op,			X
						1	1602 Plas	به خــ		V	V	,
							As reguested by JPW			<u>ل</u>		
						<u> </u>	telan (1/27 3:00)			H13L		
Relinquished I	ry: L.S.L.	Date/Time /1/27/00	Received b	Y://	10	<u> </u>	Data/Time (1/27)	Receiv	ed by:		Date/Ti	me
	State Project: Yes N	0			Requested A	nalys	es					
I pH	6	TKN	11	Total S	olids	16	Sulfate	21	1664 TPH/FOG	26 27	8270 PAH PP13 Metals	
2 Chlorid	le 7	Total P		TSS		17	Coliform (Specify)	22	8015 GRO 8015 DRO	28	RCRA8 Metal	ls
3 Ammor		Total Diss. P	13	TDS Turbid		19	8021B	24	8260/8260B	29		L-747
4 Nitrite		BOD Alkalinity	15	Condu	<u> </u>	20	8010/8020	25	8270 B/N or Acid	30		
5 Nitrate	(As Is, Total, Diss.) Ag, Al	As. B. Ba. Be.	Ca, Cd.	Co, C	Cr, Cu, Fe, H			Ia, Ni,	Pb, Sb, Se, Si, S	Sr, Ti, T l, V	/,Zn	
	(Specify: volatiles, semi					33						
34 Other	(Opening, Fordingo, John		<u> </u>			•						



LABORATORY REPORT

Laboratory Services

160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

Heindel & Noyes PO Box 64709, Burlington, VT 05406-4709

Durington, VI 05400

Attn: Jeff Noyes

PROJECT: OMYA-Dogleg Spill

ORDER ID: 10502

RECEIVE DATE: December 5, 2000 REPORT DATE: December 7, 2000

Enclosed please find the results of the analyses performed for the samples referenced on the attached chain of custody. Different groups of analyses may be reported under separate cover.

All samples were prepared and analyzed by requirements outlined in the referenced methods and within the specified holding times.

All instrumentation was calibrated with the appropriate frequency and verified by the requirements outlined in the referenced methods.

Blank contamination was not observed at levels affecting the analytical results.

Analytical method precision and accuracy was monitored by laboratory control standards which include matrix spike, duplicate and quality control analyses. These standards were determined to be within established laboratory method acceptance limits, unless otherwise noted.

Asterisk in results column indicates calculated value based on single point calibration using external standard calibration.

NA in results column indicates data was not available, surrogate compounds not added to sample matrix.

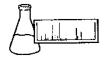
TBQ in results column indicates trace below quantitation limit.

1/1/

Reviewed by,

Harry B. Locker, Ph.D. Laboratory Director

Enclosures





160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT SW 8270C

CLIENT: Heindel & Noyes PROJECT: OMYA-Dogleg Spill

SITE: Well 96-1

DATE RECEIVED: December 5, 2000 REPORT DATE: December 7, 2000 ANALYSIS DATE: December 6, 2000 ORDER ID: 10502

REFERENCE NUMBER: 166509 DATE SAMPLED: December 4, 2000

TIME SAMPLED: 12:45 PM

	Result		Result
Parameter	ug/L	Parameter	ug/L
Acenaphthene	< 2.0	1-Methylnaphthalene	< 2.0
Acenaphthylene	< 2.0	2-Methylnaphthalene	< 2.0
Aniline	< 10.0	Naphthalene	< 2.0
Anthracene	< 2.0	1-Naphthylamine	< 10.0
Azobenzene	< 5.0	2-Naphthylamine	< 10.0
Benzidine	< 10.0	2-Nitroaniline	< 20.0
Benzo(a)anthracene	< 2.0	3-Nitroaniline	< 20.0
Benzo(b&k)fluoranthene	< 2.0	4-Nitroaniline	< 20.0
Велго(а)ругене	< 2.0	Nitrobenzene	< 5.0
Benzo(g,h,i)perylene	< 2.0	N-Nitroso-di-n-butylamine	< 5.0
Bis(2-chloroethyl)ether	< 5.0	N-Nitrosodiphenylamine	< 5.0
Bis(2-chloroethoxy)methane	< 5.0	N-Nitrosodimethylamine	< 10.0
Bis(2-ethylhexyl)phthalate	< 10.0	N-Nitrosodi-n-propylamine	< 10.0
Bis(2-chloroisopropyl)ether	< 10.0	N-Nitrosopiperidine	< 10.0
4-Bromophenyl phenyl ether	< 2.0	Phenanthrene	< 2.0
Butyl benzyl phthalate	< 10.0	Pyrene	< 2.0
Carbazole	< 10.0	Pyridine	< 10.0
4-Chloroaniline	< 5.0	1,2,4-Trichlorobenzene	< 2.0
1-Chloronaphthalene	< 2.0	Benzyl alcohol	< 10.0
2-Chloronaphthalene	< 2.0	4-Chloro-3-methylphenol	< 10.0
4-Chlorophenyl phenyl ether	< 2.0	2-Chlorophenol	< 5.0
Chrysene	< 2.0	2,4-Dichlorophenol	< 5.0
Dibenzofuran	< 2.0	2,6-Dichlorophenol	< 5.0
Dibenzo(a,h)anthracene	< 2.0	2,4-Dimethylphenol	< 5.0
Di-n-butylphthalate	< 10.0	4,6-Dinitro-2-methylphenol	< 50.0
1,2-Dichlorobenzene	< 2.0	2,4-Dinitrophenol	< 10.0
1,3-Dichlorobenzene	< 2.0	2-Methylphenol (o-cresol)	< 5.0
1,4-Dichlorobenzene	< 2,0	3&4-Methylphenol	< 5.0
3,3'-Dichlorobenzidine	< 5.0	2-Nitrophenol	< 10.0
Diethyl phthalate	< 10.0	4-Nitrophenol	< 10.0
Dimethyl phthalate	< 10.0	Pentachlorophenol	< 15.0
2,4-Dinitrotoluene	< 5.0	Phenol	< 5.0
2,6-Dinitrotoluene	< 5.0	2,4,5-Trichlorophenol	< 10.0
Di-n-octylphthalate	< 10.0	2,4,6-Trichlorophenol	< 10.0
Fluoranthene	< 2.0	Acid Surrogate !	NA
Fluorene	< 2.0	Acid Surrogate 2	NA
Hexachlorobenzene	< 5.0	Acid Surrogate 3	NA
Hexachlorobutadiene	< 5.0	Base/Neutral Surrogate 1	NA
Hexachlorocyclopentadiene	< 20.0	Base/Neutral Surrogate 2	NA
Hexachloroethane	< 5.0	Base/Neutral Surrogate 3	NA
Indeno(1,2,3-cd)pyrene	< 2.0	UIP's	0.
Isophorone	< 2.0	ortho-Phenyl phenol	< 5.0
Page 2 of 4			



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

Result

LABORATORY REPORT SW 8270C

CLIENT: Heindel & Noyes
PROJECT: OMYA-Dogleg Spill

SITE: Well #2

DATE RECEIVED: December 5, 2000 REPORT DATE: December 7, 2000 ANALYSIS DATE: December 6, 2000 ORDER ID: 10502

REFERENCE NUMBER: 166510 DATE SAMPLED: December 4, 2000

TIME SAMPLED: 3:45 PM

	Result		Kesutt
Danamatan	ug/L	Parameter	ug/L
Parameter	< 2.0	1-Methylnaphthalene	< 2.0
Acenaphthene	< 2.0	2-Methylnaphthalene	< 2.0
Acenaphthylene	< 10.0	Naphthalene	< 2.0
Aniline	< 2.0	I-Naphthylamine	< 10.0
Anthracene	< 5.0	2-Naphthylamine	< 10.0
Azobenzene	< 10.0	2-Nitroaniline	< 20.0
Benzidine	< 2.0	3-Nitroaniline	< 20.0
Benzo(a)anthracene	< 2.0	4-Nitroaniline	< 20.0
Benzo(b&k)fluoranthene	< 2.0	Nitrobenzene	< 5.0
Benzo(a)pyrene	< 2.0	N-Nitroso-di-n-butylamine	< 5.0
Benzo(g,h,i)perylene	< 5.0	N-Nitrosodiphenylamine	< 5.0
Bis(2-chloroethyl)ether	< 5.0 < 5.0	N-Nitrosodimethylamine	< 10.0
Bis(2-chloroethoxy)methane	< 10.0	N-Nitrosodi-n-propylamine	< 10.0
Bis(2-ethylhexyl)phthalate	< 10.0	N-Nitrosopiperidine	< 10.0
Bis(2-chloroisopropyl)ether	< 2.0	Phenanthrene	< 2.0
4-Bromophenyl phenyl ether	< 10.0	Pyrene	< 2.0
Butyl benzyl phthalate	< 10.0	Pyridine	< 10.0
Carbazole	< 5.0	1,2,4-Trichlorobenzene	< 2.0
4-Chloroaniline	< 2.0	Benzyl alcohol	< 10.0
1-Chloronaphthalene	< 2.0	4-Chloro-3-methylphenol	< 10.0
2-Chloronaphthalene	< 2.0	2-Chlorophenol	< 5.0
4-Chlorophenyl phenyl ether	< 2.0	2,4-Dichlorophenol	< 5.0
Chrysene	< 2.0	2.6-Dichlorophenol	< 5.0
Dibenzofuran	< 2.0 < 2.0	2,4-Dimethylphenol	< 5.0
Dibenzo(a,h)anthracene		4,6-Dinitro-2-methylphenol	< 50.0
Di-n-butylphthalate	< 10.0	2.4-Dinitrophenol	< 10.0
1,2-Dichlorobenzene	< 2.0	2-Methylphenol (o-cresol)	< 5.0
1,3-Dichlorobenzene	< 2.0	3&4-Methylphenol	< 5.0
1,4-Dichlorobenzene	< 2.0	2-Nitrophenol	< 10.0
3,3'-Dichtorobenzidine	< 5.0	4-Nitrophenol	< 10.0
Diethył phthalate	< 10.0	Pentachlorophenol	< 15.0
Dimethyl phthalate	< 10.0	Pheno!	< 5.0
2,4-Dinitrotoluene	< 5.0	2,4,5-Trichlorophenol	< 10.0
2,6-Dinitrotoluene	< 5.0	2,4,6-Trichlorophenol	< 10.0
Di-n-octylphthalate	< 10.0	Acid Surrogate 1	NA
Fluoranthene	< 2.0	Acid Surrogate 2	NA
Fluorene	< 2.0	Acid Surrogate 3	NA
Hexachlorobenzene	< 5.0	Base/Neutral Surrogate 1	NA
Hexachlorobutadiene	< 5.0	Base/Neutral Surrogate 2	NA
Hexachlorocyclopentadiene	< 20.0	Base/Neutral Surrogate 3	NA
Hexachloroethane	< 5.0		2.
Indeno(1,2,3-cd)pyrene	< 2.0	UIP's	< 5.0
Isophorone	< 2.0	ortho-Phenyl phenol	- 2.0
Page 3 of 4			
· • ·			



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT SW 8270C

CLIENT: Heindel & Noyes PROJECT: OMYA-Dogleg Spill SITE: Dolomite Quarry

Page 4 of 4

DATE RECEIVED: December 5, 2000 REPORT DATE: December 7, 2000 ANALYSIS DATE: December 6, 2000 ORDER ID: 10502

REFERENCE NUMBER: 166511 DATE SAMPLED: December 4, 2000

TIME SAMPLED: 3:30 PM

	Result		Result
Damen atom	ug/L	Parameter	ug/L
Parameter	< 2.0	1-Methylnaphthalene	< 2.0
Acenaphthene	< 2.0	2-Methylnaphthalene	< 2.0
Acenaphthylene	< 10.0	Naphthalene	< 2.0
Aniline	< 2.0	1-Naphthylamine	< 10.0
Anthracene	< 5.0	2-Naphthylamine	< 10.0
Azobenzene	< 10.0	2-Nitroaniline	< 20.0
Benzidine	< 2.0	3-Nitroaniline	< 20.0
Benzo(a)anthracene	< 2.0	4-Nitroaniline	< 20.0
Benzo(b&k)fluoranthene	< 2.0	Nitrobenzene	< 5.0
Benzo(a)pyrene	< 2.0	N-Nitroso-di-n-butylamine	< 5.0
Benzo(g,h,i)perylene	< 5.0	N-Nitrosodiphenylamine	< 5.0
Bis(2-chloroethyl)ether	< 5.0	N-Nitrosodimethylamine	< 10.0
Bis(2-chloroethoxy)methane	< 10.0	N-Nitrosodi-n-propylamine	< 10.0
Bis(2-ethylhexyl)phthalate	< 10.0	N-Nitrosopiperidine	< 10.0
Bis(2-chloroisopropyl)ether	< 2.0	Phenanthrene	< 2.0
4-Bromophenyl phenyl ether	< 10.0	Pyrene	< 2.0
Butyl benzyl phthalate	< 10.0	Pyridine	< 10.0
Carbazole	< 5.0	1.2.4-Trichlorobenzene	< 2.0
4-Chloroaniline	< 2.0	Benzyl alcohol	< 10.0
1-Chloronaphthalene	< 2.0 < 2.0	4-Chloro-3-methylphenol	< 10.0
2-Chloronaphthalene	< 2.0	2-Chlorophenol	< 5.0
4-Chlorophenyl phenyl ether	< 2.0	2,4-Dichlorophenol	< 5.0
Chrysene	< 2.0 < 2.0	2,6-Dichforophenol	< 5.0
Dibenzofuran	< 2.0 < 2.0	2,4-Dimethylphenol	< 5.0
Dibenzo(a,h)anthracene	< 10.0	4.6-Dinitro-2-methylphenol	< 50.0
Di-n-butylphthalate	< 10.0 < 2.0	2,4-Dinitrophenol	< 10.0
1,2-Dichlorobenzene	< 2.0	2-Methylphenol (o-cresol)	< 5.0
1,3-Dichlorobenzene	< 2.0	3&4-Methylphenol	< 5.0
1,4-Dichlorobenzene	< 5.0	2-Nitrophenol	< 10.0
3,3'-Dichlorobenzidine	< 10.0	4-Nitrophenol	< 10.0
Diethyl phthalate	< 10.0 < 10.0	Pentachlorophenol	< 15.0
Dimethyl phthalate	< 10.0 < 5.0	Pheno!	TBQ < 5.0
2,4-Dinitrotoluene	< 5.0 < 5.0	2,4,5-Trichlorophenol	< 10.0
2,6-Dinitrotoluene		2,4,6-Trichlorophenol	< 10.0
Di-n-octylphthalate	< 10.0	Acid Surrogate 1	NA.
Fluoranthene	< 2.0 < 2.0	Acid Surrogate 2	NA
Fluorene		Acid Surrogate 3	NA.
Hexachlorobenzene	< 5.0	Base/Neutral Surrogate 1	NA
Hexachlorobutadiene	< 5.0		NA
Hexachlorocyclopentadiene	< 20.0	Base/Neutral Surrogate 2	NA NA
Hexachloroethane	< 5.0	Base/Neutral Surrogate 3	7.
Indono(1,2,3-cd)pyrene	< 2.0	UIP's	7. 4,990.*
Isophorone	< 2.0	ortho-Phenyl phenol	4,770.



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT

CHARACTERIZATION OF UNIDENTIFIED PEAKS

CLIENT: Heindel & Noyes PROJECT NAME: OMYA

REPORT DATE: December 6, 2000 DATE RECEIVED: December 5, 2000

ORDER ID: 10502 SAMPLER: AP

Unidentified peak characterization is achieved by direct comparison of sample and library spectral data:

REFERENCE NUMBER: 166,510

SITE: Well #2 NUMBER UIPs: 2

CHARACTERIZATION:

	Best Library match	Estimated Concentration ug/L
· 1	Dimethyl-1,4-Dioxane	15.
2	2-Methyl-3-buten-2-ol	5.



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT

CHARACTERIZATION OF UNIDENTIFIED PEAKS

CLIENT: Heindel & Noyes PROJECT NAME: OMYA

REPORT DATE: December 6, 2000 DATE RECEIVED: December 5, 2000

ORDER ID: 10502 SAMPLER: AP

Unidentified peak characterization is achieved by direct comparison of sample and library spectral data:

REFERENCE NUMBER: 166,511

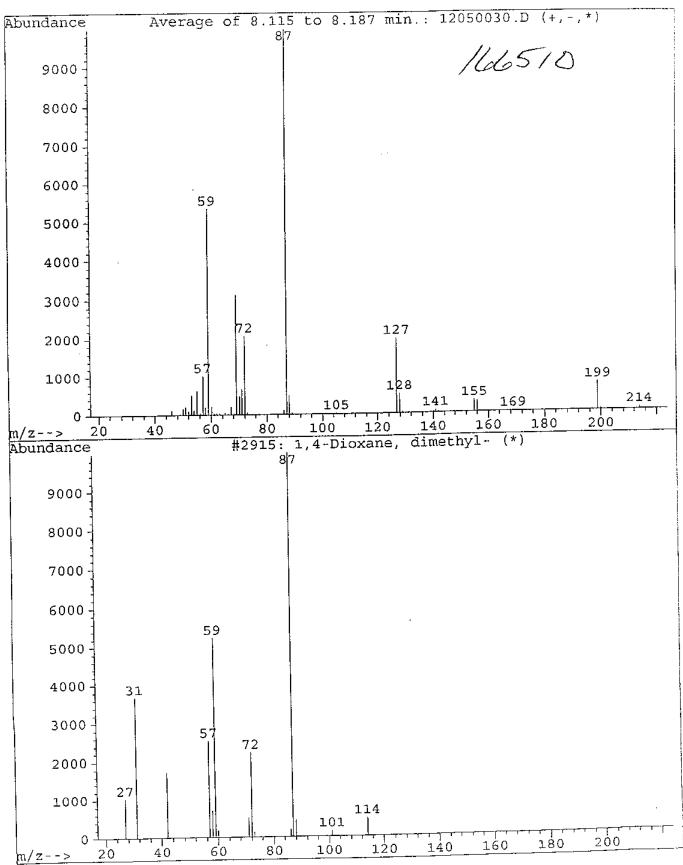
SITE: Dolomite Quarry NUMBER UIPs: 7

CHARACTERIZATION:

-	Best Library match	Estimated Concentration ug/L
- 1	Crotonic Acid	2.
2	2-Heptanone	2.
3	2-Piperidinone	3.
4	1-Phenyl ethanone	1.
5	2-Dibenzofuranol	2.
6	3-Amino Phenol	5.
7	UnKnown (no match)	10.
'		

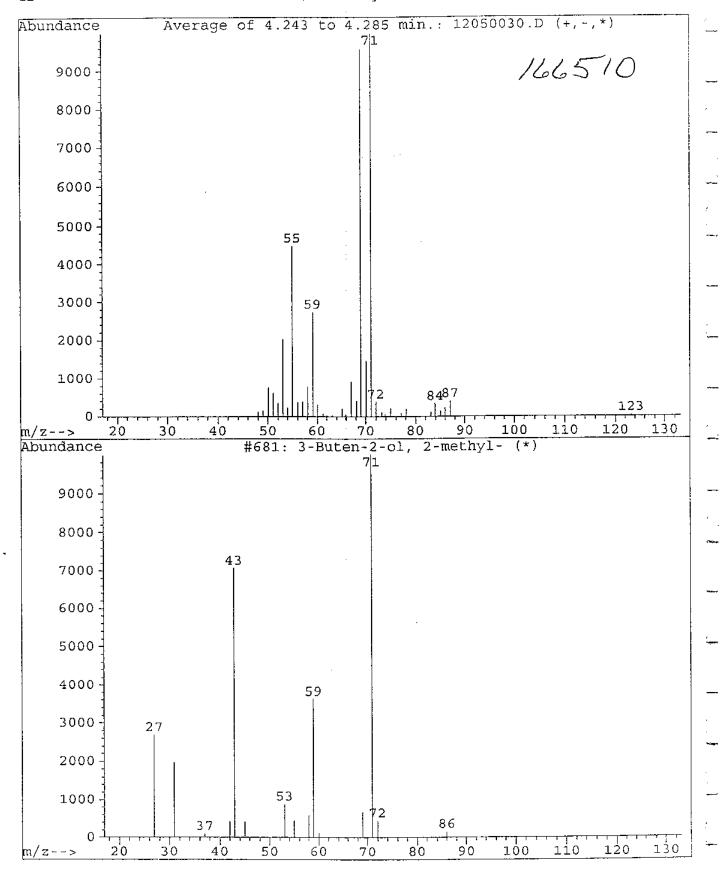
Quality : 59

ID : 1,4-Dioxane, dimethyl-

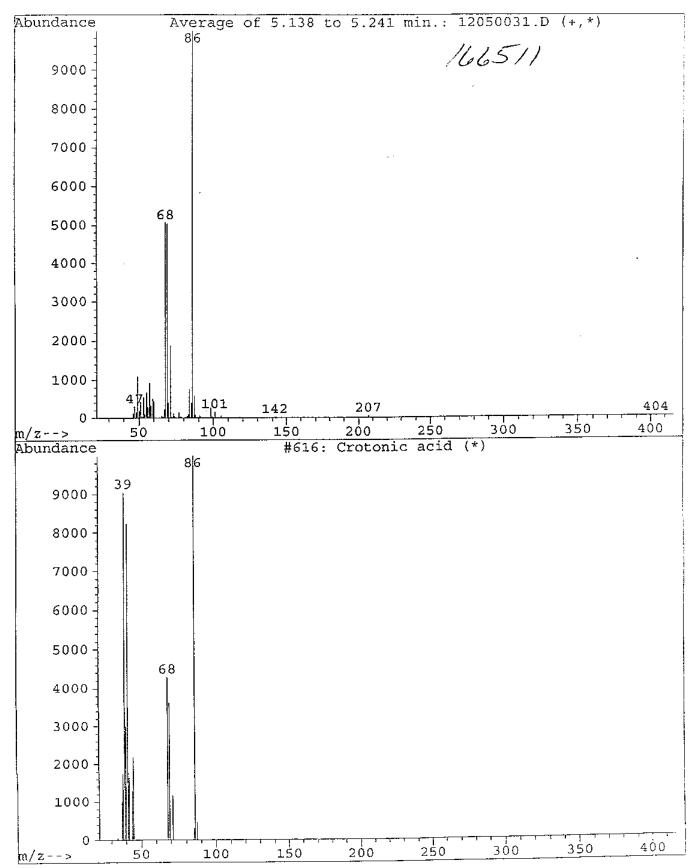


Quality ID

: 3-Buten-2-ol, 2-methyl-



ID : Crotonic acid

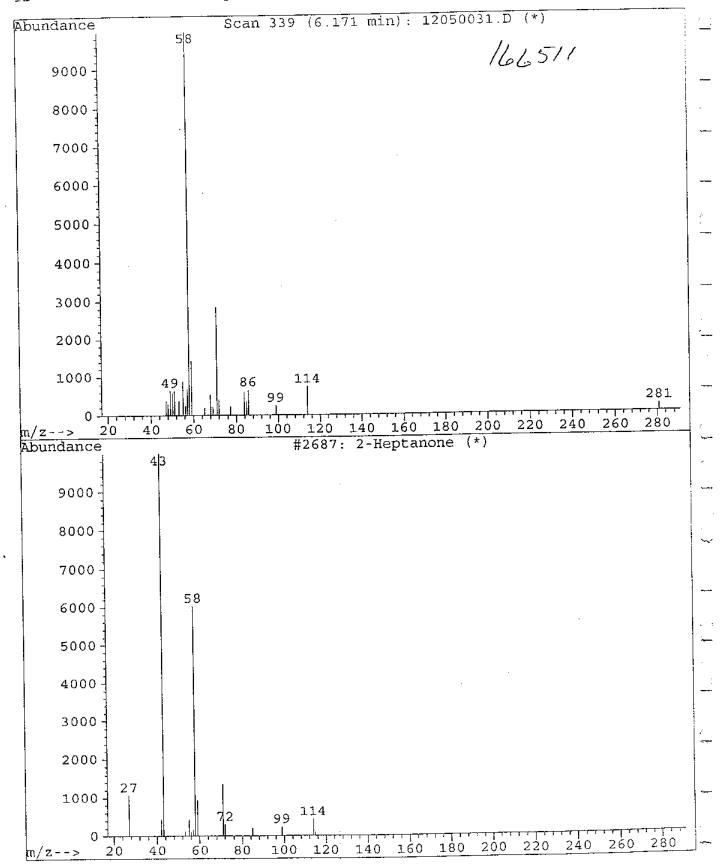


Quality

: 80

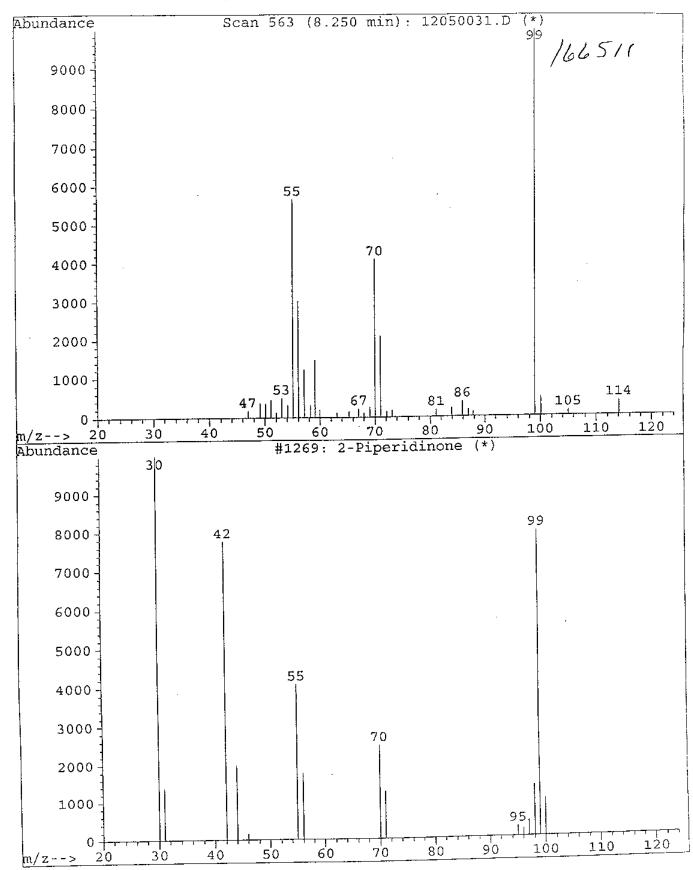
ĨD

: 2-Heptanone



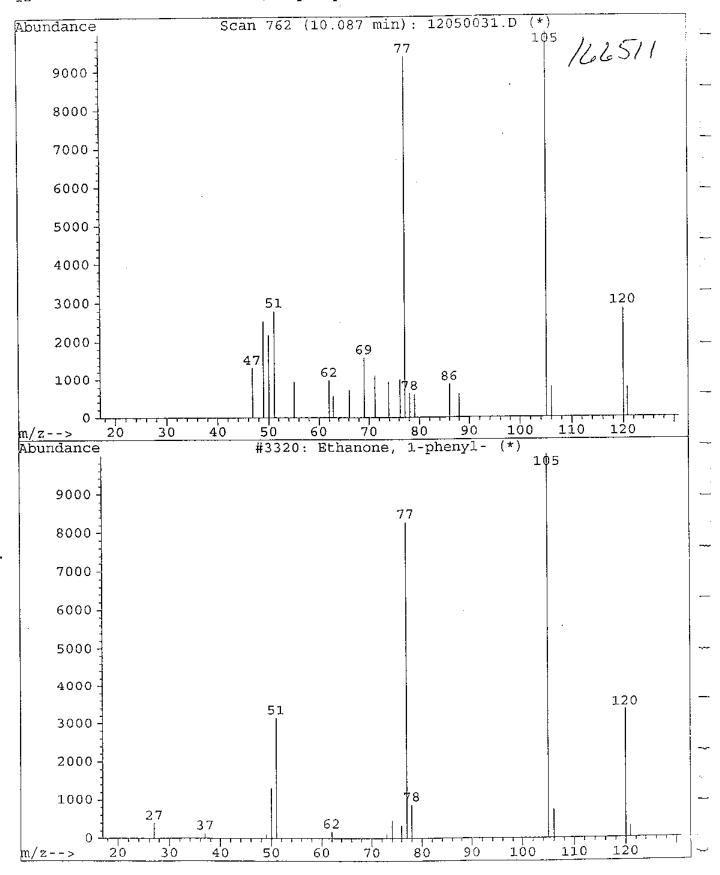
Quality : 64

ID : 2-Piperidinone



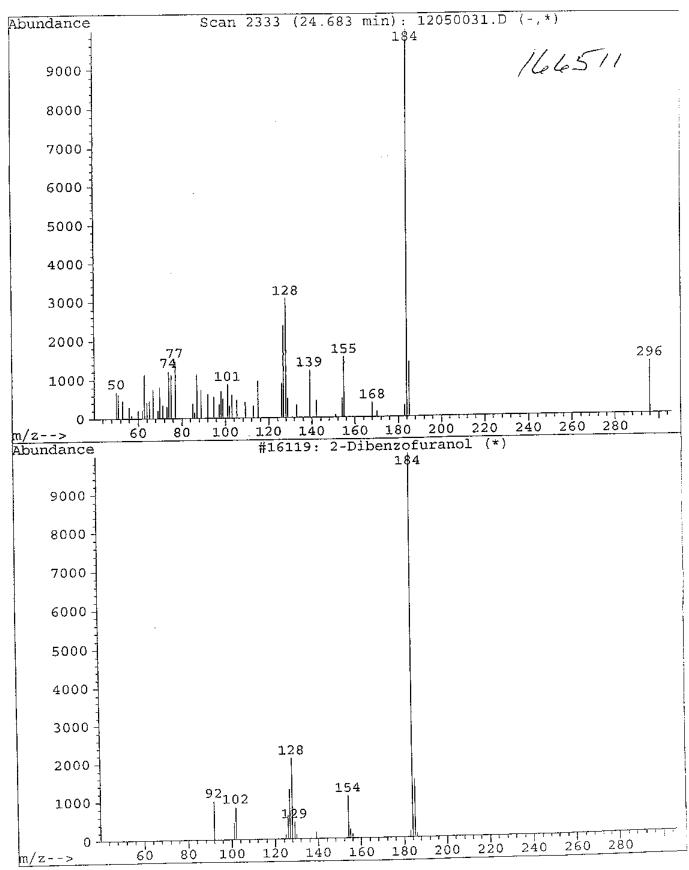
Quality ID

: Ethanone, 1-phenyl-



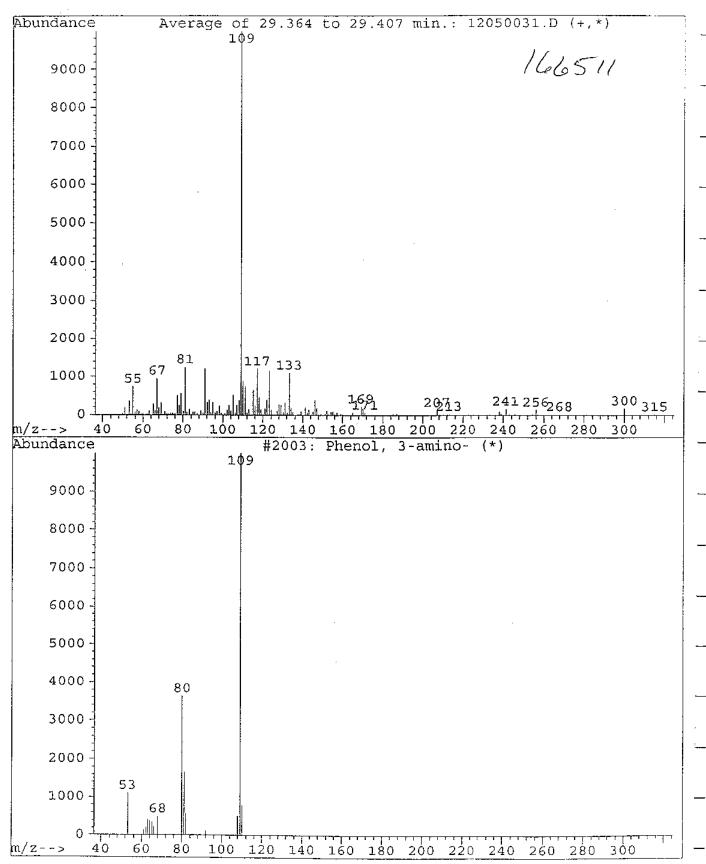
Quality : 53

ID : 2-Dibenzofuranol



Quality : 47

ID : Phenol, 3-amino-



: C:\HPCHEM\1\DATA\12050031.D File

Operator

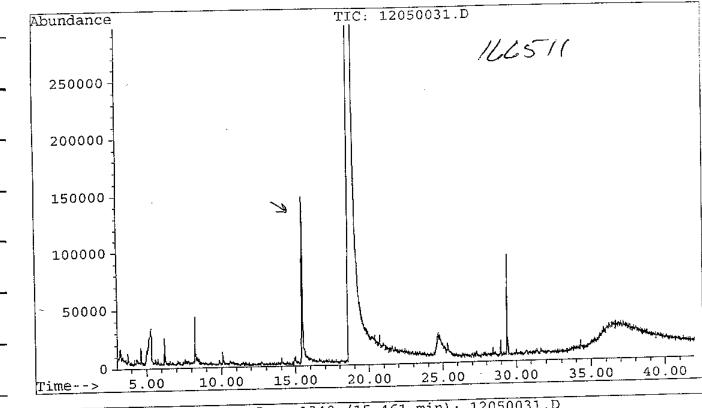
1:42 pm using AcqMethod AQ-82-2.M 6 Dec 100 Acquired

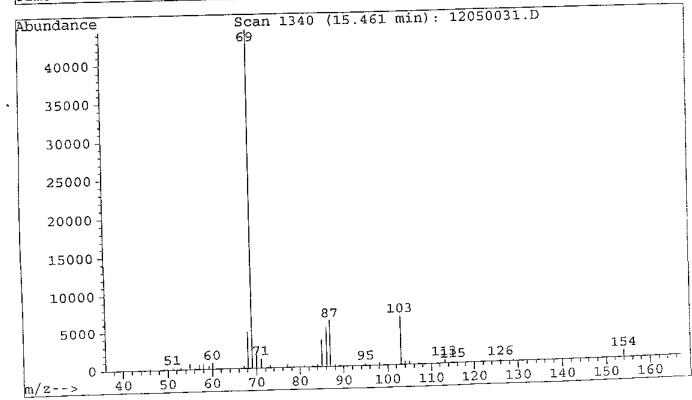
5971 - In Instrument :

Sample Name: 166511 925ul 910ml 12/5/00

Misc Info

Vial Number: 31





41139

160 James Brown Drive Williston, Vermont 05495 (802) 879-4333

	act Nom	2-AYMC	Dogl	eg Spil			ing Add	H)-1(Hba		#0017	<u> </u>
11 1	yne Ordo Use On	er ID:	502	23	7-I C	Company: HMW Contact Name/Phone #: Seff Noyes					- 1	Sampler Name: Amadda Ferry Phone #: 658-0820						
	ef# se Only)	Sample	Identific	ation	Matrix	C N M	C O M P	Date/Time	Sai		Containers Type/Size	Ę	ield Re	sults/Remarks		alysis quired	Sample Preservation	Rush
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1600	510	Well #	<u> </u>			14)	15:45	1					 		<u> </u>		
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1 1	рH		6	TKN		11	Total Sc	lids	16	Su	lifate	2	1 16	64 TPH/FOG		26	8270 PAH	
2	Chloride		7	Total P		12	TSS		17	Col	liform (Specify)	2	2 80	15 GRO		27	PP13 Metals	
3 Ammonia N 8 Total Diss. P 13 TDS 18 COD 23 8015 DRO								28	RCRA8 Metals									
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اللبارا																		
11																		
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	(White, Yellow, Pink Copy - Laboratory / Goldenrod Copy - Client)																	



LABORATORY REPORT

Laboratory Services

160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

Heindel & Noyes PO Box 64709,

Burlington, VT 05406-4709

Attn: Jeff Noyes

PROJECT: OMYA-Dogleg Spill

ORDER ID: 10502

RECEIVE DATE: December 5, 2000 REPORT DATE: December 7, 2000

Enclosed please find the results of the analyses performed for the samples referenced on the attached chain of custody. Different groups of analyses may be reported under separate cover.

All samples were prepared and analyzed by requirements outlined in the referenced methods and within the specified holding times.

All instrumentation was calibrated with the appropriate frequency and verified by the requirements outlined in the referenced methods.

Blank contamination was not observed at levels affecting the analytical results.

1/4/

Analytical method precision and accuracy was monitored by laboratory control standards which include matrix spike, duplicate and quality control analyses. These standards were determined to be within established laboratory method acceptance limits, unless otherwise noted.

Reviewed by,

Harry B. Locker, Ph.D. Laboratory Director

Enclosures

Page 1 of 2



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT

CLIENT: Heindel & Noyes
PROJECT: OMYA-Dogleg Spill
REPORT DATE: December 7, 2000

ORDER ID: 10502

DATE RECEIVED: December 5, 2000

SAMPLER: AP

Ref. Number: 166509	Site: Well 96-1		Date Sampled: Dec	ember 4, 2000	Time: 12:45 PM
Parameter	Result	<u>Unit</u>	Method	Analysis Dat	e <u>Analyst</u>
ortho-Phenyl phenol	< 0.001	mg/L	EPA 642	12/5/00	515
Ref. Number: 166510	Site: Well #2		Date Sampled: Dec	ember 4, 2000	Time: 3:45 PM
Parameter	Result	<u>Unit</u>	Method	Analysis Dat	e <u>Analyst</u>
ortho-Phenyl phenol	< 0.001	mg/L	EPA 642	12/5/00	515
Ref. Number: 166511	Site: Dolomite Quarry	:	Date Sampled: Dec	ember 4, 2000	Time: 3:30 PM
Parameter	Result	<u>Unit</u>	Method	Analysis Dat	<u>e Analyst</u>
ortho-Phenyl phenol	4.75	mg/L	EPA 642	12/6/00	515

= EIVD I IVE, IIVG. 160 James Brown Drive

160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 41139

(802) 879-4333		n.	namina Ada	least			Billi	ng Address:			أحنيها	
Project Name: OM VA - Docal	a spil		porting Add	iress: HJ-N)		2		HUN	-	±0017	
Endyne Order ID:	3		mpany:	HWW)			·	pler Name:	Ameddia	r te	ud	
(Lab Use Only) 10502	_	7-1 Co	ontact Name	HVW /Phone #: 5e	EE 1	Voyes	Phor	ie#: GE	<u>8-08</u>	3,2,c)	<u> </u>	
Ref# Sample Identifice	tion	Matrix	G C O M P	Date/Time	Samp	ole Containers Type/Size	Field l	Results/Remarks	An Re	alysis quired	Sample Preservation	Rush
(Lab Use Only)	·	110	B P	12:45	1eau	2 40ml 2 28 G				34	GO HOLAN	U
66509 Well 967	·· ····	1/20	101-	15:45	i	7.2.46						
000.510 0000		1),	+₩-	15:30						/		L
666511 Dolomite Qua	172	V	 M	1,5-50	- 1	- W						
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MAMALIN			///	tanul	/	11:50 An				<u></u>		
New York State Project/Yes			7-/	Requested	Analy	ses						, "
1 pH 6	TKN		11 Total	Solids	16	Sulfate	21	1664 TPH/FOG		26	8270 PAH	
2 Chloride 7	Total P		12 TSS		17	Coliform (Specify)	22	8015 GRO		27	PP13 Metals RCRA8 Metals	.16
3 Ammonia N 8	Total Diss. P		13 TDS		18	COD	23	8015 DRO		28	RCRA6 MEI	
4 Nitrite N 9	BOD	*	14 Turb	dity	19	8021B	24	8260/8260B	<u> </u>	30		
5 Nitrate N 10	Alkalinity			uctivity	20	8010/8020	25	8270 B/N or A			↓ V 7n	
31 Metals (As Is, Total, Diss.) Ag, A	, As, B, B	a, Be, Ca	, Cd, Co,	Cr, Cu, Fe, I	Hg, K,	Mg, Mn, Mo,	Na, Ni	, Pb, Sb, Se			V, ZII	
32 TCLP (Specify: volatiles, semi	-volatiles, n	netals, pesti	icides, herb	icides)	33							<u> </u>
34 Other O. P. J. m	HOLC		and	Phead	5	by 8270)	and	Vars	<u> </u>	<u>826 (</u>)
	<u> </u>	(White, Yello			/ Goldenrod Copy	- Client)			•		
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LABORATORY REPORT

Laboratory Services

160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

Heindel & Noyes PO Box 64709, Burlington, VT 05406-4709

Attn: J Noyes

PROJECT: OMYA ORDER ID: 10366

RECEIVE DATE: November 27, 2000 REPORT DATE: December 7, 2000

Enclosed please find the results of the analyses performed for the samples referenced on the attached chain of custody. Different groups of analyses may be reported under separate cover.

All samples were prepared and analyzed by requirements outlined in the referenced methods and within the specified holding times.

All instrumentation was calibrated with the appropriate frequency and verified by the requirements outlined in the referenced methods.

Blank contamination was not observed at levels affecting the analytical results.

6747

Analytical method precision and accuracy was monitored by laboratory control standards which include matrix spike, duplicate and quality control analyses. These standards were determined to be within established laboratory method acceptance limits, unless otherwise noted.

Reviewed by,

Harry B. Locker, Ph.D. Laboratory Director

Enclosures

Page 1 of 2



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT

CLIENT: Heindel & Noyes

PROJECT: OMYA

REPORT DATE: December 7, 2000

ORDER ID: 10366

DATE RECEIVED: November 27, 2000

SAMPLER: DB/CA

Ref. Number: 166047	Site: SS-1(core) 15-18'		Date Sampled: November 21, 2000 Time: NI					
Parameter	Result	<u>Unit</u>	Method	Analysis Date	Analyst			
ortho-Phenyl phenol	168.	mg/L	EPA 642	12/2/00	515			
Ref. Number: 166051	Site: SS-1(core) 0-4"		Date Sampled: Nov	vember 21, 2000 Tim	ie: NI			
Parameter ortho-Phenyl phenol	<u>Result</u>	<u>Unit</u>	<u>Method</u>	Analysis Date	Analyst			
	42.3	mg/Kg	EPA 642	12/2/00	515			

160 James Brown Drive
Williston, Vermont 05495
(802) 879-4333

Project Name: OMYA	Reporting	Address:	H+N	Billing Address:	HAN	·
Endyne Order ID: 1-0 (Lab Use Only) / 03/08 1-1	Company:	: # V Iame/Phone #: J.	46	Phone #	B+CA	
-5		J.	Nohr?	65	T 0820	
Ref # Sample Identification Ma	atrix & &	Date/Time	Sample Containers No. Type/Size	Field Results/Remarks	Anatysis Required	Sample Preservation Rush
166047 SS-1 (lone) 15-18' H	120 X		6 40ml V	94 8760	+ HPLC Z	HELL = NOVE (SI
166048 SS-3 (one) 6-9'			3 4	7/	1 2	- 8260 HCL
1106049 SS-4 (love) 25-28'			2 4	νl)	-HCL -None
1(06050 SS-5 (core) 17-15'	\forall		3 4	4	<u> </u>	-Hcl -vone
1(06,05, SS-1 ((one) 0-4" Slu	odge		4	8260+HPL	c+%Solids	None
16605255-3 (one 0-4"						
16/2033 SS-4 Come 0-4"						
166054 SS-5 (one 0-4"	VV	V	↓ ∨	<u> </u>	V	$ \psi$
				<u> </u>	<u> </u>	
Retriquished by: Date Time 1/2700 1/St	Received by:	// Lr	Deter time	Received by:		Date/Time
New York State Project: Yes No		Requested A	nalyses			
1 pH 6 TKN	11 To	otal Solids	16 Sulfate	21 1664 TPH/FOG	26	8270 PAH
2 Chloride 7 Total P	12 TS	ss	17 Coliform (Specify) 22 8015 GRO	27	PP13 Metals
3 Ammonia N 8 Total Diss. P	13 TI	DS	18 COD	23 8015 DRO		RCRA8 Metals
4 Nitrite N 9 BOD	- -	urbidity	19 8021B	24 8260/8260B	29	
5 Nitrate N 10 Alkalinity	11 1	11	20 8010/8020	25 8270 B/N or Acid	30	~
31 Metals (As Is, Total, Diss.) Ag, Al, As, B, Ba, Be,	, Ca, Cd, Co			o, Na, Nt, Pb, Sb, Se, Si, S	5r, T1, TI, V	Zn
32 TCLP (Specify: volatiles, semi-volatiles, metals, p	pesticides, he	rbicides)	33		<u> </u>	
34 Other						



LABORATORY REPORT

Laboratory Services

160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

Heindel & Noyes PO Box 64709,

Burlington, VT 05406-4709

Attn: J Noyes

PROJECT: OMYA ORDER ID: 10376

RECEIVE DATE: November 27, 2000 REPORT DATE: December 7, 2000

Enclosed please find the results of the analyses performed for the samples referenced on the attached chain of custody. Different groups of analyses may be reported under separate cover.

All samples were prepared and analyzed by requirements outlined in the referenced methods and within the specified holding times.

All instrumentation was calibrated with the appropriate frequency and verified by the requirements outlined in the referenced methods.

Blank contamination was not observed at levels affecting the analytical results.

Analytical method precision and accuracy was monitored by laboratory control standards which include matrix spike, duplicate and quality control analyses. These standards were determined to be within established laboratory method acceptance limits, unless otherwise noted.

Reviewed by,

ved by,

Harry B. Locker, Ph.D. Laboratory Director

Enclosures

Page 1 of 2



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT

CLIENT: Heindel & Noyes

PROJECT: OMYA

REPORT DATE: December 7, 2000

ORDER ID: 10376

DATE RECEIVED: November 27, 2000

SAMPLER: DB/CA

Ref. Number: 166075	Site: SS #1		Date Sampled: Nov	vember 21, 2000 Tin	ne: 11:30 AM
Parameter	Result	<u>Unit</u>	Method	Analysis Date	<u>Analyst</u>
ortho-Phenyl phenol	97.0	mg/L	EPA 642	12/2/00	515
Ref. Number: 166076	Site: SS #2	:	Date Sampled: Nov	vember 21, 2000 Tin	ne: 12:10 PM
<u>Parameter</u>	Result	<u>Unit</u>	<u>Method</u>	Analysis Date	Analyst
ortho-Phenyl phenol	5.32	mg/L	EPA 642	12/2/00	515

160 James Brown Drive

Williston, Vermont 05495 (802) 879-4333

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(802) 879-4333																
Proje	ct Name	· OMYa			Rep	oortin	g Addr	ress:	Billing Address: H5, N							
1 -	ne Order Use Onl		37		-	Company: H 'N Contact Name/Phone #:						Sampler Name: DB 7 C A Phone #: 658 - 0820				
Re (Lab Us		Sample Ide	ntifica	tion	Matrix	G R A B	Б М Р	Date/Time	Samp No.	le Containers Type/Size			esults/Remarks	Analysis Required		
1/0/	2075	- SS #	1		150	V		11.30	2	ILame	<u> </u>		HPLC opp	Horg	None	24/1
7 427.42	<u></u>								2	40ml	Voa		8260			
										1602 F	Plush	`				15/
160	076	55 #	2	i i.	Hao	\overline{V}		12:10	2	11206	<u></u>		FPLC op			1
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											75	requested !	1/ J	PLY _	<u> </u>	
									telan 11/27 3:00 HBL			3/_	 _			
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Reling	uished by:	; 2h_		Date/Time /// 27/	/	ceived l	y;/	16	,/	Data/Time	127	Receive	d by:		Date/	Time
L		ate Project: Yes	N	0				Requested A	nalys	es				· <u>-</u>		
	pH			TKN		11	Total S	olids	16	Sulfate	1	21	1664 TPH/FOG	2	6 8270 PAH	
1	Chłoride	····-	7	Total P		12	TSS		17	Coliform (Specify))	22	8015 GRO	2		
	Ammonia l	N	8	Total Diss. P		13	TD\$		18	COD		23	8015 DRO	2	8 RCRA8 Me	tals
┟┝┷┷┼	Nitrite N		9	BOD		14	Turbid	ity	19	8021B		24	8260/8260B		9	
5	5 Nitrate N 10 Alkalinity 15 Conductivity				20 8010/8020 25 8270 B/N or Acid 30											
31	Metals (A.	s Is, Total, Diss.) Ag	, Al	, As, B, Ba	, Be, Ca,	Cd,	Co, C	Cr, Cu, Fe, H	g,K,	Mg, Mn, M	0, Na	ı, Ni,	Pb, Sb, Se, Si,	Sr, Ti, T	1, V, Zn	.
1		pecify: volatiles,							33							
11	Other															
11 1																

(White, Yellow, Pink Copy - Laboratory / Goldenrod Copy - Client)

53

Tested for OPP Compiled 12/8/00

Date	Location	Result	pН
11/20/00	CDP On-Site	< 40 ppb	
11/21/00	CDP On-Site	< 40 ppb	
11/21/00	Town	< 40 ppb	
11/22/00	LaFlamme	<40 ppb	7.44
11/22/00	O'Keefe	<40 ppb	7.56
11/22/00	Sandillo	<40 ppb	7.61
11/22/00	CDP On-Site	<40 ppb	7.36
11/22/00	Town	<40 ppb	7.51
11/23/00	Town	<40 ppb	7.6
11/24/00	Town	<40 ppb	8.26
11/25/00	Town	<40 ppb	7.49
11/26/00	Town	<40 ppb	7.57
11/27/00	CDP On-Site	<40 ppb	7.43
11/29/00	CDP On-Site	<40 ppb	7.38
11/30/00	CDP On-Site	<40 ppb	7.38
11/30/00	Town	<40 ppb	7.55
12/1/00	CDP On-Site	<40 ppb	7.34
12/1/00	Town	<40 ppb	7.57

⁴⁰ ppb is the lowest concentration detectable by Verichem testing.

Water Sample Data Tested for OPP Compiled 12/8/00

Verichem Analysis	Dolomite Quarry Discharge	: Ha	East Quarry (PIQ) Discharge	pH	11K Tank Process Waler	Ha	Dog Leg Quarry South End (Plant Discharge End)	pΗ	Dog Leg Quarry North End (Plant Intake End)	рН	Primary Discharge West Settling Cell	Hq
Sample Date	< 40 ppb	8.02	< 40 ppb	7,92	< 40 ppb	8,13	650 ррт	9.81	950 ppm	10.04		
11/18/00		7.94	< 40 ppb	8	< 40 ppb	8.1	550 ppm	9.64	950 ppm	10.06		
11/19/00	< 40 ppb	7.77	< 40 ppb	7.84	< 40 ppb	8.04	108 ppm	7.76	391 ppm	9,14		
11/20/00	< 40 ppb	1.77	< 40 ppb	7.8	< 40 ppb	8.12	360 ppm	9,59	518 ppm	9,59		
11/21/00	67 ppb		< 40 ppb	7.74	1 40 PPP		187 ppm	7.77	167 ppm	9.27	11 ppm	7.58
11/22/00		7.00	< 40 ppb	7,79	< 40 ppb	8.1	100 ppm	7.59	153 ppm	7.44	< 40 ppb	7.63
11/23/00	< 40 ppb	7.53	, ,	7.9	- 40 ppo						\	
11/23/00			< 40 ppb	7.71	< 40 ppb	8.1	55 ppm	7.36	145 ppm	7,36	< 40 ppb	7.63
11/24/00	< 40 ppb	7.54	< 40 ppb			8.02	123 ppm	7.54	47 ppm	7.43	6 ррт	7.59
11/25/00			< 40 ppb	7.78	< 40 ppb	8,26	86 ppm	7.52	57 ppm	7.53	9 ppm	7.81
11/26/00					< 40 ppb	8,20 8,12	19 ppm	7.35	37 ppm	7.51	25 ppm	7.49
11/27/00	< 40 ppb	7.8	< 40 ppb	7.86	11 ppm		• •	7.42	51 ppm	7.9	12 ppm	7.82
11/28/00	< 40 ppb	7.37	< 40 ppb	7.84	9 ppm	8.09	40 ppm	7.43	26 ppm	7.55	11 ppm	7.78
11/29/00	< 40 ppb	7,45	< 40 ppb	7,8	≺ 40 ppb	6.08	< 40 ppb	7.45	25 ppm	7,33	6 ppm	7.94
11/30/00	10 ppm	7.86	< 40 ppb	7.86			11 ppm			7.31	< 40 ppb	7,71
12/1/00	< 40 ppb	7,59	< 40 ppb	7.82	< 40 ppb	7.97	< 40 ppb	7.53	14 ppm	1.01	< 40 ppb	* , * 1
12/4/00	< 40 ppb		< 40 ppb		< 40 ppb		< 40 ppb		7 ppm		add nu a	

Ed Nohus of Verichem reports that occasionally a sample will be cloudy one day and clear the next. Plant personnel observe this same effect when much washing is performed in the plant one day and not the next.

Endyne Analysis Sample Date 12/4/00	Dolomile Quarry Discharge 4.75 ppm	East Quarry (PIQ) Discharge	CDP Well	Well # 2 Near Traller Drop Area < 1 ppb	Well # 96-1 Near 500K Tank < 1 ppb	
12/6/00	3.57 ppm	1.96 ppm	< 2 ppb			

- 1. A small amount of material actually passed through flotation and out to the doloralte quarry the night of the spill.
- It took a couple days before the material passed out of the Dolomite quarry in a concentration high enough to measure, i.e. 67 ppb.
- 2. Dog leg samples on 11/20 show less concentration than on 11/21. This may be due to either precipitation or fresh water from relicar rinse.
- 3. Concentration levels in the dog leg quarry continued to fluctuate slightly, again either due to precipitation or fresh water from railcar rinse.
- 4. The 11/22 sample for the west settling cell was taken after the flotation waste/dog leg quarry water mixture began entering the cell.
- 5. On 11/28, water balance issues caused the plant to switch the flotation waste/dog leg quarry water mixture from the west settling cell to the dolomite quarry, resulting in a value on 11/30,
- Flow was switched back to the west settling cell on 12/4. 6. The two west settling cell samples on 11/23 & 11/24 were delayed in transit by Fed Ex to the lab. We question the validity of these two samples due to this delay.

OMYA Verpol Water -- Bacteria and pH Data

Date		g Quarry ntake End		eg Quarry Discharge End	Dolomite Quarry Discharge			
	pН	Aerobic Bacteria (cfu/ml)	pН	Aerobic Bacteria (cfu/ml)	pН	Aerobic Bacteria (cfiv/ml)		
11/20/00	8.22	510000	9.29	12000	8.13	720000		
11/21/00	8.09	2500000	9,73	900	8.07	600000		
11/22/00	8.19	40000	9.12	120000	8.01	630000		
11/23/00	8.03	80000	8.08	520000	7.95	960000		
11/24/00	7.97	370000	7.94	700000	7.91	250000		
11/25/00	7.78	400000	8.05	Data Questionable (10 ⁵ - 10 ⁶)	8.01	300000		
11/26/00	7.58	>1000000	7.86	800000	7.90	270000		
11/27/00	7.53	Data Questionable (10 ⁴ - 10 ⁶)	7.66	Data Questionable (10 ⁵ - 10 ⁶)	7.8	90000		
11/28/00	7.85	250000	7.82	Data Questionable (10 ⁵ - 10 ⁶)	7.95	160000		
11/29/00	7.39	650000	7.55	380000	8.02	190000		
11/30/00	7.91	800000	7.65	860000	7.96	90000		
12/1/00	7.79	8000000	7.67	1250000	7.94	140000		
12/2/00	7.73	4000000	7.67	2000000	7.95	330000		
12/3/00	7.54	1500000	7.5	600000	7.69	390000		

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Wagner, Heindel, and Noyes, Inc.

Consulting Geologists

434 Shelburne Rd. Burlington, VT 05401

	(56) Page No.
)F	

PAGE _____ OF _____

PROJECT: <u>OPM</u>

DATE: <u>II 30 /2000</u>

Determine Permenbilly of Lagoon Sediment

use constant-head. Permeameter test:

Darcy's law: Q = K JH A, where

Q = discharge

K = Permeability

dH = gradient

 $A = C \cos 5 - s \cot i \text{ or } c$ $A = \left(\frac{4.267}{2}\right)^2 \text{ cm} \times 27$ dH dL s odd ment core Q Q dLZ RowL

 $A = \frac{(4.267)^2 \text{cm} \times 29}{2} = 14.23 \text{ cm}^2, \text{ where conditioned or } 4.267 \text{cm}$ 4H = 96.5 cm = constant field 4L = 86.4 cm = constant

Qat = 38.77 milder more rate Qout = 5.44 milder average (See next page)

solve for K: $K = \frac{Q}{A \frac{dH}{dL}}$

- Majimum: $K = \frac{38.77 \text{ ml/den}}{(14.23 \text{ cm}^2) \times (96.5/86.4)} = 2.44 \text{ ml/des/cm}^2$

2.44 mc/day/cm2 +3785 mc/gal x 929 cm2/ft2 +7.48 gal/ft3 = 0.080 ft/day

MAX K = 0.080 felday

- average: $k = \frac{5.44 \text{ mL/day}}{(14.23 \text{ cm}^2) \times (96.5/86.4)} = 0.342 \text{ mL/day/cm}^2$

0.342 ml)day cm2 = 3785 ml)gol ×929 cm2/ft2 = 7.48 gol)ft3 = 0.0112 ft/dog

NERAGE K = 0.0112 ft/day



Wagner, Heindel, and Noyes, Inc.

Consulting Geologists

434 Shelburne Rd. Burlington, VT 05401

		Page No.
PAGE	_ OF	
PROJECT:		<u></u>
DATE:		

SEDIMENT CORE PERMEABILITY TEST CONSTANT-HEAD TEST: FLOW MEASUREMENTS

date/time	Elapsed Days	vol out (mL)	Q (mL/day)
11/29/00 9:20	0.000	0	
11/29/00 10:20	0.042	0	0.000
11/29/00 11:20	0.083	0	0.000
11/29/00 12:40	0,139	0	0.000
11/29/00 14:35	0.219	1.5	18.783
11/29/00 16:45	0,309	3.5	38.769
11/30/00 7:50	0.938	0.1	0,159
TOTAL / AVERA	GE:	5,1	5,440

Z Highest flow me during test

L Accorde Haw inte for entire test



LABORATORY REPORT

Laboratory Services

160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

Heindel & Noyes PO Box 64709,

Burlington, VT 05406-4709

Attn: Jen

PROJECT: OMYA ORDER ID: 10622

RECEIVE DATE: December 13, 2000 REPORT DATE: December 15, 2000

Enclosed please find the results of the analyses performed for the samples referenced on the attached chain of custody. Different groups of analyses may be reported under separate cover.

All samples were prepared and analyzed by requirements outlined in the referenced methods and within the specified holding times.

All instrumentation was calibrated with the appropriate frequency and verified by the requirements outlined in the referenced methods.

Blank contamination was not observed at levels affecting the analytical results.

941

Analytical method precision and accuracy was monitored by laboratory control standards which include matrix spike, duplicate and quality control analyses. These standards were determined to be within established laboratory method acceptance limits, unless otherwise noted.

Reviewed by,

Harry B. Locker, Ph.D. Laboratory Director

Enclosures

Page 1 of 4



160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT SW 8270C

CLIENT: Heindel & Noyes PROJECT: OMYA

SITE: P1Q Quarry

DATE RECEIVED: December 13, 2000 REPORT DATE: December 15, 2000 ANALYSIS DATE: December 14, 2000 ORDER ID: 10622

REFERENCE NUMBER: 166934
DATE SAMPLED: December 12, 2000

TIME SAMPLED: 10:10 AM

SAMPLER: JPW ANALYST: 917

	Result		Result
Parameter	ug/L	Parameter	ug/L
Acenaphthene	< 20.0	I-Methylmaphthalene	< 20.0
Acenaphthylene	< 20.0	2-Methylnaphthalene	< 20.0
Aniline	< 100.	Naphthelene	< 20.0
Anthracene	< 20.0	I-Naphthylamine	< 100.
Azobenzene	< 50.0	2-Naphthylamine	< 100.
Benzidine	< 100.	2-Nitroaniline	< 200.
Benzo(a)anthracene	< 20.0	3-Nitromiline	< 200.
Benzo(b&k)fluoranthene	< 20.0	4-Nitroaniline	< 200.
Benzo(a)pyrene	< 20.0	Nitrobenzene	< 50.0
Benzo(g,h,i)perylene	< 20.0	N-Nitroso-di-n-butylamine	< 50.0
Bis(2-chloroethyl)ether	< 50.0	N-Nitrosodiphenylamine	< 50.0
Bis(2-chloroethoxy)methane	< 50.0	N-Nitrosodimethylamine	< 100.
Bis(2-cthylhexyl)phthalate	< 100.	N-Nitrosodi-n-propylamine	< 100.
Bis(2-chloroisopropyl)ether	< 100.	N-Nitrosopiperidine	< 100.
4-Bromophenyl phenyl ether	< 20.0	Phenanthrene	< 20.0
Butyl benzyl pinhalate	< 100.	Ругене	< 20.0
Carbazole	< 100.	Pyridine	< 100,
4-Chloroaniline	< 50.0	1,2,4-Trichlorobenzene	< 20.0
1-Chloronaphthalene	< 20.0	Benzyl alcohol	< 100.
2-Chloronaphthalene	< 20.0	4-Chloro-3-methylphenol	< 100.
4-Chlorophenyl phenyl ether	< 20.0	2-Chlerophenol	< 50.0
Chrysene	< 20.0	2,4-Dichlorophenel	< 50.0
Dibenzofuran	< 20.0	2,6-Dichlorophenel	< 50.0
Dibenzo(a,h)anthracene	< 20.0	2,4-Dimethylphenol	< 50.0
Di-n-butylphthulate	< 100.	4,6-Dinitro-2-methylphenol	< 500.
1,2-Dichlerebenzene	< 20.0	2,4-Dinitrophenol	< 100.
1,3-Dichlorobenzene	< 20.0	2-Methylphenol (o-cresol)	< 50.0
1,4-Dichlombenzene	< 20.0	3&4-Methylphenol	< 50.0
3,3'-Dichlorobenzidine	< 50,0	2-Nitrophenol	< 100.
Diethyl phthalate	< 100.	4-Nitrophenol	< 100.
Dimethyl phthalate	< 100.	Pentachloropheno!	< 150.
2,4-Dinitrotolucne	< 50.0	Phenol	< 50.0
2,6-Dinitrotoluene	< 50.0	2,4,5-Trichiorophenol	< 100.
Di-n-octylphthalate	< 100.	2,4,6-Trichlorophenol	< 100.
Fluoranthene	< 20.0	Acid Surrogate I	45.%
Pluorene	< 20.0	Acid Surrogate 2	41.%
Hexachiorobenzene	< 50.0	Acid Surrogate 3	81.%
Hexachlorobutadione	< 50.0	Base/Neutral Surrogate 1	81.%
Hexachlorocyclopentadiene	< 200.	Base/Neutral Surrogate 2	81.%
Hexachloroethano	< 50.0	Base/Neutral Surrogate 3	103.%
Indeno(1,2,3-cd)pyrene	< 20.0	UIP's	٥
Isophorone	< 20.0	orthochenylphenol	2,620,
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160 James Brown Drive Williston, Vermont 05495 (802) 879-4333 FAX 879-7103

LABORATORY REPORT SW 8270C

CLIENT: Heindel & Noyes PROJECT: OMYA

SITE: Dolomite Quarry

DATE RECEIVED: December 13, 2000 REPORT DATE: December 15, 2000 ANALYSIS DATE: December 14, 2000 ORDER ID: 10622

REFERENCE NUMBER: 166933
DATE SAMPLED: December 12, 2000

TIME SAMPLED: 9:45 AM

SAMPLER: JPW ANALYST: 917

	Result		Result
Parameter .	ug/L	Parameter	<u>ue/L</u>
Acenaphthene	< 2.0	1-Methylnaphthalene	< 2.0
Accnaphthylene	< 2.0	 2-Methylnaphthalene 	< 2,0
Aniline	< 10.0	Naphthalene	< 2.0
	< 2.0	1-Naphthylamine	< 10.0
Anthrecence	< 5.0	2-Naphthylamine	< 10.01 >
Henzidine	< 10.0	2-Nitroaniline	< 20.0
	< 2.0	3-Nitroaniline	< 20.0
Benzo(a)anthracene	< 2.0	4-Nitrozniline	< 20,0
Benzo(b&k)fluoranthene	< 2.0	Nitrobenzene	< 5.0
Henzo(a)pyrene	< 2.0	N-Nitroso-di-n-butylamine	< 5.0
Benzo(g,h,i)perylene	< 5.0	N-Nitrosodiphenylamine	< 5.0
Bis(2-chloroethyl)ether		N-Nitrosodimethylamine	< 10.0
Bis(2-chloroethoxy)methane	< 5.0 < 10.0	N-Nitrosodinamyamane	< 10.0
Bis(2-ethylhexyl)phthalate	< [0,0] >	N-Nitrosomeridiae	< 10.0
Bis(2-chloroisopropyl) the	< 10.0	N-Nitrosopiperioine Phenanthrenc	< 2.0
4 Bromophenyl phenyl ether	< 2.0		< 2.0
Butyl benzyl phthalate	< 10.0	Pyrene	< 10.0
Carbazole	< 10.0	Pyridine	< 2.0
4-Chloroaniline	< 5.0	1,2,4-Trichiorobenzeac	< 10.0
1-Chloronaphthulene	< 2.0	Benzyl alcohol	< 10.0
2-Chioronaphthalene	< 2.0	4-Chloro-3-methylphenol	
4 Chlorophenyl phenyl ether	< 2.0	2-Chlorophenol	< 5.0
Chrysene	< 2.0	. 2,4-Dichlerophenol	< 5.0
Dibenzofuran	< 2.0	2,6-Dichlorophenol	< 5.0
Dibenzo(a,h)anthuseenc	< 2.0	2,4-Dimethylphenol	< 5.0
Di-n-butylphthalale	< 10.0	4,6-Dinitro-2-methylphenol	< 50.0
1,2-Dichlorobenzene	< 2.0	2.4-Dinitrophenol	< 10.01
1.3-Dichlorobenzene	< 2.0	2-Methylphonol (a-cresol)	< 5.0
1,4-Dichlorobenzene	< 2.0	3&4-Methylphenol	< 5.0
3,3'-Dichlorobenzidine	< 5.0	2-Nitrophenol	< 10.0
Diethyl phthalate	< 10.0	4-Nitrophenol	< 10,0
Directly1 phthalate	< 10.0	Pentachlorophenol	< 15.0
2.4-Dinitratoluene	< 5,0	Phenol	< 5.0
2,6-Dinimatoliume	< 5.0	2,4,5-Trichlorophenol	< 10.0
Di-n-octylphthalute	< 10.0	2.4.6-Trichlorophenol	< 10.0
Fluoranthene	< 2.0	Acid Surrogate 1	46.%
	< 2.0	Acid Surrogate 2	34.%
Fluorene	< 5.0	Acid Surrogate 3	129.%
Hexachlorobenzene	< 5.0	Base/Neural Surregate 1	88.%
Hexachlorobutadiene	< 20.0	Base/Neutral Surrogate 2	85.%
Hexachlorocyclopentudiene	< 20.0 < 5.0	Base/Neutral Surrogate 3	99.%
Hexachloroethanc		Cibs	2.
Indeno(1.2,3-cd)pyrene	< 2.0		75,360
Isophorone	< 2.0	orthophenylphenol	
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160 James Brown Drive Williston, Vermont 05495 (602) 879-4333 FAX 879-7103

LABORATORY REPORT SW 8270C

CLIENT: Heindel & Noyes

PROJECT: OMYA SITE: Dogleg Quarry

DATE RECEIVED: December 13, 2000 REPORT DATE: December 15, 2000 ANALYSIS DATE: December 14, 2000 ORDER ID: 10622

REFERENCE NUMBER: 166935.
DATE SAMPLED: December 12, 2000

TIME SAMPLED: 5:53 PM

SAMPLER: JPW ANALYST: 917

	Result		Result
Parameter .	ue/L	Parameter	ug/L
Accrephthene	< 20.0	I-Methylnaphthalene	< 20.0
Acemphthylene	< 20,0	2-Mcthylnaphthalcnc	< 20.0
Apiline.	< 100,	Naphihalene	< 20.0
Anthracene	< 20,0	1-Naphthylamino	< 100.
Azobenzene	< 50.0	2-Naphthylamine	< 100.
Benzidine	< 100,	2-Nitroaniline	< 200.
Benzo(a)anthracene	< 20.0	3-Nitrosniline	< 200.
Benzo(b&k)fluoranthene	< 20.0	4-Nitroaniline	< 200.
Benzo(a)pyrene	< 20,0	Nitrobenzeno	< 50.0
Benzo(g,h,i)perylene	< 20.0	N-Nitroso-di-n-butylamine	< 50.0
Bis(2-chlomethyl)ether	< 50.0	N-Nitrosodiphenylamine	< 50.0
Bis(2-chloroethoxy)methane	< 50.0	N-Nitrosodimethylamine	< 100.
Bis(2-ethylhexyl)phthalate	< 100.	N-Nitrosodi-n-propylamine	< 100.
Bis(2-chloroisopropyl)ether	< 100.	N-Nitrosopiperidine	< 100.
4-Bromophenyl phenyl ether	< 20.0	Phenauturene	< 20.0
Butyl benzyl phihalate	< 100.	Рутеле	< 20.0
Carbazole	< 100.	Pyridine	< 100.
4-Chloroaniline	< 50,0	1,2,4-Trichlorobenzene	< 20.0
1-Chioronaphthalene	< 20.0	Benzyl alcohol	< 10 0 .
2-Chloronaphthalene	< 20.0	4-Chloro-3-methylphenol	< 100.
4-Chlorophenyl phenyl ether	< 20.0	2-Chlorophenol	< 50.0
Chrysono	< 20.0	2,4-Dichlerophenol	< 50.0
Dibenzoluran	< 20.0	2,6-Dichlorophenol	< 50.0
Dibenzo(a,h)anthracene	< 20.0	2,4-Dimethylphenol	< 50.0
Di-n-butylphthalate	< 100.	4,6-Dinitrn-2-methylphenol	< 500.
1,2-Dichlorobenzene	< 20.0	2,4-Dinitrophenol	< 100.
I,3-Dichlorobenzene	< 20.0	2-Methylphenol (n-cresol)	< 50.0
I,4-Dichlorohenzene	< 20.0	3&4-Methylphenol	< 50.0
3,3'-Dichlorobenzidine	< 50.0	2-Nitrophenol	< 100,
Diethyl phthelate	< 100.	4-Nitrophenol	< 100.
Dimethyl phthalate	< 100.	Pentachlorophenol	< 150.
2,4-Dinitratelucae	< \$0.0	Phenol	< 50.0
2,6-Dinitrotoluene	< 50.0	2,4,5-Trichlorophenol	< 100.
Di-n-netylphthalate	< 100.	2,4,6-Trichlorophenol	< 100.
Fluoranthene	< 20.0	Acid Surrogate 1	48.%
Fluorene	< 20.0	Acid Surrogate 2	40.%
Hexuchlorobenzene	< 50.0	Acid Surrogate 3	99.%
Hexachlorobutadicne	< 50.0	Base/Neutral Surrogate 1	90.%
Hexachlorocyclopentaliene	< 200.	Base/Neutral Surrogate 2	87.%
Hexachloroethane	< 50.0	Buxe/Neutral Surrogate 3	93.%
Indeno(1,2,3-ed)pyrene	< 20.0	UlP3	0.
Isophorone	< 20,0	orthophenylpheno!	283.
Page 4 of 4		* **	

PHOTO LOG: 11/21/00

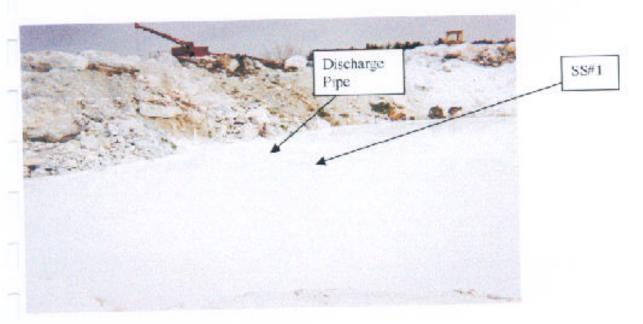


Photo #1: Discharge location into first lagoon.



Photo #2: View of lagoon #1 and southern extent of lagoon #2



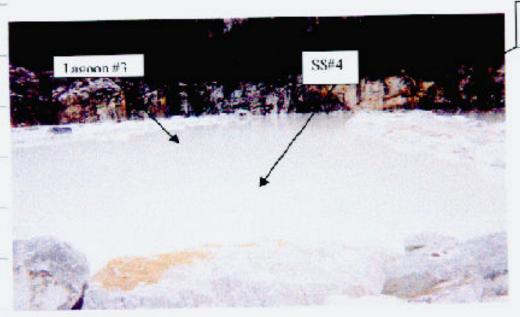
Marble sidewalls of lagoon #2

Photo #3: North/South view of Lagoon #2



Marble sidewalls

Photo #4: Northern extent of lagoon #2.



Marble sidewalls

Photo #5: Lagoon #3



Photo #6: Lagoon #6

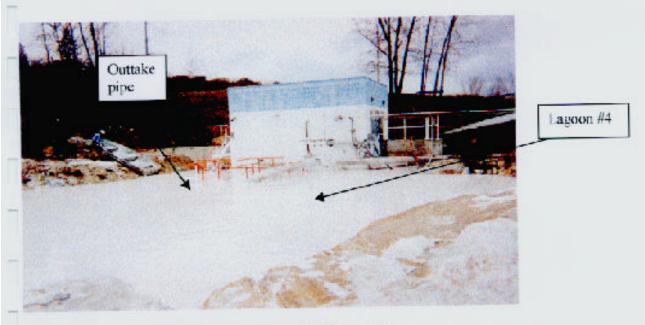


Photo #7: Outtake pipe location at northern end of lagoon #4.