

# New York State Department of Environmental Conservation **MEMORANDUM**

TO:

Jerry Rider

FROM:

Will Welling Welf Velling Love Canal, Site I.D.No. 932020

SUBJECT:

DATE:

March 24, 1995

#### THE LONG-TERM MONITORING PROGRAM

The Long Term Monitoring Program examines hydrological and chemical data from the Love Canal area in order to evaluate the overall effectiveness of the containment system. In 1994, one round of samples was collected from 34 long term monitoring wells that surround the site. Approximately bi-monthly groundwater elevations are taken in six groups of piezometers located around the site.

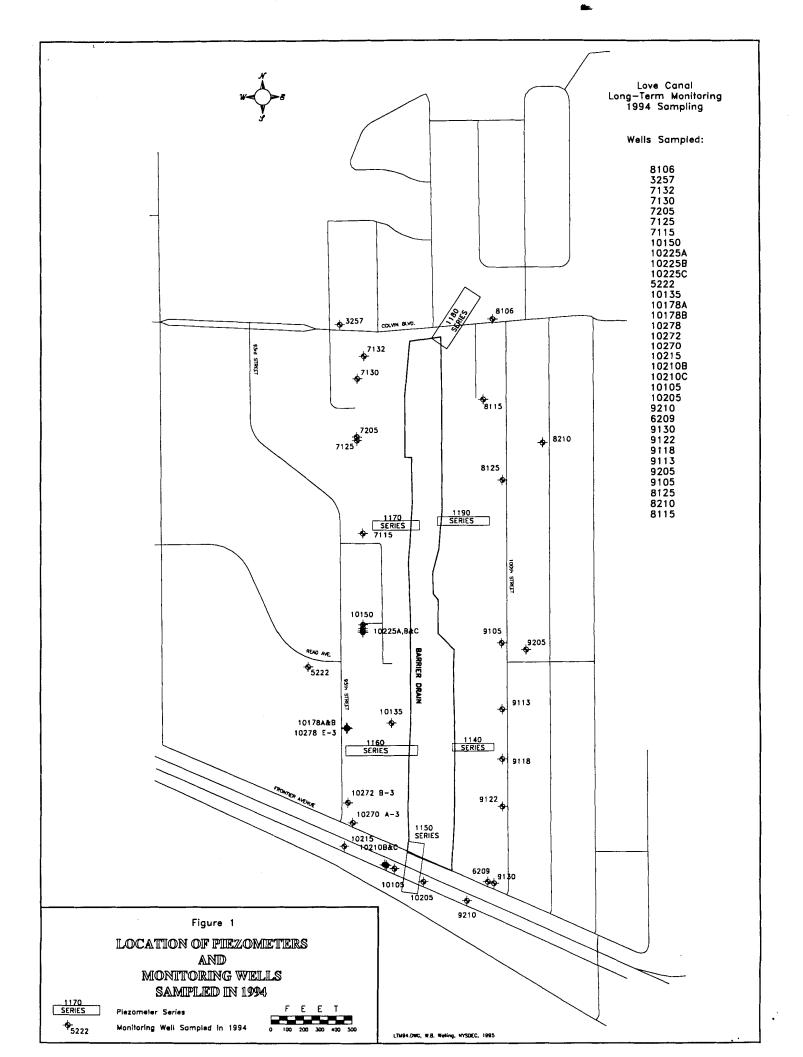
The basic conclusion from the 1994 data is that it is similar to previous data gathered from 1989 to 1993, and that the barrier drain is functioning as designed.

This conclusion is supported by both the hydrological and chemical evidence.

### Results of Groundwater Monitoring **Hydrology**

Readings were taken on six dates during the year from the six series of piezometers that were strategically placed to monitor groundwater elevations along varying cross-sections of the Canal. Figure 1 identifies these piezometers and the shows the locations of monitoring wells sampled for chemical parameters in 1994. Figure 2 is a cross-sectional flow diagram representing groundwater conditions in April 1994 at the 1170 series of piezometers at right angles to the barrier drain. Groundwater flow is toward the leachate collection system. The vertical scale on the cross-section is exaggerated 5 times over the horizontal scale, to aid in interpretation.

No abnormal groundwater conditions were noted during the hydrologic sampling events of 1994 compared to 1993. The barrier drain is capturing all leachate migrating horizontally outward from the Canal, as well as pulling groundwater, which is outside the barrier drain, back toward the drain. Vertical groundwater migration is limited by layers of extremely low permeability clay and glacial till, which underlays all of the site.



580

Figure 2

570

560

550

FEET FROM BARRIER DRAIN

# Results of Groundwater Monitoring Chemistry

In addition to the piezometers, the Love Canal site has a system of overburden and bedrock wells designed primarily for monitoring the chemical quality of groundwater on both sides of the barrier drain. These wells are known as long-term monitoring wells (LTM wells) because they are located in a pattern designed to optimize their usefulness in detecting any failure that might occur in the effectiveness of the barrier drain system over time. In 1993 we began a plan of sampling half of the overburden wells around the perimeter of the site every two years. Half of the LTM wells that are screened in the overburden were sampled in 1993 and the other half were sampled in 1994. Bedrock wells are sampled every year. During 1994, samples were collected from 34 wells and 18 blanks were submitted for quality control purposes. Figure 1 indicates the location of the wells.

In 1994 the New York State Department of Health (DOH) provided analytical services for the long-term monitoring activities at the Love Canal site. Appendix A (attached) is a listing of analytical results. As in previous years, the analytical results were characterized by a predominance of non-detect ("ND") values. When the laboratory has any indication that a compound is definitely present in a sample, that fact is reported. Each time a laboratory analysis is performed there is a level for each compound called the "detection limit." If a compound is present below this level the precise quantity of the compound cannot be accurately determined. In such a case the compound is said to be "present at or below the detection limit" and the result is reported with a "J" qualifier. This reporting protocol was followed by the New York State Department of Health for the 1994 analyses.

In 1994 tentatively identified compounds were not evaluated but they are listed in Appendix A. In 1994 more tentatively identified compounds were detected than in 1993. 152 TICs were listed in 1994 compared to 61 TICs in 1993.

As in previous years, many of the samples showed relatively low levels of phthalates, particularly bis-(2-Ethylhexyl)Phthalate. It was also present in many of the associated blanks. Phthalates are common constituents of plastics used in sampling and laboratory tubing.

Each instance where contaminants were detected at or above the laboratory detection limit is discussed below. The relatively low number of contaminants detected and the fact that the levels detected are close to the detection limits does not indicate a need for action at this time. However, future results will continue to be tracked with special attention to any evidence of a change in groundwater quality.

#### **Well-by-Well Summaries**

The next section of this document describes the wells exhibiting organic chemical contamination at concentrations which are significantly greater than detection levels. Instances of laboratory blank contamination and its impact upon sampling data for an associated round of

samples has been carefully examined by a chemist from Investigation Support Section. This documentation is included as Appendix B. Chemical compounds detected in 1994 are listed immediately after the well number to which they pertain. The units for values presented are micrograms per liter which is analogous to parts per billion. All analyses were performed by the New York State Department of Health.

#### DATA QUALIFIERS FOR RESULTS WHICH FOLLOW

"P" indicates that there is a greater than 25% difference for detected concentrations between the two GC columns in the pesticide/Aroclor target analyte. The lower of the two values is reported.

"E" indicates that the concentration for this compound exceeded the calibration range of the GC/MS instrument for that specific analysis.

"B" indicates that the compound was also found to be present in the laboratory blank for the analytical run. Data before 1993 has not been validated to determine whether such contamination is laboratory related. In 1993 "B" flagged results which are attributable to lab contamination are not presented in the well-by-well listings.

"C" indicates that the value was confirmed either by mass spectrographic analysis or re-examination of the analytical data.

Well #5222 (DOH analysis)	
bis(2-ethylhexyl)phthalate	22
Well #9105 (DOH analysis)	
bis(2-ethylhexyl)phthalate	160
Well #9113 (DOH analysis)	
bis(2-ethylhexyl)phthalate	23
Well #9122 (DOH analysis)	
bis(2-ethylhexyl)phthalate	81
Well #10105 (DOH analysis)	
bis(2-ethylhexyl)phthalate	11

W-II #5000 (DOII analysis)

## Well #10115 (DOH analysis)

bis(2-ethylhexyl)phthalate 12

Well #10205 (DOH analysis)

bis(2-ethylhexyl)phthalate 13

Well #10210B (DOH analysis)

bis(2-ethylhexyl)phthalate 12

Well #10210C (DOH analysis)

19

## Well #10135 (DOH analysis)

acetone

100
91
87
610
24PC
24.4CE
7.5CE
0.15P
61
650E
120
14
51
140
32
91
6000E
. 12
2500
23000

As noted in previous reports, Well #10135 is the only well in the Long Term Monitoring network intentionally installed into an area of known contamination. It monitors groundwater close to the leachate collection system. This well has consistently shown elevated levels of chemical compounds. The results from well 10135 are used as a baseline for comparison with findings from the other monitoring wells. If the Canal were to leak, similar compounds at similar levels would begin to be detected in other wells.

Independent hydraulic evidence indicates that Well #10135 is within the hydraulic influence of the barrier drain system, thus groundwater in the vicinity is likely to be flowing back toward the leachate collection system. The well is approximately 85 feet outside the barrier drain, which is closer to the drain than other long term monitoring wells. (Refer to Figure 1 for its location).

Well #10278 (DOH analysis)

Acetone

33

#### **Summary of 1994 Results:**

- 34 wells were sampled along with 18 blanks.
- A small concentration of acetone was present in samples from wells 10210C and 10278.
- Bis(2-ethylhexyl)phthalate was present in samples from eight wells. Phthalates are common chemical ingredients in plastics and may have been introduced into the sample through contact with plastic tubing and gloves.
- Significant amounts of twenty chemical compounds were found in the sample from well 10135. From year to year, this well is expected to be contaminated since it is within the known contaminated zone. Well 10135 is sampled as a comparison well.
- Groundwater contamination is not migrating from Love Canal based upon the 1994 hydrologic and chemical data.

cc:w/att.

B. Sadowski

G.D. Foster

cc:w/o att.

E. Barcomb

bcc: w/att.

M. Moore

W. Welling B. Loredo

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	TIC's(tentatively identified compounds) - HALOGENS Chlorobenzoic Acid	2-Chbroethykinylether 2-Chbroethykinylether Tetrachbroethene Chbrobenzene Total Chbrotokene Total Dichlorobenzene	Bromdorm Trichbroethene Ditromochloromethane 1,1,2-Tichbroethene is-1,3-Dichbroethene	1,1,2,2-Tetrachloroethane 1,2-Dichloropropane trans-1,3-Dichloropropene	T, T, T-T ICINOTORINANE Carbontetrachloride Bromodichloromethane	Chloroform 1,2-Dichloroethane	1,1-Dichloroethene (total)	Chloroethane Methylene chloride 1 1-Dichloroethane	HALOGEN-VOLATILE Chloromethane Bromomethane Vinvi chloride	TIC's (Tentatively identified compounds) 2.4.6-tribromo-pherod deriva (F1=2 90) Benzene, 1-chioro-2-methyl- (F1=3 95) Benzene, 1-chioro-2-methyl- (F1=3.95) 2.3-deribrono-lumen (F1=5.95) 2.3-deribrono-lumen (F1=5.96) Benzone-methanol, 2-chioro- (F1=6.59) Benzone-methanol, 2-chioro- (F1=8.59) Benzone-audi, 4-chioro- (F1=8.59)	Endosultan Sulfate 4,4*-DDT Methoxychlor	Endnin Aldehyde Endosultan il 44'-DDD	4,4-00E Dieldrin Endrin	Heptachlor Epoxide Endosulfan I		3.3-Dichlorobenzidine Beta&Gamma-BHC(sum of isomers)	Alpha-BHC Hexachlorobenzene Perilachlorobenzene	Chlorophenyl-phenylether     Heromophenyl-phenylether     Heromophenyl-phenylether	ַ מַּמִּ	<ul> <li>4-Chloro-3-Methylphenol</li> <li>Hexachlorocyclopentadiene</li> <li>2,4,6-Trichlorophenol</li> </ul>	4-Chlore Hexach	2,4-Dichlorophenol 1,2,3-Trichlorobenzene 1,2,4-Trichlorobenzene	Hexachloroethane bis(2-chloroethoxy)Methane	1,4-Dichlorobenzene 1,2-Dichlorobenzene bis(2-chlorolsopropyt)Ether	2-Chlorophenol Bis(2-Chloroethyl)Ether 1,3-Dichlorobenzene	B=FOUND IN BLANK DATE SAMPLED LAB HAI OGEN-SEMI-VOJ ATII F	J=BELOW DETECTION LIMITS WELL#	1994 LOVE CANAL LONG TERM MONITORING DATA
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		<b>4</b> 7	<b>X X</b> :	<b>X X</b>	죾죾	<b>4</b> 4	<b>F F</b>	ა¥ Z	<b>*</b>	졲졲	¥ 4	<u> </u>	9 <sup>2</sup> Z	<b>4</b>	¥ 7	<b>F F</b>	., ~		¥ ₹	K K	¥.	Z Z	¥,	¥.			X X	¥.	<b>4</b> 4	<b>F F</b>	졲주	3 BJN	<u> </u>	Ä	12 J /	BJN N	Z Z	2 BJN	X X	Ž,	X X	ĸ.	¥ ¥	, Z	¥ 3	K 4	¥ 4	<b>X</b>	¥,	K 4	독 6 도	, 독	•	•	۲.,
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26 4	130 JN	<b>독</b> 죾	N o	¥.	K K	10 JN	¥ 7	Z Z	죾주	<b>4 4 3</b>	Z Z	¥.	- - - -	죾	ZZ	¥ 4	Z Z	Z Z	Z Z	4 BJN	Ž Z	i i	¥.	Z Z	4 BJN	Z Z	ž,	Z Z	ZZ	¥.	Z Z	Z :	Z Z	ŽĘ,	7 I	4 <sup>2</sup>	, Z	Z Z	Z,	Z 2	<b>.</b> ~	<b>X</b> 7	Z	Z,	ZZ	Z	- - - - -	Ž	Z Z	Z	N W	Ž	05/11/94 DOH	8210	
N N A BJN	ᆩᆩ	죾죾	죾주	<b>4</b>	K K	졲 돆	¥ ?	K =	죾주	<b>A</b>	Ž Z	¥.	<b>K K</b>	짂	Z Z	¥ 4	Z Z	Z 7			Z Z	, Z	짂	Z Z	Z N	Z Z	, N	<b>X</b> ¥	ž Ž	목		3 BJN	ž 7			6 T	Ž Ą	Y Y	Ž,	Z 7	i Z	Z 7	ž Ž	Z,	Ž Z	¥,	o Z	¥,	Z Z	Z,	Z Z	Z,	05/18/94 DOH	9105	
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TiC s(tertatively identified compounds) - NON HALOGENS Methyl Sulfide Methaneithid (RT = 0.91) Methane. Thiobis (RT = 1.42) Silandi, finishity - (RT = 3.39) Ethane. (methythio) - (RT = 3.11) Dimethyl risulfide Dimethyl risulfide Dimethyl (RT = 10.08) Benzene, 1-methyl-3(1-methyl (RT = 10.22) Beryclo(2,2,1) heptan-2-one (RT = 19.21) Molecular sulfur (RT = 28.39) Sulfur, mol. (S8) (RT = 22.61)	Acetone Carbon discrete Cathon discrete 2-Butanone Vinyl acetate Benzene 2-Hezanone 4-Methyl-2-pentanone Total Xylenes 3-Methyl pentane	J-BELOW DETECTION LIMITS WELL# B-FOUND IN BLANK DATE SAMPLED LAB Heplacosane (RT=18.97) 9-octadecenamide, (27- (RT=19.50) Encamide (RT=19.51) Heplacosane (RT=19.51) Heplacosane (RT=19.61) Cotacosane (RT=19.61) 12.6, 10.14, 18.22-lertacosahezane (RT=19.82) Nonecosane (RT=20.83) Oktane, 2.2*(1,4-bufanedy (RT=21.10) Henriscontane (RT=22.05) Cholest-5-en-3-ol (RT=22.05) Trinscontane (RT=22.05) Trinscontane (RT=23.02)
61260 09 61261 09 61262 09 61263 09 61263 09 61260 09	00420 09 30020 09 30022 09 30037 09 00344 09 30031 09 30011 09 30011 09 00392 09 00510 09 30054 09 00545 09	PID UNITCODE
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1994 LOVE CANAL LONG TERM MONITORING DATA

TIC's(tentatively identified compounds) - HALOGENS Chlorobenzoic Acid	Trichloroethene Dibromochloromethane 1,1,2-Trichloroethane isi-1,3-Dichloropropene 2-Chloroethylwinyether Tetrachloroethane Chlorobenzene Total Chlorobenzene Total Dichlorobenzene	1,2-Dictionelhane 1,1-1-Tichbroelhane Carbonieirachloride Bromodichoromelhane 1,2-2-Teirachlorocethane 1,2-Dichloropropane 1rans-1,3-Dichloropropene Bromodom	HALOGEN-VOLATILE Chhoronethane Bromomethane Winyl chloride Chlorothane Methylene chloride 1,1-Dichloroethene 1,1-Dichloroethene 1,1-Dichloroethene 1,2-Dichloroethene 1,2-Dichloroethene 1,2-Dichloroethene 1,2-Dichloroethe	TIC's (rentatively identified compounds) 2.4,6-ti/bromo-phenol deriva (RT=12.90) Benzene, 1-chloro-2-methyl, (RT=3.75) Benzene, 1-chloro-2-methyl, (RT=3.85) 2.3-dichlorool-teme (RT=5.95) Phenol, 4-chloro- (RT=6.36) Benzenemethanol, 2-chloro- (RT=6.56) Benzole acid, 4-chloro- (RT=8.66) Benzole acid, 4-chloro- (RT=8.66)	Big(2-Chioroshir)(Ether Big(2-Chioroshir)(Ether 1,3-Dichlorosherzene 1,2-Dichlorosherzene 1,2-Dichlorosherzene big(2-chioroshorzene big(2-chioroshory))(Ether big(2-chioroshory))(Ether big(2-chioroshory)(Ether big(2-chioroshory)(Ether 1,2,3-Tichlorosherzene 1,2,3-Tichlorosherzene 1,2,3-Tichlorosherzene 1,2-A-Tichlorosherzene 1,2-A-Tichlorosher	J=BELOW DETECTION LIMITS WELL# B=FOUND IN BLANK DATE SAMPLED LAB HALOGEN-SEMI-VOLATILE
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¥,	*********	38888888	6888888888	~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~	~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~	9130 05/25/9 <b>4</b> DOH
죾	*********	38888888	888838888 E	<b>ਸ਼</b> ਸ਼ਸ਼ਸ਼ਸ਼ਸ਼	~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~	9205 05/11/94 DOH
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Z,	<b>33838888</b>	58888888		<del>~</del> ~~~~~	& & & & & & & & & & & & & & & & & & &	10115 05/25/94 DOH
Y Ti	NA 8 8 8 8 8 8 8 8 8 8 8 8 8 8 8 8 8 8 8	5885588 <b>1</b> 8	ND 8 B	NF 2700 JN 1600 JN 470 JN 470 JN 1600 JN 220 JN 21000 JN	888888 <sup>0</sup> ,088888 <sup>0</sup> ,088888 <sup>0</sup> ,08888 <sup>0</sup> ,0888 <sup>0</sup> ,08888 <sup>0</sup> ,08888 <sup>0</sup> ,08888 <sup>0</sup> ,08888 <sup>0</sup> ,08888 <sup>0</sup> ,08888 <sup>0</sup> ,0888 <sup>0</sup> ,08888 <sup>0</sup> ,0888 <sup>0</sup> ,08888 <sup>0</sup> ,0888 <sup>0</sup> ,08888 <sup>0</sup> ,08888 <sup>0</sup> ,08888 <sup>0</sup> ,08888 <sup>0</sup> ,08888 <sup>0</sup> ,08888 <sup>0</sup> ,0888 <sup>0</sup> ,08888 <sup>0</sup> ,0888	10135 06/22/9 <b>4</b> DOH
¥,	*****	5888888	888828888 E	<u> </u>	&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&	10150 06/08/94 DOH
¥	X	8888888	\ 888828888 E	<b></b>	&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&	10176B 05/25/94 DOH
	*****	888 <b>88</b> 88	~	~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~	&&&&&\$\$&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&&	10178A 05/25/94 DOH
죾	N N 1 0 N N N N N N N N N N N N N N N N	8888888			66668866666666666666666666666666666666	10205 05/25/94 DOH
Ž,	XX28X88828			<b>ﯧ</b> ﯧﯧﯧﯧ	* ************************************	10210В 06/15/94 DOH
	********	8888888	888828888 B	****	88888¥88888888888888888888888¥888¥	10210C 06/08/94 DOH
졲	**********	8888888	**	~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~	88888888888888888888888888888888888888	10215 05/25/94 DOH

Orio Swit, C (URCU (H = 3.7e) Discomylene (RT=3.79) Cyclohexane, 1-methyl-3-prop (RT=3.62)	. 0 4 9	Unknown, C9H1802 (RT=3.48) Cyclohexane, 1.2.3,-trimethyl (RT=3.50)	Unknown, C10H20 (H1 = 3.43) Unknown, C10H20 (RT = 3.45) Unknown, C9H160 (RT = 3.46)	2-hexanone, 3-hydroxy-3,5-di (RT=3.33) Unknown, C9H160 (RT=3.43)	3-penten-2-one (RT=3.25) 1.2.3.4.5-pentamethyl-cyclopentane (RT=3.29)	Methane, sulfonyible: (RT=3.05) Acetaldehyde, 2-butenyihydra (RT=3.24) (ZT-2.3.4.5-tetramethyl-3-he (RT=3.25)	Unknown, C10H20 (RT=2.36) Oycobpentane, 1.1.3.4-leftam (RT=2.38)	TICs(lentalively identified compounds) - NON-HALOGEN Cyclohexanol (RT-2.19) Unknown, C10H20 (RT-2.28) 2-hexene, 2.3-dimethyl- (RT-2.33) 3-hexene, 2.3-dimethyl- (RT-2.33)	Benzo(a)Pyrene Indeno(1,2,3-cd)Pyrene Dibenz(a,h)Anthracene Benzo(g,h,j)Penylene	Di-N-Octyl Phihalate Benzo(b)Fluoranthene Benzo(k)Fluoranthene	Benzo (a) Anthracene bis(2-Ethylhexyl)Phthalate	Pyrene Butylbenzylphthalate Chrysene	Carbazole Di-N-Burylphthalate	4 Nitroanline Phenanihrene Anthraene	c.+Diminologierie Diethyphthalate 2-Methyl 4,6-Dinitrophenol N-Nitrosodiphenylamine	Dibenzoluran 4-Nitrophenol 2-4-Dinitrophenol	Acenaphhene Dimethyl Phthalate 2 4-Dintrophenol	3-Nitroaniline Acenaphthylene 2,6-Dinitrotoluene	Naphthalene Benzoic acid 2-Methylnaphthalene	Isophorone 2-Nitrophenol 2,4-Dimethylphenol	4-Methylphenol N-Nitroso-di-propylamine Nitrobenzene	Aniline Benzyl Akohol 2-Methylphenol	NON HALOGEN-SEMI VOLATILE	1,2,3-tric 1,2,4-tric 1,2,4-tric 1,2,3,4-t	Chlorobenzenemethanol(tot) Chlorobenzyl Asorhol Chlorobenzyl Asorhol Benzene, 1-chloro-2-methyl- (RT=16.59) Benzene, 1-chloro-2-methyl- (RT=16.74)	2,3-Dichlorotoluene (RT=19.87) 3,5-Dichlorotoluene (RT=20.33) Dichlorophenol	J=BELOW DETECTION LIMITS WELL# B=FOUND IN BLANK DATE SAMPLED LAB	1994 LOVE CANAL LONG TERM MONITORING DATA
¥ 2 ¥	<b>.</b>	<b>4</b> 4	N S BUN		13 BJN	X		<u></u> * * * * *	8888	0.07 BJ	14 N 8	0.05 0.68 E	5.7 B	3 8 8 8	5 6 8 8 E	5888	52 B 8	388	8 ₹ 8	888	888	N N N	5	<u> </u>	X	· 짂짂짂	9118 05/18/94 DOH	
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N 2 N	<u> </u>	NF 2 BJN	NF 4 BJN		1 K 4	* <del>*</del> * * :	<u> </u>		8888	0.2 BJ	∞ K S	0.09 EJ	2 N S S	3333	S 8 3 8	5668	888	388	888	888	888	8888	5	<b>ፈ</b> ዷዷዷ	X	ጃ <b>ጃ</b> ጃ	9205 05/11/94 DOH	
7 BJN 2 BJN	4 4 4	<b>4</b> 4	NE BIN		5 <b>4</b> 1 4	<b>.</b>	<u> </u>	N 3 BJN	0.06 J 0.1 J 0.1 J	0.05 J	<sub>9</sub> Β ₹	0.05 0.05 0.05 0.05 0.05 0.05 0.05 0.05	0.4 N	5888	5 6 S S	5 6 6 8	588	588	8 8 8	888	888	8 8 8 8	5	<u> </u>	K	<u> </u>	9210 05/25/94 DOH	
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N S N	ᆩᆩᆩ	<b>주</b> 주	<b>X</b>	폮 닦 .	, F, F	<b>.</b>	<b>.</b>	N S N	88888	568	13 8 8 80	5 S S S	0.4 N B	5 6 6 E	E E E	888	§ 8 8	588	8 8 8 8	3 <b>3</b> 8 8	888	5	į	<u> </u>	<b>.</b>	¥ ጙ ጙ	10115 05/25/94 DOH	
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Chancown (H = 14.54)  1-propartone: I, phenyl, (related (RT=14.57) Unknown CHO (RT=14.59)  Benzole acid, 14-benzoy, (rel (RT=14.62) 1.4-dibutoxybutane-related (RT=15.17) Octadecanoic acid (RT=15.29) Hevadecanoic acid (RT=15.20) Phenol, 4.4-(1-methylethyl (RT=15.30) Phenologian (RT=17.41) defiv of 1.2-propanediol, mo (RT=17.41) defiv of 1.2-propanediol, mo (RT=17.59) Pentacosane (RT=17.64) Phosphine oxide, diphenyl, (RT=18.38) Benzene, bis-(phenoxymethyl) (RT=18.08) Unknown (RT=18.09) Phosphine oxide, injohenyl-(RT=18.12) Benzene, bis-(phenoxymethyl) (RT=18.15) Phosphine oxide, diphenyl (RT=18.15) Phosphine oxide, diphenyl (RT=18.15)	Unknown hydrocarbon (RT-8.85) Unknown hydrocarbon (B-7.1 MW (RT-9.32) Unknown hydrocarbon (B-7.1 MW (RT-9.40) Dodecanose acid (RT-9.68) Benzole acid:methy-2-(Pg (RT-9.73) 1.4-butlanedoi, monobenzoate (RT-10.03) Unknown hydrocarbon (B-7.1 (RT-10.13) Unknown hydrocarbon (B-7.1 (RT-10.12) Unknown hydrocarbon (B-7.1 (RT-10.27) 1.4-butlanedoi, monobenzoate (RT-11.52) Unknown benzole acid (RT-11.61) 1.4-butlanedoi, monobenzoate (RT-11.55) Tefradecanole acid (RT-11.61) 1.5-butlanedoi, monobenzoate (RT-11.55) 1.6-butlanedoi, monobenzoate	Berzole acid (RT=6.25) Berzole acid (RT=6.30) Beryole(3.30)cct (12)-en-3- (RT=6.30) I-melfyl-2-yanov-2-piperidai (RT=6.30) Unknown C10H20 (RT=6.41) Unknown C9H180 (RT=6.59) Unknown, C9H180 (RT=6.59) Unknown, C9H180 (RT=6.62) Dimelhyl letrasulphide (RT=6.62) I-vyclopenten-1-y-propane (RT=6.62) I-vyclopenten-1-y-propane (RT=6.71) I-vyclopenten-1-y-propane (RT=6.72) Berzole acid (RT=6.27) Berzole acid (RT=6.27) Berzoneacetic acid (RT=6.26) Berzoneacetic acid (RT=6.27) Berzoneacetic acid (RT=7.02) Nonanocic acid (RT=7.02) Nonanocic acid (RT=7.02) Nonanocic acid (RT=7.03) Berzaldehyde, 4-hydroxy-3-me (RT=7.87) Berzaldehyde, 4-hydroxy-3-me (RT=8.32) Unknown hydroathoria-1-1, MW (RT=8.67)	J=BELOW DETECTION LIMITS WELL# B=FOUND IN BLANK DATE SAMPLED LAB  Cyclohexane, 1-ethyl-2-methy (RT=3.76) Cyclohexane, 1-methyl-2-prop (RT=3.82) 4-hepten-3-one, 2.6-dimethyl (RT=3.82) 5-nonene-4-one (RT=3.83) Benzene, 1-ethyl-2-methyl-(RT=3.89) Trisuffice, dimethyl-(RT=3.89) Trisuffice, dimethyl-(RT=4.32) Unknown, C9H1802 (RT=4.51) 2.4-hosadicine, 3.4-dimethyl-(RT=4.2) 1-hexanol, 2-ethyl-(RT=4.84) 2.3-dimethyl-2-cyclopenten-1 (RT=4.93) Unknown hydrocarbon (RT=6.31) Ethanone, 1.2-dt-2-furamyl-2 (RT=5.64) Phosphoric acid, 1-ethyl-(RT=5.80) Chanonic acid, 2-ethyl-(RT=5.80) Chanonic acid, 2-ethyl-(RT=5.80)	1994 LOVE CANAL LONG TERM MONITORING DATA
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Suffur, mol. (S8) (RT=32.61)	Molecular suffur (RT=28.39)	Bicycloi2 2 1 heptan-2-one (RT=19 21)	Benzene, 1-methyl-3-(1-methyl (RT=18.02)	Dimethyl Disuffide (RT=10.08)	Dimethyl Insulfide	Ethane, (methythio)- (RT=3.11)	Silanoi, trimetriyi- (FI =2.39)	Citarol trimothy (DT-2 20)	Methane, Thiobis (RT±1.42)	Methanethiol (RT=0.91)	Methyl Sunide	TIC's (tentalively identified compounds) - NON HALOGEN	3-Methyl peniane	Total Aylenes	Styreine	Shrana	Ethylbenzene	Toluene	4-Methyl-2-pentanone	z-Hexanone	Contraine	Benzene	Vinyl acetate	2-Butanone	Carbon disulfide	Acetone	NON HALOGEN-VOLATILE	Oxirane, 2,2-[1,4-butanedly (HT=25.02)	i di	Tetratriacontana (RT-23.02)	Trinacontane (RT-23 04)	Dolnacontane (BT=22.22)	Cholest-5-en-3-ol (RT=22.06)	Hentnacontane (RT=21.50)	Oxirane, 2,2'-[1,4-butanedly (RT=21.10)	Triacontane (RT=20.83)	Nonacosane (RT=20.21)	2,6,10,14,18,22-tetracosahexane (RT=19.82)	Octamballe (MIE19.01)	Octobrano (DT_10 61)	Hentarlecane (BT=19.60)	Encamide (RT=19.51)	9-octadecenamide, (Z)- (RT=19.50)	Heplacosane (RT=18.97)	LAB	B=FOUND IN BLANK DATE SAMPLED	J=BELOW DETECTION LIMITS WELL#		1994 LOVE CANAL LONG TERM MONITORING DATA
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TIC's (tentatively identified compounds) - HALOGENS Chlorobenzoic Acid	1.2 Dichloroethane (Idal) Chlordom 1.2 Dichloroethane 1.1.1-1 richloroethane 1.1.1-1 richloroethane 1.1.1.2 richloroethane 1.1.2 richloroethane 1.1.2 richloroptane 1.2 Dichloroptane 1.2 Dichloroptane 1.2 Dichloroptane 1.2 Tichloroethane 1.2 Dichloroethane	HALOGEN-VOLATILE Chloromethane Bornomethane Vinyl chloride Chloroethane Melitylene (chloride 1,1-Dichloroethane	TIC's (Tentatively identified compounds) 2,46-th/bromo-phenol deriva (RT=12.90) Benzene, 1-chloro-2-melthy-(RT=3.75) Benzena, 1-chloro-4-melthy-(RT=3.85) 2,3-dichlorotoluene (RT=5.96) Phenol, 4-chloro- (RT=6.36) Benzenemelhanol, 2-chloro- (RT=6.85) Benzole add, 4-chloro- (RT=8.85) Benzole add, 2-chloro- (RT=8.85)	HALOGEN-SEMI-VOLATILE 2.Chlorophenol Bis(2.Chlorophenol Bis(2.Chlorophenol 1.2.Dichloroberzene 1.2.Dichloroberzene 1.2.Dichloroberzene 1.2.Chlorophenol 1.2.3.Tichloroberzene 1.2.4.Tirchloroberzene 1.2.4.Tirchloroberzene 1.2.4.Tirchloroberzene 1.2.4.Tirchloroberzene 1.2.4.Tirchloroberzene 1.2.4.Tirchloroberzene 1.2.4.Tirchloroberzene 1.2.4.Tirchloroberzene 1.2.4.Tirchlorophenol 1.2.4.Tirchlorophenol 1.2.4.Tirchlorophenol 1.2.4.Tirchlorophenol 1.2.4.Tirchlorophenol 1.2.4.Tirchlorophenol 1.2.4.Tirchlorophenol 1.2.4.Tirchlorophenol 1.2.4.Tirchlorophenol 1.3.Dichloroberzene 4.Chlorophenol 2.Chlorophenol 2.Chlorophenol 2.Chlorophenol 3.Dichloroberzene 4.Chlorophenol	1994 LOVE CANAL LONG TERM MONITORING DATA J-BELOW DETECTION LIMITS WELL# B-FOUND IN BLAKK DATE SAMPLED
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•	2.4 - chruran, octahydro. (R1 = 3.50) Cyclohezanol, 3-methy (R1 = 3.62) Butane, 2-chloro-2-methyl. (R1 = 3.63) Unknown, C10H20 (R1 = 3.78) Discamylene (R1 = 3.78) Cyclohezane, 1-methyl-3-prop (R1 = 3.62)	Unknown, C10H20 (H1=3.45) Unknown, C9H160 (R1=3.46) Unknown, C9H160 (R1=3.46) Cyclohexane, 1,2,3,-trimethyl (R1=3.50)	2-nevanone, 3-nydroxy-3,5-di (H = 3.33) Unknown, C9H160 (RT = 3.43) Unknown, C10H20 (RT = 3.43)	T TD:	Methane, sulfonytbis- (RT=3.05) Acetaldehyde, 2-butlenythydra (RT=3.24) (Z)-2.3.4.5-fortamethyd-3-he (RT=3.25)	3-hexene, 2,3-dimethyt- (RT=2,36) Unknown, C10H20 (RT=2,36) Cycloperfane, 1,1,3,4-lefram (RT=2,38)	Cyclohexanic (RT-2.19) Unknown, C10H20 (RT-2.28) 2-hexane, 2,3-dimethyl- (RT-2.33)	Benzo(g,h,j)Perylene	Indenot (2, 3-cd) Pyrene Dibenz (a, h)Anthracene	Benzo(k) Fluoranihene Benzo(k) Fluoranihene Benzo(a) Borene	ois(z-einyinexyi)Phinalale Di-N-Octyl Phihalate Botzoth Eliopanhago	Chrysene Benzo (a) Anthracene	Pyrene Butylbenzylphthalate	Di-N-Butylphthalate Fluoranthene	Anthracene Carbazole	A-Nitroaniine  Phenanthrene	Dietry Iphthalate 2-Methyl 4,6-Dinitrophenol N. Närssockishandhaniss	4-Nitrophenol 2,4-Dinitrotoluene	2,4-Dintrophenol Dibenzofuran	Acenaphhene Dimethyl Phihalate	Acenaphthylene Acenaphthylene 2 6-Dipitrolohene	2-Methylaphthalene 3-Nitroanline	A,4-Unneurypneno: Naphthalene Banzok soid	Isophorone 2-Nitropherol	N-Nitroso-di-propylamine Nitrobenzene	2-Methylphenol 4-Methylphenol	Anilina Benzyl Akohol	NON HALOGEN-SEMI VOLATILE	Benzene, 1,2,4-inchloro-3-methyl (RT=23,03) Benzene, 1,2,3,4-tetrachloro (RT=24,18)	Benzene, 1,2,4-trichloro-3-methyl (RT=20.83) Benzene, 1,2,4-trichloro-3-methyl (RT=22.85)	Chlorobenzyl Alcohol  Benzene, 1-chloro-2-methyl- (RT=16.59)	2.3-Uchlorololuene (RT=19.87) 3,5-Dichlorololuene (RT=20.33) Dichlorophenol Chlorobaranemethanol(fot)	TE SAMPLED	MONITORING DATA	
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Hexanediok acid, diocryl ester (RT=16.91) Tetracosane (RT=16.83) deriv of 1.2 propanaciol, mo (RT=17.41) deriv of 1.2 propanaciol, mo (RT=17.50) deriv of 1.2 propanaciol, mo (RT=17.59) Pentacosane (RT=17.64) Phosphine oxide, diphenyt- (RT=17.88) Bentane, bs-(phenoxymethy) (RT=17.89) Oxirane, 2.2-(1.4 butanediy (RT=18.08) Unknown (RT=18.09) Unknown (RT=18.09) Phosphine oxide, tiphenyt- (RT=18.12) Bentane, bs-(phenoxymethyt) (RT=18.16) Phosphine oxide, tiphenyt- (RT=18.16) Phosphine oxide, diphenyt (RT=18.51)	Hexadecanoic acid (RT=13.45) Sulfur, mol. (Sib) (RT=14.35) 1.4-dibutoxybutane-related (RT=14.43) 1-ocidadecano (TR=14.48) Tinuvin P (RT=14.51) Unknown (RT=14.51) 1-propanone, 1-phenyl-(related (RT=14.57) Unknown (RT=14.54) 1-propanone, 1-phenyl-(related (RT=14.57) Unknown (RT=14.54) 1-dibutoxybutane-related (RT=15.17) Octadecanoic acid, thenyl est (RT=15.30) Phenol, 4.4-(1-methylethyl (RT=15.33) Octadecanoic acid, buyl est (RT=15.33) Octadecanoic acid, buyl est (RT=16.82)	1.4-butanediol, monobenzoate (RT=10.03) Unknown hydrocarbon, B+71 (RT=10.19) Unknown hydrocarbon, B+71 (RT=10.19) Unknown hydrocarbon, B+71 (RT=10.27) 1.3-propanediol, monobenzoate (RT=11.12) Benzock acid, 1-melhylethy (RT=11.22) Benzock acid, 1-melhylethy (RT=11.23) 1.4-butanediol, monobenzoate (RT=11.22) Unknown benzote acid (RT=11.55) 1-teradecanolo acid (RT=11.51) 1-teradecanolo acid (RT=11.51) 1-poxadecanol (RT=12.73) 3.9-dezatitrycolof (3.0.0(3. (RT=13.38))	droc droc 4	Unknown C10+20 (Řͱ6, 41) Unknown, C9H 80 (Řͱ6, 58) 1*cyckpolarien*-1-vlypotpane-(Řͱ6, 68) 2,6-octadiene, 2,4-dimelhyl- (Řͱ6, 71) 1*cyckbutlen*-1-yl-butlane-1, (Řͱ6, 71) 1*cyckbutlen*-1-yl-butlane-1, (Řͱ6, 74) Benziote acid (Řͱ6, 89) Benzensecelic acid (Řͱ6, 95) 2-duranaceik acid. alpha. (Řͱ7, 02) Nonanoic acid (Řͱ7, 703)	Cycloheanaire,	1994 LOVE CANAL LONG TERM MONITORING DATA J-BELOW DETECTION LIMITS WELL# B-FOUND IN BLANK DATE SAMPLED LOSS CONTROL OF THE SAMPLED CONTRO
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Sultur, mol. (S8) (RT=32.61)	Molecular sultur (RT=28.39)	Bicyclo[2.2.1]heptan-2-one (RT=19.21)	Benzene, 1-methyl-3-(1-methyl (RT=18.02)	Dimethyl Disuttide (RT=10.08)	Dimethyl trisumide	Custing, (manymino)- (mi =0.11)	Thank (methythio) (DT-3-11)	Silanol trimethyl- (AT=2.39)	Methane, Thiobis (RT=1.42)	Methanethiol (RT=0.91)	Methyl Sulfide	TIC's(tentatively identified compounds) - NON HALOGEN	3-Methyl pentane	Total Xytenes	Styrene	Elhylbenzene	Toluene	4-Methyl-2-pentanone	2-Hexanone	Benzene	Vinyt acetate	2-Butanone	Carbon disulfide	Acetone	NOW HAT OCEN, WOLATHE	Oxirane, 2,2'-[1,4-butanediy (RT=25.02)	Tetratriacontane (RT=23.92)	Tritriacontane (RT=23.04)	Dotnacontane (RT=22.22)	Cholest-5-en-3-ol (RT=22.06)	Hentriacontane (RT=21.50)	Oxirane, 2,2'-[1,4-butanediv (RT=21.10)	Triacontane (RT=20.83)	E,o, 10:14, 10,66 tenacosalexale (F1 = 19.06)	Octowoode (D1=19.01)	Polacocano (PT-10 61)	Endederane (DT=10.60)	S-comids (DT=10.51)	Decladesenamida (7)- (BT-10.50)	Jeplemene (BT-19 07)	æ	B=FOUND IN BLANK DATE SAMPLED	LEELOW DETECTION LIMITS WELL*	1994 LOVE CANAL LONG TERM MONITORING DATA
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'ilC's(tentatively identified compounds) - HALOGENS Chlorobenzoic Acid	Total Chlorotoluene Total Dichlorobenzene	Tetrachloroethene Chlorobenzene	1,1,2-Trichloroethane cis-1,3-Dichloropropene	Trichloroethene Dibromochloromethane	mans-1,3-Denoropropene	1,2-Dehloropropane	Bromodichloromethane	1,1,1-Trichloroethane Carbontetrachloride	1,2-Dichloroethane	1,2-Dichloroethene (total)	1,1-Dichloroethene	Chlorethane Methylene chloride	Bromoethane Vinyl chloride	HALOGEN-VOLATILE Choromethane			benzene, 1-cnioro-4-metnyl- (H ) =3.85) 2,3-dichlorotoluene (HT=5.96)	2,4,6-tribromo-phenol deriva (RT=12.90) Benzene, 1-chloro-2-methyl- (RT=3.76)	Methoxychlor	Endosulfan Sulfate 4,4'-DDT	Endosullan ii 4,4'-DDD	Endrin Aldehyde	4,4 COCE	Endosullan I	Aldrin Hanlachlor Enoxide	Deta-BHC Heptachlor	3,3-Dichlorobenzidine Beta&Gamma-BHC(sum of isomers)	Hexachlorobenzene Pentachlorophenol	Alpha-BHC	4-Chlorophenyl-phenylether	2-Chloronaphthalene	2.4.6-Trichkorphenol	4-Chloro-3-Methylpheno! Hexachlomous-bonentadione	4-Chloroaniline	1,2,3-Trichlorobenzene	bis(2-chloroethoxy)Methane 2,4-Dichlorophenol	bis(2-chloroisopropyl)Ether Hexachloroethane	1,2-Dichlorobenzene	i.3-Dichkorobenzene		LAB	J=BELOW DETECTION LIMITS WELL#	
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TIC's (entatively identified compounds) • NONHALOGEN Cyclobeanol (RT-2.19) Unknown. C10H20 (RT-2.28) + hexene. 2.3-dimethy. (RT-2.35) Jhexene. 2.3-dimethy. (RT-2.35) Unknown. C10H20 (RT-2.36) Unknown. C10H20 (RT-2.36) Unknown. C10H20 (RT-2.36) Cyclopeatane. 1,13.4 steffam (RT-2.28) Methane, suffonylbis. (RT-3.26) Cyclopeatane. (RT-3.28) L. 3.4.5-pentiamethyl-3-the (RT-3.29) L. 3.4.5-pentiamethyl-3-the (RT-3.39) Unknown. C10H20 (RT-3.46) Unknown. C10H20 (RT-3.46) Unknown. C9H160 (RT-3.46) Unknown. C9H160 (RT-3.46) Unknown. C9H160 (RT-3.50) C2-bitamethyl-(RT-3.50) C3-bitamethyl-(RT-3.50) C3-bitamethyl-(RT-3.50) Discamylene (RT-3.79) Unknown. C10H20 (RT-3.79) Unknown. C10H20 (RT-3.39) Discamylene (RT-3.39) Discamylene (RT-3.39) Discamylene (RT-3.39)	2.3-Dichlorototuene (RT=19.67) 3.5-Dichlorototuene (RT=20.33) Dichlorophenol (RT=20.33) Dichlorophenol (RT=20.33) Dichlorophenol (RT=20.33) Dichlorophenol (RT=20.33) Benzene, 1.2-Alrichloro-3-methyl (RT=20.33) Benzene, 1.2.4-Richloro-3-methyl (RT=20.33) Benzene, 1.2.4-Richloro-3-methyl (RT=20.33) Benzene, 1.2.3-Richloro-3-methyl (RT=20.33) Benzenehhitene  3-Alirophenol A-Nitrophenol A-Nitrophenol Benzelorian A-Nitrophenol Benzelorian A-Nitrophenol Benzelorian A-Nitrosodishenylamine Benzelorian Benzelorian-Ben	1994 LOVE CANAL LONG TERM MONITORING DATA J-BELOW DETECTION LIMITS WELL! B-FOUND IN BLANK DATE SAMPLED
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Tinum P (RT=14.54) Unknown (RT=14.54) Unknown (RT=14.53) 1-progenome. 1-phenyl-(related (RT=14.57) Unknown CHO (RT=14.59) Beruck actd, 4-berucyl-(rel (RT=4.82) 1.4-dibutoxybutane-related (RT=15.31) Cotadecanole actd (RT=15.30) Prenol. 4.4-(1-methylethyl (RT=15.30) Prenol. 4.4-(1-methylethyl (RT=15.30) Prenol. 4.4-(1-methylethyl (RT=15.30) Obtadecanole actd, bufy est (RT=15.30) Obtadecanole actd, bufy est (RT=16.9) Tetracosane (RT=16.93) denv of 1.2-propanediol, mo (RT=17.50) denv of 1.2-propanediol, mo (RT=17.50) denv of 1.2-propanediol, mo (RT=17.50) denv of 1.2-propanediol, mo (RT=17.80) Prosphine oxide, diphenyl (RT=17.80) Prosphine oxide, diphenyl (RT=18.90) Unknown (RT=18.09) Prosphine oxide, triphenyl (RT=18.12) Beruces, bis-(phenoxymethyl) (RT=18.12) Brozzosane (RT=18.32) Phosphine oxide, diphenyl (RT=18.15) Phosphine oxide, diphenyl (RT=18.15) Phosphine oxide, diphenyl (RT=18.15) Phosphine oxide, diphenyl (RT=18.15)	Unknown hydrocarbon (8-71 MW (RT-9-32) Unknown hydrocarbon (8-71 (RT-9-40)) Dodecanic acid (RT-9-68) Benzoic acid (RT-9-68) 1,4-butanedol, monobenzoate (RT-9-73) 1,4-butanedol, monobenzoate (RT-10-03) 1,4-butanedol, monobenzoate (RT-10-03) 1,4-butanedol, monobenzoate (RT-11-12) 1,4-butanedol, monobenzoate (RT-11-12) 1,4-butanedol, monobenzoate (RT-11-12) 1,3-propanedol (RT-11-13) 1,4-butanedol, monobenzoate (RT-11-15) 1,5-detadecanoic acid (RT-11-12-73) 1,5-detadecanoic acid (RT-11-12-73) 1,5-detadecanoic acid (RT-11-13-14) 1,4-debutoxybutane-related (RT-14-43) 1,4-debutoxybutane-related (RT-14-43)	2)-en-3- (RT= 2)-en-3- (RT= RT=6.61) RT=6.62) de (RT=6.62) propane- (RT imethyl- (RT= Litane-1, (RTE- RZ) (RT=6.95) (RT=6.95) Apha (RTE- 7.03) Apha (RTE- 7.03) Apha (RTE- 4.04) Apha (RTE- 4.05) Apha (RTE- 4.05)	Oyclohexane, 1-ethyl-2-methy (RT-3.76) Cyclohexane, 1-methyl-2-prop (RT-3.82) A-pepin-3-one, 2-edimethyl (RT-3.82) A-pepin-3-one, 2-edimethyl (RT-3.82) Benzene, 1-ethyl-2-methyl-(RT-3.89) Trisulide, dimethyl-(RT-3.89) Benzene, 1-2,4-trimethyl-(RT-3.89) Benzene, 1-2,4-trimethyl-(RT-4.32) Unknown, CSH1962 (RT-4.51) 2-4-hexanol, 2-ethyl-(RT-4.84) Unknown hydrocarbon (RT-4.83) Unknown hydrocarbon (RT-4.83) 2-3-dimethyl-2-cycloperten-1 (RT-5.31) Ethamone, 1-2-di-2-trimyl-2 (RT-5.64) Phosphocia cadd, triethyl ester (RT-5.72) Hexanoc cedd, 2-ethyl-(RT-5.80) Cotlanoic acid (RT-6.12) Benzolic acid (RT-6.12) Benzolic acid (RT-6.12)	1994 LOVE CANAL LONG TERM MONITORING DATA J-BELOW DETECTION LIMITS WELL* B-FOUND IN BLANK DATE SAMPLED
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TIC's(tentatively identified compounds) - NON HALOGEN Methyl Sulfide Methyl Sulfide Methanethol (RT=0.91) Methanethol (RT=0.91) Methane, Thiobis (RT=1.42) Silanol, timethyl- (RT=3.39) Ethane, (methythio)- (RT=3.31) Dimethyl Insulfide Dimethyl Insulfide Dimethyl Insulfide Dimethyl Insulfide (RT=10.08) Benzene, 1-methyl-3-(1-methyl (RT=18.02) Bicyclof(2.2.1)heplan-2-one (RT=19.21) Molecular sulfur (RT=28.38) Sulfur, mol. (88) (RT=32.61)	NON HALOGEN-VOLATILE Aceitone Carbon disulfide 2-Butanone Vinyl acetate Benzene 2-Hexanone 4-Methyl-2-pentanone Total Xylenes 3-Methyl pentane	Heptacosane (RT=18.97) 9-octadecenamicia, (Z)(RT=19.50) Encamide (RT=19.51) Heptadecane (RT=19.51) Heptadecane (RT=19.51) Octacosane (RT=19.61) 2.6, 10.14, 18.22-etracosahexane (RT=19.82) Nonacosane (RT=20.21) Triacontane (RT=20.85) Oxirane, 2.2-(1.4-butanedly (RT=21.10) Hentiacontane (RT=22.06) Cholest-5-en-3-of (RT=22.06) Cholest-5-en-3-of (RT=23.04) Tritriacontane (RT=23.04) Tritriacontane (RT=23.04) Tritriacontane (RT=23.04) Tritriacontane (RT=23.04) Tritriacontane (RT=23.04)	1994 LOVE CANAL LONG TERM MONITORING DATA  J-BELOW DETECTION LIMITS WELL#  B-FOUND IN BLANK DATE SAMPLED  LAB
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