#### **Exhibit 6**

**Current BCEE 2014 Detection Limits** 

Test America MEthod 625 RC

Aneyres	(2) The second s		uall
-Bromophenyl phenyl ether	101-55-3	10	ug/L
-Chloro-3-methylphenol	59-50-7	10	ug/L
-Chloroaniline	106-47-8	10	ug/L
-Chlorophenyl phenyl ether	7005-72-3	10	ug/L
-Methylphenol	106-44-5	10	ug/L
-Nitroaniline	100-01-6	20	ug/L
-Nitrophenol *	100-02-7	30	ug/L
Acenaphthene	83-32-9	10	ug/L
Acenaphthene-d10	15067-26-2		ug/L
Acenaphthylene	208-96-8	10	ug/L
Acetophenone	98-86-2	10	ug/L
Aniline	62-53-3	10	ug/L
Anthracene	120-12-7	10	ug/L
Atrazine	1912-24-9	10	ug/L
Benzaldehyde	100-52-7	10	ug/L.
Benzidine	92-87-5	10	ug/L
Benzo[a]anthracene	56-55-3	1	ug/L
Benzo[a]pyrene	50-32-8	1	ug/L
Benzo[b]fluoranthene	205-99-2	1	ug/L
Benzo[g,h,i]perylene	191-24-2	10	ug/L
Benzo[k]fluoranthene	207-08-9	1	ug/L
Benzoic acid	65-85-0	50	ug/L
Benzonitrile	100-47-0	20	ug/L
Benzyl alcohol	100-51-6	10	ug/L
Bis(2-chloroethoxy)methane	111-91-1	10	ug/L
Bis(2-chloroethyl)ether	111-44-4	1	ug/L
Bis(2-ethylhexyl) phthalate	117-81-7	10	ug/L
Butyl benzyl phthalate	85-68-7	10	ug/L
Caprolactam	105-60-2	10	ug/L
Carbamazepine	298-46-4	20	ug/L
Carbazole	86-74-8	10	ug/L
Chrysene	218-01-9	10	ug/L
Chrysene-d12	1719-03-5		ug/L
Coumarin	91-64-5	10	ug/L
DFTPP	5074-71-5		ug/L
Dibenz(a,h)anthracene	53-70-3	1	ug/L
Dibenzofuran	132-64-9	10	ug/L
Diethyl phthalate	84-66-2	10	ug/L
Dimethyl phthalate	131-11-3	10	ug/L

MEHOD GLE MOL'S

METADO ULE		Linite.	
4-Chloroaniline	106-47-8	2.0	ug/L
4-Chlorophenyl phenyl ether	7005-72-3	2.5	ug/L
4-Methylphenol	106-44-5	1.6	ug/L
4-Nitroaniline	100-01-6	5.8	ug/L
1-Nitrophenol	100-02-7	6.7	ug/L
Acenaphthene	83-32-9	2.7	ug/L
Acenaphthene-d10	15067-26-2		ug/L
Acenaphthylene	208-96-8	2.7	ug/L
Acetophenone	98-86-2	2.7	ug/L
Aniline	62-53-3	1.8	ug/L
Anthracene	120-12-7	2.8	ug/L
Atrazine	1912-24-9	3.0	ug/L
Benzaldehyde	100-52-7	2.0	ug/L
Benzidine	92-87-5	0.89	ug/L
Benzo[a]anthracene	56-55-3	0.27	ug/L
Benzo[a]pyrene	50-32-8	0.14	ug/L
Benzo[b]fluoranthene	205-99-2	0.26	ug/L
Benzo[g,h,i]perylene	191-24-2	2.0	ug/L
Benzo[k]fluoranthene	207-08-9	0.26	ug/L
Benzoic acid	65-85-0	50	ug/L
Benzonitrile	100-47-0	1.8	ug/L
Benzyl alcohol	100-51-6	1.5	ug/L
Bis(2-chloroethoxy)methane	111-91-1	2.6	ug/L
Bis(2-chloroethyl)ether	111-44-4	0.28	ug/L
	117-81-7	2.0	ug/L
Bis(2-ethylhexyl) phthalate	85-68-7	2.5	ug/L
Butyi benzyl phthalate	105-60-2	2.5	ug/L
Caprolactam		1.6	ug/L
Carbamazepine	298-46-4		ug/L
Carbazole	86-74-8	3.2	ug/L
Chrysene	218-01-9	3.1	_
Chrysene-d12	1719-03-5		ug/L
Coumarin	91-64-5	3.0	ug/L
DFTPP	5074-71-5		ug/L.
Dibenz(a,h)anthracene	53-70-3	0.09	ug/L
Dibenzofuran	132-64-9	2.8	ug/L
Diethyl phthalate	84-66-2	2.9	ug/L
Dimethyl phthalate	131-11-3	2.8	ug/L
Di-n-butyl phthalate	84-74-2	2.9	ug/L
Di-n-octyl phthalate	117-84-0	1.5	ug/L

		-
8270	SIM	RL

8270 SIM RL			a Dinies	
4,6-Dinitro-2-methylphenol	534-52-1	0.5	ug/L	RL
Acenaphthene	83-32-9	0.05	ug/L	RL.
Acenaphthylene	208-96-8	0.05	ug/L	RL
Anthracene	120-12-7	0.05	ug/L	RL
Benzo[a]anthracene	56-55-3	0.05	ug/L	RL
Benzo[a]pyrene	50-32-8	0.05	ug/L	RL
Benzo[b]fluoranthene	205-99-2	0.05	ug/L	RL.
Benzo[g,h,i]perylene	191-24-2	0.05	ug/L	RL.
Benzo[k]fluoranthene	207-08-9	0.05	ug/L	RL
Bis(2-chloroethyl)ether	111-44-4	0.02	ug/L	RL
Chrysene	218-01-9	0.05	ug/L	RL
Dibenz(a,h)anthracene	53-70-3	0.05	ug/L	RL
Fluoranthene	206-44-0	0.05	ug/L	RL
Fluorene	86-73-7	0.05	ug/L	RL
Hexachlorobenzene	118-74-1	0.02	ug/L	RL
indeno[1,2,3-cd]pyrene	193-39-5	0.05	ug/L	RL
Naphthalene	91-20-3	0.05	ug/L	RL
N-Nitrosodimethylamine	62-75-9	0.2	ug/L	RL
Pentachlorophenol	87-86-5	0.2	ug/L	RL.
Phenanthrene	85-01-8	0.05	ug/L	RL
Pyrene	129-00-0	0.05	ug/L	RL



#### 8270 SIM MDL

4. Analyse				
1,4-Dioxane	123-91-1	0.1517	ug/L	MDL
4,6-Dinitro-2-methylphenol	534-52-1	0.2	ug/L	MDL
Acenaphthene	83-32-9	0.0283	ug/L	MDL
Acenaphthylene	208-96-8	0.0238	ug/L	MDL
Anthracene	120-12-7	0.0283	ug/L	MDL
Benzo[a]anthracene	56-55-3	0.0350	ug/L	MDL
Benzo[a]pyrene	50-32-8	0.0475	ug/L	MDL
Benzo[b]fluoranthene	205-99-2	0.0299	ug/L	MDL
Benzo[g,h,i]perylene	191-24-2	0.0299	ug/L	MDL
Benzo[k]fluoranthene	207-08-9	0.0153	ug/L	MDL
Bis(2-chloroethyl)ether	111-44-4	0.0181	ug/L	MDL
Chrysene	218-01-9	0.0283	ug/L	MDL
Dibenz(a,h)anthracene	53-70-3	0.0363	ug/L	MDL
Fluoranthene	206-44-0	0.0283	ug/L	MDL
Fluorene	86-73-7	0.0283	ug/L	MDL
Hexachlorobenzene	118-74-1	0.0168	ug/L	MDL
Indeno[1,2,3-cd]pyrene	193-39-5	0.0283	ug/L	MDL
Naphthalene	91-20-3	0.0238	ug/L	MDL
N-Nitrosodimethylamine	62-75-9	0.0350	ug/L	MDL
Pentachlorophenol	87-86-5	0.14	ug/L	MDL
Phenanthrene	85-01-8	0.0283	ug/L	MDL
Pyrene	129-00-0	0.0336	ug/L	MDL

027	n.	0!	10
02.II	и.	RΙ	_ 2ò

8270 RL's Analytes		A Salar Electrica	e Subirs	Init in
3,5-di-tert-butyl-4-hydroxytol	128-37-0	10	ug/L	RL
3-Nitroaniline	99-09-2	10	ug/L	RL
4,6-Dinitro-2-methylphenol	534-52-1	20	ug/L	RL
4-Bromophenyl phenyl ether	101-55-3	10	ug/L	RL
4-chloro-2-methylaniline	95-69-2	10	ug/L	RL
4-Chloro-3-methylphenol	59-50-7	10	ug/L	RL
4-Chloroaniline	106-47-8	10	ug/L	RL
4-Chlorophenyl phenyl ether	7005-72-3	10	ug/L	RL
4-Methylphenol	106-44-5	10	ug/L	RL
4-Nitroaniline	100-01-6	10	ug/L	RL
4-Nitrophenol	100-02-7	20	ug/L	RL
Acenaphthene	83-32-9	10	ug/L	RL.
Acenaphthylene	208-96-8	10	ug/L	RL
Acetophenone	98-86-2	10	ug/L	RL
Aniline	62-53-3	10	ug/L	RL
Anthracene	120-12-7	10	ug/L	RL
Atrazine	1912-24-9	10	ug/L	RL
Benzaldehyde	100-52-7	10	ug/L	RL
Benzidine	92-87-5	20	ug/L	RL.
Benzo[a]anthracene	56-55-3	1	ug/L	RL
Benzo[a]pyrene	50-32-8	1	ug/L	RL
Benzo[b]fluoranthene	205-99-2	1	ug/L	RL.
Benzo[g,h,i]perylene	191-24-2	10	ug/L	RL
Benzo[k]fluoranthene	207-08-9	1	ug/L	RL
Benzoic acid	65-85-0	50	ug/L	RL
Benzyl alcohol	100-51-6	10	ug/L	RL
Bis(2-chloroethoxy)methane	111-91-1	10	ug/L	RL
Bis(2-chloroethyl)ether	111-44-4	1	ug/L	RL
Bis(2-ethylhexyl) phthalate	117-81-7	10	ug/L	RL
Bisphenol-A	80-05-7	30	ug/L	Ri.
Butyl benzyl phthalate	85-68-7	10	ug/L	RL
Caprolactam	105-60-2	10	ug/L	RL.
Carbamazepine	298-46-4	20	ug/L	RL
Carbazole	86-74-8	10	ug/L	RL
Chrysene	218-01-9	10	ug/L	RL
Coumarin	91-64-5	10	ug/L	RL
Dibenz(a,h)anthracene	53-70-3	1	ug/L	RL
Dibenzofuran	132-64-9	10	ug/L	RL



8270	
X / / II	DATE 21

8270 MDL Analyes				
4-Chloroaniline	106-47-8	2	ug/L	MDL
4-Chlorophenyi phenyi ether	7005-72-3	2.5	ug/L	MDL
4-Methylphenol	106-44-5	1.6	ug/L	MDL.
4-Nitroaniline	100-01-6	5.8	ug/L	MDL
4-Nitrophenol	100-02-7	6.7	ug/L	MDL
Acenaphthene	83-32-9	2.7	ug/L	MDL
Acenaphthylene	208-96-8	2.7	ug/L	MDL
Acetophenone	98-86-2	2.7	ug/L	MDL
Aniline	62-53-3	1.8	ug/L	MDL
Anthracene	120-12-7	2.8	ug/L	MDL
Atrazine	1912-24-9	3	ug/L	MDL
Benzaldehyde	100-52-7	2	ug/L	MDL
Benzidine	92-87-5	0.89	ug/L	MDL
·	56-55-3	0.27	ug/L	MDL
Benzo[a]anthracene	50-32-8	0.14	ug/L	MDL
Benzo[a]pyrene Benzo[b]fluoranthene	205-99-2	0.26	ug/L	MDL
	191-24-2	2	ug/L	MDL
Benzo[g,h,i]perylene	207-08-9	0.26	ug/L	MDL
Benzo[k]fluoranthene	65-85-0	50	ug/L	MDL
Benzoic acid	100-51-6	1.5	ug/L	MDL
Benzyl alcohol	111-91-1	2.6	ug/L	MDL
Bis(2-chloroethoxy)methane	111-44-4	0.28	ug/L	MDL.
Bis(2-chloroethyl)ether	117-81-7	2	ug/L	MDL.
Bis(2-ethylhexyl) phthalate	80-05-7	3.5	ug/L	MDL
Bisphenol-A	85-68-7	2.5	ug/L	MDL.
Butyl benzyl phthalate	105-60-2	2.5	ug/L	MDL
Caprolactam	298-46-4	1.6	ug/L	MDL.
Carbamazepine	86-74-8	3.2	ug/L	MDL.
Carbazole	218-01-9	3.1	ug/L	MDL
Chrysene	91-64-5	3	ug/L	MDL
Coumarin	53-70-3	0.09	ug/L	MDL
Dibenz(a,h)anthracene	132-64-9	2.8	ug/L ug/L	MDL
Dibenzofuran	84-66-2	2.9	ug/L	MDL
Diethyl phthalate		2.8	ug/L	MDL
Dimethyl phthalate	131-11-3	2.9	ug/L	MDL
Di-n-butyl phthalate	84-74-2	1.5	ug/L	MDL
Di-n-octyl phthalate	117-84-0 206-44-0		ug/L	MDL
Fluoranthene		3.2		MDL
Fluorene	86-73-7	2.8	ug/L	1716-

# Hampton Clarke

# SEMIVOLATILES (Aqueous) RL / MDL

MOI

		• •		KL	MUL	
						1 14 1 3 2
625-8270	Aqueous	Bis(2-Chloroethoxy)Methane		2	0.33051	ppb
625-8270	Aqueous	bis(2-Chloroethyl)ether		0.5	0.26639	ppb
625-8270	Aqueous	bis(2-chloroisopropyl)ether		2	0.19748	ppb
625-8270	Aqueous	bis(2-Ethylhexyl)phthalate		2	0.26216	ppb
625-8270	Aqueous	Butylbenzylphthalate		2	0.26991	ppb
625-8270	Aqueous	Caprolactam		2	0.32768	ppb
625-8270	Aqueous	Carbazole		2	0.21483	ppb
625-8270	Aqueous	Chrysene		2	0.26109	ppb
625-8270	Aqueous	Coumarin		2	0.24791	ppb
625-8270	Aqueous	Dibenzo[a,h]anthracene		2	0.27244	ppb
625-8270	Aqueous	Dibenzofuran		0.5	0.25412	ppb
625-8270	Agueous	Diethylphthalate		2	0.21768	ppb
625-8270	Aqueous	Dimethylnaphthalenes		2	0.17503	ppb
625-8270	Aqueous	Dimethylphthalate		2	0.20053	ppb
625-8270	Aqueous	Di-n-Butylphthalate		0.5	0.22526	ppb
625-8270	Aqueous	Di-n-octylphthalate		2	0.18227	ppb
625-8270	Aqueous	Diphenyl Ether		2	0.29104	ppb
625-8270	Aqueous	Fluoranthene		2	0.25699	ppb
625-8270	Aqueous	Fluorene		2	0.24931	ppb
625-8270	Aqueous	Hexachlorobenzene		2	0.54546	ppb
625-8270	Aqueous	Hexachlorobutadiene		2	0.58974	ppb
625-8270	Aqueous	Hexachlorocyclopentadiene		2	0.46359	ppb
625-8270	Aqueous	Hexachloroethane		2	0.50682	ppb
625-8270	Aqueous	Indeno[1,2,3-cd]pyrene		2	0.22197	ppb
625-8270	Aqueous	Isophorone		2	0.25581	ppb
625-8270	Aqueous	Methylnaphthalenes	•	2	0.27167	ppb
625-8270	Aqueous	Naphthalene		0.5	0.22385	ppb
625-8270	Aqueous	n-Decane	•	2	0.39496	ppb
625-8270	Aqueous	Nitrobenzene		2	0.34116	ppb
625-8270	Aqueous	N-Nitrosodimethylamine		10	0.46142	ppb
625-8270	Aqueous	N-Nitroso-Di-n-propylamine		0.5	0.33776	ppb
625-8270	Aqueous	n-Nitrosodiphenylamine		2	0.35492	bbp
625-8270	Aqueous	n-Octadecane		2	0.16261	ppb
625-8270	Aqueous	Pentachloroethane		2	0.23695	ppb
625-8270	Aqueous	Pentachlorophenol		10	3.40831	ppb
625-8270	Aqueous	Phenanthrene		2	0.22577	ppb
625-8270	Aqueous	Phenol		2	0.29395	ppb
625-8270	Aqueous	Pyrene		2	0.20070	ppb
625-8270	Aqueous	Pyridine		10	0.68442	ppb

Reporting Limit for All Methods (Soil and Aqueous) is Based on Concentration of Low Standard in Calibration Range



## Hampton Clarke

8270 SIM (Aqueous) RL / MDL

RL

MOL

			KL	LIOT	
8270 SIM	Aqueous	1,2-Diphenylhydrazine	0.05	0.01678	ppb .
8270 SIM	Aqueous	2,4,5-Trichlorophenol	0.5	0.09597	ppb
8270 SIM	Aqueous	4,6-Dinitro-2-methylphenol	0.2	0.29705	ppb
8270 SIM	Aqueous	4-Chloroaniline		0.01149	ppb
8270 SIM	Aqueous	Acenaphthene	0.02	0.01943	ppb
8270 SIM	Aqueous	Acenaphthylene	0.2	0.03668	ppb
8270 SIM	Aqueous	Anthracene	0.02	0.01867	ppb
8270 SIM	Aqueous	Benzo[a]anthracene	0.02	0.01335	ppb
8270 SIM	Aqueous	Benzo[a]Pyrene	0.02	0.019	ppb
8270 SIM	Aqueous	Benzo[b]fluoranthene	0.02	0.01403	ppb
8270 SIM	Aqueous	Benzo[g,h,i]perylene	0.02	0.01596	ppb
8270 SIM	Aqueous	Benzo[k]fluoranthene	0.02	0.00773	ppb
8270 SIM	Aqueous	bis(2-Chloroethyl)ether	0.2	0.0345	ppb
8270 SIM	Aqueous	Carbazole	0.02	0.00603	ppb
8270 SIM	Aqueous	Chrysene	0.02	0.00513	ppb
8270 SIM	Aqueous	Dibenzo[a,h]anthracene	0.02	0.00848	ppb
8270 SIM	Aqueous	Dibenzofuran			ppb
8270 SIM	Aqueous	Fluoranthene	0.02	0.01466	ppb
8270 SIM	Aqueous	Fluorene	0.02	0.01479	ppb
8270 SIM	Aqueous	Hexachlorobenzene			ppb
8270 SIM	Aqueous	Hexachlorobutadiene			ppb
8270 SIM	Aqueous	Hexachlorocyclopentadiene			ppb
8270 SIM	Aqueous	Hexachloroethane			ppb
8270 SIM	Aqueous	Indeno[1,2,3-cd]pyrene			ppb
8270 SIM	Aqueous	Naphthalene			ppb
8270 SIM	Aqueous			* - *	ppb
8270 SIM	Aqueous				ppb
8270 SIM	Aqueous	· ·			ppb
8270 SIM	Aqueous	Phenanthrene			ppb
8270 SIM	Aqueous	Pyrene	0.02	0.01433	ppb
	8270 SIM 8270 SIM	8270 SIM         Aqueous           8270 SIM         Aqueous	8270 SIM Aqueous 2,4,5-Trichlorophenol 8270 SIM Aqueous 4,6-Dinitro-2-methylphenol 8270 SIM Aqueous 4-Chloroaniline 8270 SIM Aqueous Acenaphthene 8270 SIM Aqueous Acenaphthylene 8270 SIM Aqueous Anthracene 8270 SIM Aqueous Benzo[a]anthracene 8270 SIM Aqueous Benzo[a]Pyrene 8270 SIM Aqueous Benzo[b]fluoranthene 8270 SIM Aqueous Benzo[b]fluoranthene 8270 SIM Aqueous Benzo[k]fluoranthene 8270 SIM Aqueous Benzo[k]fluoranthene 8270 SIM Aqueous Benzo[k]fluoranthene 8270 SIM Aqueous Carbazole 8270 SIM Aqueous Chrysene 8270 SIM Aqueous Dibenzo[a,h]anthracene 8270 SIM Aqueous Fluoranthene 8270 SIM Aqueous Fluoranthene 8270 SIM Aqueous Fluoranthene 8270 SIM Aqueous Hexachlorobenzene 8270 SIM Aqueous Hexachlorocyclopentadiene 8270 SIM Aqueous Indeno[1,2,3-cd]pyrene 8270 SIM Aqueous Naphthalene 8270 SIM Aqueous N-Nitrosodimethylamine 8270 SIM Aqueous N-Nitrosodimethylamine 8270 SIM Aqueous Pentachlorophenol	8270 SIM         Aqueous         1,2-Diphenylhydrazine         0.05           8270 SIM         Aqueous         2,4,5-Trichlorophenol         0.5           8270 SIM         Aqueous         4,6-Dinitro-2-methylphenol         0.2           8270 SIM         Aqueous         4-Chloroaniline         0.02           8270 SIM         Aqueous         Acenaphthene         0.02           8270 SIM         Aqueous         Acenaphthylene         0.2           8270 SIM         Aqueous         Anthracene         0.02           8270 SIM         Aqueous         Benzo[a]anthracene         0.02           8270 SIM         Aqueous         Benzo[a]Pyrene         0.02           8270 SIM         Aqueous         Benzo[a]n,i]perylene         0.02           8270 SIM         Aqueous         Benzo[g]n,i]perylene         0.02 </td <td>8270 SIM         Aqueous         1,2-Diphenylhydrazine         0.05         0.01678           8270 SIM         Aqueous         2,4,5-Trichlorophenol         0.5         0.09597           8270 SIM         Aqueous         4,6-Dinitro-2-methylphenol         0.2         0.29705           8270 SIM         Aqueous         4-Chloroaniline         0.02         0.01149           8270 SIM         Aqueous         Acenaphthene         0.02         0.01943           8270 SIM         Aqueous         Acenaphthylene         0.2         0.03668           8270 SIM         Aqueous         Anthracene         0.02         0.01867           8270 SIM         Aqueous         Benzo[a]anthracene         0.02         0.01867           8270 SIM         Aqueous         Benzo[a]Pyrene         0.02         0.019           8270 SIM         Aqueous         Benzo[a]Hyrene         0.02         0.019           8270 SIM         Aqueous         Benzo[a]Hyrene         0.02         0.019           8270 SIM         Aqueous         Benzo[a]Hyrene         0.02         0.01403           8270 SIM         Aqueous         Benzo[a]Hyrene         0.02         0.01403           8270 SIM         Aqueous         Benzo[a]Hyrene</td>	8270 SIM         Aqueous         1,2-Diphenylhydrazine         0.05         0.01678           8270 SIM         Aqueous         2,4,5-Trichlorophenol         0.5         0.09597           8270 SIM         Aqueous         4,6-Dinitro-2-methylphenol         0.2         0.29705           8270 SIM         Aqueous         4-Chloroaniline         0.02         0.01149           8270 SIM         Aqueous         Acenaphthene         0.02         0.01943           8270 SIM         Aqueous         Acenaphthylene         0.2         0.03668           8270 SIM         Aqueous         Anthracene         0.02         0.01867           8270 SIM         Aqueous         Benzo[a]anthracene         0.02         0.01867           8270 SIM         Aqueous         Benzo[a]Pyrene         0.02         0.019           8270 SIM         Aqueous         Benzo[a]Hyrene         0.02         0.019           8270 SIM         Aqueous         Benzo[a]Hyrene         0.02         0.019           8270 SIM         Aqueous         Benzo[a]Hyrene         0.02         0.01403           8270 SIM         Aqueous         Benzo[a]Hyrene         0.02         0.01403           8270 SIM         Aqueous         Benzo[a]Hyrene

Reporting Limit for All Methods (Soil and Aqueous) is based on concentration of Low Standard in Calibration Range

#### Accutest

Compound List Report

Product: AB8270NJPPTCL11+ PP/TCL Semivolatiles +25, NJ

Matrix: AQ Aqueous Feb 06, 2014 08:43

Method List: AB8270 / Method Ref: SW846 8270D

Report List: ABNJPPT ABN Special List

RL/MDL Factor: 1.0

LJ43649 LJ28154

						Rev: 7/13		
Compound CAS No.	RL	MDL	Units	MS/MSD	RPD	BS	DUP	
2-Chloroph 95-57-8	5	0.97	ua/l	30-106	•	33 42-110		10
4-Chloro-3-59-50-7	5		ug/l	40-130		28 49-120		10
2,4-Dichlor 120-83-2	2		ug/l	27-128		34 47-114		10
2,4-Dimeth 105-67-9	5		ug/l	32-145		31 48-129		10
2,4-Dinitro; 51-28-5	20		ug/l	10-162		43 26-153		10
4,6-Dinitro- 534-52-1	20		ug/l	10-147		42 44-132		10
2-Methylph 95-48-7	2		ug/l	29-111		29 39-110		10
3&4-Methy	. 2			26-114		29 36-110		10
2-Nitrophei 88-75-5	5		ug/l	27-121		35 45-110		10
4-Nitropher 100-02-7	10		ug/l	10-135		37 10-110		10
Phenol 108-95-2	2	1.3	ug/l	10-110		33 10-110		10
2,3,4,6-Tet 58-90-2	5		ug/l	30-128		33 47-119		10
2,4,5-Trich 95-95-4	5		ug/l	35-128	•	30 52-119		10
2,4,6-Trich 88-06-2	5		ug/l	31-130		34 51-118		10
Acetophen 98-86-2	2			24-142		31 39-125		10
Atrazine 1912-24-9	2		ug/l	42-160		26 58-153		10
Benzidine 92-87-5	20		ug/l	10-110		39 10-110		10
Benzaldeh 100-52-7	5		ug/l	17-130		33 32-119		10
4-Bromoph 101-55-3	2		-	51-122		26 55-120		10
Butyl benz <sub>)</sub> 85-68-7	2		_	37-145		29 42-138		10
1,1'-Bipher 92-52-4	1		ug/l	40-122		30 41-123		10
2-Chlorona 91-58-7	2		ug/l	36-108		30 38-107		10
4-Chloroan 106-47-8	5		_	24-110		32 38-110		10
Carbazole 86-74-8	1			47-128		25 55-121		10
Caprolacta 105-60-2	2		_	10-110		44 10-113		10
bis(2-Chlor 111-91-1	2			43-120		31 48-117		10
bis(2-Chlor 111-44-4	2			33-115		34 40-108		10
bis(2-Chlor 108-60-1	2			28-114		33 37-107		10
4-Chloroph 7005-72-3	2			46-118		29 48-117		10
1,2-Dichlor 95-50-1	1	0.29	_	22-110		34 23-110		10
1,2-Diphen 122-66-7	1	0.29		43-133		27 49-130		10
1,3-Dichlor 541-73-1	1	0.38		19-110		34 20-110		10
1,4-Dichlor 106-46-7	1	0.36	•	20-110		34 21-110		10
2,4-Dinitrot 121-14-2	1	0.43	-	49-129		27 53-129		10
2,6-Dinitrot 606-20-2	1	0.46	-	52-129		26 57-126		10
3,3'-Dichlor91-94-1	2		_	14-119		33 41-109		10
Dibenzofur 132-64-9	5		_	44-114		28 46-112		10
Di-n-butyl r 84-74-2	2		_	47-131 47-149		27 54-127		10
Di-n-octyl r 117-84-0	2		-	47-148		29 55-142		10
Diethyl phtl 84-66-2	2		-	32-131		29 35-130		10
Dimethyl pl 131-11-3	2	0.28	ug/I	26-132		33 26-132		10

#### Accutest

Compound List Report
Product: AB625NJTCL11+ TCL Semivolatiles +20, NJ

Matrix: AQ Aqueous

Aug 29, 2014 11:44 an

Method List: AB625 A(Method Ref: EPA 625 Report List: ABNJTCL ABN TCL List without PAH LJ39692 LJ32414

RL/MDL Factor: 1.0

	Compound CAS No.	RL	MDL	Units	Control Limits (%) MS/MSD RPD	Rev: 1/14 BS	DUP	
	2-Chloroph 95-57-8	5	0.95	ug/l	20-122	36 36-112		0
	4-Chloro-3-59-50-7	ε	1.2	ug/l	35-133	26 43-122		0
	2,4-Dichlor 120-83-2	5	i 1.6	ug/l	25-132	35 40-126		0
	2,4-Dimeth 105-67-9	5	1.6	ug/i	36-138	27 44-131	1	0
	2,4-Dinitro; 51-28-5	20	0.89	ug/l	10-152	44 17-161		0
	4.6-Dinitro- 534-52-1	10	0.72	ug/l	10-141	43 34-142		0
	2-Methylph 95-48-7	2	1.4	ug/l	24-123	26 35-108		0
	3&4-Methy	5	i 1.3	ug/l	29-117	27 31-107		25
	2-Nitropher 88-75-5	2		ug/l	13-133	30 35-128		0
	4-Nitrophei 100-02-7	10	0.84	ug/l	10-132	30 10-117		0
	Phenol 108-95-2	2		ug/l	10-110	34 10-110		0
	2,3,4,6-Tet 58-90-2	5	5 1.1	ug/l	23-139	34 46-127		0
	2,4,5-Trich 95-95-4	2	2 1.9	ug/l	30-134	31 51-124		0
	2,4,6-Trich 88-06-2	2	2 1.3	ug/l	24-139	32 49-127		0
	Acetophen 98-86-2	{		ug/l	34-139	29 41-131		0
	Atrazine 1912-24-			ug/l	61-160	27 64-156		0
	Benzaldeh 100-52-7			' ug/l	15-163	33 26-152		0
	1,1'-Bipher 92-52-4	5		ug/l	43-130	27 46-126		0
	Caprolacta 105-60-2			ug/l	10-110	32 10-108		0
	4-Bromoph 101-55-3			ug/i	57-121	26 55-124		0
	Butyl benzy 85-68-7	2		ug/l	53-136	27 45-142		0
	2-Chlorona 91-58-7			ug/l	42-111	29 38-115		0
	4-Chloroan 106-47-8	2	0.4	· ug/l	11-100	32 19-102		0
	Carbazole 86-74-8	2		ug/l	60-118	25 55-124		2
	bis(2-Chlor 111-91-1	2	0.65	ug/l	42-124	29 42-126		0
(	bis(2-Chlor 111-44-4	2	2 0.53	ug/l	38-119	27 38-120		0
	bis(2-Chlor 108-60-1		0.74	ug/l	34-117	29 33-121		0
	4-Chloroph 7005-72-	3 2		ug/l	53-117	26 50-119		0
	2,4-Dinitrot 121-14-2	2		ug/l	52-129	24 56-133		0
	2,6-Dinitrot 606-20-2	2	0.56	ug/l	57-129	27 56-135		0
	3,3'-Dichlor 91-94-1	2	1.2	ug/l	10-135	31 24-126		0
	Dibenzofur 132-64-9			ug/l	49-113	26 47-115		18 17
	Di-n-butyl r 84-74-2	2		ug/l	55-130	25 50-135		
	Di-n-octyl r 117-84-0	2		ug/l	58-140	25 50-146		0
	Diethyl phtl 84-66-2	2		ug/l	47-126	24 43-129		10 10
	Dimethyl pl 131-11-3			ug/l	45-126	27 43-129		
	bis(2-Ethyll 117-81-7			i ug/l	60-135	22 51-143 31 15-114		22 10
	Hexachlor 87-68-3			ug/l	15-111			10
	Hexachlorc 77-47-4	10		ug/l	10-115	39 10-127 30 16 108		10
	Hexachlorc 67-72-1	į		ug/l	10-110	30 16-108 29 40-124		10
	Isophoron∈ 78-59-1	. 2	2 0.59	ug/l	38-125	29 40-124		Ü

#### Method Blank Summary

#### Laboratory DRAFT - Not for Release

 $\begin{array}{c}QF\\1^{a}\end{array}$ 

\*

Sample OP77597A-MB1	File ID 3M49296.D	DF 1	Analyzed 09/03/14	By SP	Prep Date 08/28/14	Prep Batch OP77597A	Analytical Batch E3M2238

The QC reported here applies to the following samples:

Method: SW846 8270D BY SIM

JB74466-3, JB74724-13, JB74764-1, JB74764-3, JB74764-4, JB74764-5, JB74764-6, JB74777-1, JB74777-2, JB74777-3, JB74777-5, JB74777-7, JB74834-1, JB74834-2, JB74834-3, JB74834-4, JB74834-5

Cresol, Total ND 1.0 0.011 ug/l 108-95-2 Phenol ND 0.30 0.10 ug/l 108-95-2 Phenol ND 0.50 0.14 ug/l 83-32-9 Acenaphthene ND 0.10 0.020 ug/l 208-96-8 Acenaphthylene ND 0.10 0.024 ug/l 120-12-7 Anthracene ND 0.10 0.024 ug/l 120-12-7 Anthracene ND 0.10 0.024 ug/l 120-12-8 Benzo(a) anthracene ND 0.10 0.012 ug/l 205-99-2 Benzo(b) fluoranthene ND 0.10 0.012 ug/l 191-24-2 Benzo(g, h, i) perylene ND 0.10 0.016 ug/l 207-08-9 Benzo(k) fluoranthene ND 0.10 0.015 ug/l 218-01-9 Chrysene ND 0.10 0.012 ug/l 111-44-4 bis(2-Chloroethyl) ether ND 0.10 0.012 ug/l 206-44-0 Fluoranthene ND 0.10 0.010 0.0026 ug/l 53-70-3 Dibenzo(a, h) anthracene ND 0.10 0.017 ug/l 206-44-0 Fluoranthene ND 0.10 0.017 ug/l 86-73-7 Fluorene ND 0.10 0.017 ug/l 118-74-1 Hexachlorobenzene ND 0.10 0.017 ug/l 87-68-3 Hexachlorobenzene ND 0.10 0.017 ug/l 193-39-5 Indeno(1, 2, 3-cd) pyrene ND 0.10 0.0028 ug/l 193-39-5 Indeno(1, 2, 3-cd) pyrene ND 0.10 0.023 ug/l 91-57-6 2-Methylnaphthalene ND 0.10 0.021 ug/l 191-20-3 Naphthalene ND 0.10 0.021 ug/l 191-20-3 Naphthalene ND 0.10 0.021 ug/l 191-20-3 Naphthalene ND 0.10 0.021 ug/l 191-20-0-0 Pyrene ND 0.10 0.015 ug/l 191-20-0-0 Pyrene ND 0.10	CAS No.	Compound	Result	RL	MDL	Units	Q
87-86-5         Pentachlorophenol         ND         0.30         0.10         ug/l           108-95-2         Phenol         ND         0.50         0.14         ug/l           83-32-9         Acenaphthene         ND         0.10         0.020         ug/l           208-96-8         Acenaphthylene         ND         0.10         0.024         ug/l           120-12-7         Anthracene         ND         0.10         0.020         ug/l           56-55-3         Benzo(a)anthracene         ND         0.10         0.012         ug/l           50-32-8         Benzo(b)fluoranthene         ND         0.10         0.012         ug/l           205-99-2         Benzo(b)fluoranthene         ND         0.10         0.010         ug/l           191-24-2         Benzo(g,h,i)perylene         ND         0.10         0.016         ug/l           207-08-9         Benzo(k)fluoranthene         ND         0.10         0.015         ug/l           218-01-9         Chrysene         ND         0.10         0.012         ug/l           211-44-4         bis(2-Chloroethyl)ether         ND         0.10         0.0026         ug/l           53-70-3         Dibenzo(a,h)ant		Cresol, Total	ND cons	1.0	0.011	ug/l	
108-95-2         Phenol         ND         0.50         0.14         ug/l           83-32-9         Acenaphthene         ND         0.10         0.020         ug/l           208-96-8         Acenaphthylene         ND         0.10         0.024         ug/l           120-12-7         Anthracene         ND         0.10         0.020         ug/l           56-55-3         Benzo(a)anthracene         ND         0.10         0.012         ug/l           50-32-8         Benzo(a)pyrene         ND         0.10         0.012         ug/l           205-99-2         Benzo(b)fluoranthene         ND         0.10         0.010         ug/l           191-24-2         Benzo(g,h,i)perylene         ND         0.10         0.016         ug/l           207-08-9         Benzo(k)fluoranthene         ND         0.10         0.015         ug/l           218-01-9         Chrysene         ND         0.10         0.012         ug/l           211-4-4         bis(2-Chloroethyl)ether         ND         0.10         0.0026         ug/l           23-70-3         Dibenzo(a,h)anthracene         ND         0.10         0.017         ug/l           26-44-0         Fluoranthene <td>87-86-5</td> <td>· · · · · · · · · · · · · · · · · · ·</td> <td>- 1 (1 (1 (1 (1 (1 (1 (1 (1 (1 (1 (1 (1 (</td> <td>0.30</td> <td>0.10</td> <td></td> <td></td>	87-86-5	· · · · · · · · · · · · · · · · · · ·	- 1 (1 (1 (1 (1 (1 (1 (1 (1 (1 (1 (1 (1 (	0.30	0.10		
83-32-9         Acenaphthene         ND         0.10         0.020         ug/1           208-96-8         Acenaphthylene         ND         0.10         0.024         ug/1           120-12-7         Anthracene         ND         0.10         0.020         ug/1           56-55-3         Benzo(a)anthracene         ND         0.10         0.012         ug/1           50-32-8         Benzo(a)pyrene         ND         0.10         0.012         ug/1           205-99-2         Benzo(b)fluoranthene         ND         0.10         0.010         ug/1           191-24-2         Benzo(k)fluoranthene         ND         0.10         0.016         ug/1           207-08-9         Benzo(k)fluoranthene         ND         0.10         0.015         ug/1           218-01-9         Chrysene         ND         0.10         0.012         ug/1           218-01-9         Chrysene         ND         0.010         0.012         ug/1           211-44-4         bis(2-Chloroethyl)ether         ND         0.010         0.0026         ug/1           25-70-3         Dibenzo(a,h)anthracene         ND         0.10         0.017         ug/1           26-44-0         Fluoranthe		-	ND	0.50	0.14	ug/l	
208-96-8         Acenaphthylene         ND         0.10         0.024         ug/l           120-12-7         Anthracene         ND         0.10         0.020         ug/l           56-55-3         Benzo(a)anthracene         ND         0.10         0.012         ug/l           50-32-8         Benzo(a)pyrene         ND         0.10         0.012         ug/l           205-99-2         Benzo(b)fluoranthene         ND         0.10         0.010         ug/l           191-24-2         Benzo(g,h,i)perylene         ND         0.10         0.016         ug/l           207-08-9         Benzo(k)fluoranthene         ND         0.10         0.015         ug/l           207-08-9         Benzo(k)fluoranthene         ND         0.10         0.015         ug/l           207-08-9         Benzo(k)fluoranthene         ND         0.10         0.012         ug/l           207-08-9         Benzo(k)fluoranthene         ND         0.10         0.012         ug/l           218-01-9         Chrysene         ND         0.10         0.002         ug/l           211-14-4         bis(2-Chloroethyl)ether         ND         0.10         0.017         ug/l           206-44-0		Acenaphthene	ND	1111	0.020		
120-12-7			ND	0.10	0.024	ug/l	
56-55-3         Benzo (a) anthracene         ND         0.10         0.012         ug/l           50-32-8         Benzo (a) pyrene         ND         0.10         0.012         ug/l           205-99-2         Benzo (b) fluoranthene         ND         0.10         0.010         ug/l           191-24-2         Benzo (g,h,i) perylene         ND         0.10         0.016         ug/l           207-08-9         Benzo (k) fluoranthene         ND         0.10         0.015         ug/l           218-01-9         Chrysene         ND         0.10         0.012         ug/l           218-01-9         Chrysene         ND         0.10         0.012         ug/l           218-01-9         Chrysene         ND         0.10         0.012         ug/l           218-01-9         Chrysene         ND         0.10         0.0026         ug/l           218-01-9         Chrysene         ND         0.010         0.0026         ug/l           219-04-0         Dibenzo (a,h) anthracene         ND         0.10         0.017         ug/l           206-44-0         Fluoranthene         ND         0.10         0.017         ug/l           86-73-7         Fluorene			ND ***	0.10	0.020	ug/l	
50-32-8         Benzo(a) pyrene         ND         0.10         0.012         ug/l           205-99-2         Benzo(b) fluoranthene         ND         0.10         0.010         ug/l           191-24-2         Benzo(g,h,i) perylene         ND         0.10         0.016         ug/l           207-08-9         Benzo(k) fluoranthene         ND         0.10         0.015         ug/l           218-01-9         Chrysene         ND         0.10         0.012         ug/l           111-44-4         bis(2-Chloroethyl) ether         ND         0.010         0.0026         ug/l           53-70-3         Dibenzo(a,h) anthracene         ND         0.10         0.017         ug/l           206-44-0         Fluoranthene         ND         0.10         0.017         ug/l           86-73-7         Fluorene         ND         0.10         0.017         ug/l           118-74-1         Hexachlorobenzene         ND         0.020         0.017         ug/l           87-68-3         Hexachlorocyclopentadiene         ND         0.10         0.017         ug/l           193-39-5         Indeno(1,2,3-cd) pyrene         ND         0.10         0.014         ug/l           90-12-		Benzo(a)anthracene	ND	0.10	0.012	ug/I	
205-99-2         Benzo (b) fluoranthene         ND         0.10         0.010         ug/l           191-24-2         Benzo (g,h,i) perylene         ND         0.10         0.016         ug/l           207-08-9         Benzo (k) fluoranthene         ND         0.10         0.015         ug/l           218-01-9         Chrysene         ND         0.10         0.012         ug/l           111-44-4         bis (2-Chloroethyl) ether         ND         0.010         0.0026         ug/l           53-70-3         Dibenzo (a,h) anthracene         ND         0.10         0.017         ug/l           206-44-0         Fluoranthene         ND         0.10         0.017         ug/l           86-73-7         Fluorene         ND         0.10         0.017         ug/l           118-74-1         Hexachlorobenzene         ND         0.020         0.017         ug/l           87-68-3         Hexachlorobutadiene         ND         0.10         0.017         ug/l           193-39-5         Indeno (1,2,3-cd) pyrene         ND         0.10         0.044         ug/l           90-12-0         1-Methylnaphthalene         ND         0.10         0.023         ug/l           91			ND .	0.10	0.012	ug/l	
191-24-2   Benzo(g,h,i)perylene   ND   0.10   0.016   ug/l			ND	0.10	0.010	ug/l	
207-08-9         Benzo(k)fluoranthene         ND         0.10         0.015         ug/l           218-01-9         Chrysene         ND         0.10         0.012         ug/l           111-44-4         bis(2-Chloroethyl)ether         ND         0.010         0.0026         ug/l           53-70-3         Dibenzo(a,h)anthracene         ND         0.10         0.017         ug/l           206-44-0         Fluoranthene         ND         0.10         0.013         ug/l           86-73-7         Fluorene         ND         0.10         0.017         ug/l           118-74-1         Hexachlorobenzene         ND         0.020         0.017         ug/l           87-68-3         Hexachlorobutadiene         ND         0.10         0.017         ug/l           77-47-4         Hexachlorocyclopentadiene         ND         0.20         0.0088         ug/l           193-39-5         Indeno(1,2,3-cd)pyrene         ND         0.10         0.014         ug/l           90-12-0         1-Methylnaphthalene         ND         0.10         0.023         ug/l           91-20-3         Naphthalene         ND         0.10         0.036         ug/l           85-01-8			ND 🗀	0.10	0.016	ug/I	
218-01-9         Chrysene         ND         0.10         0.012         ug/l           111-44-4         bis(2-Chloroethyl)ether         ND         0.010         0.0026         ug/l           53-70-3         Dibenzo(a,h)anthracene         ND         0.10         0.017         ug/l           206-44-0         Fluoranthene         ND         0.10         0.013         ug/l           86-73-7         Fluorene         ND         0.10         0.017         ug/l           118-74-1         Hexachlorobenzene         ND         0.020         0.017         ug/l           87-68-3         Hexachlorobutadiene         ND         0.10         0.017         ug/l           77-47-4         Hexachlorocyclopentadiene         ND         0.20         0.0088         ug/l           193-39-5         Indeno(1,2,3-cd)pyrene         ND         0.10         0.014         ug/l           90-12-0         1-Methylnaphthalene         ND         0.10         0.023         ug/l           91-57-6         2-Methylnaphthalene         ND         0.10         0.023         ug/l           91-20-3         Naphthalene         ND         0.10         0.021         ug/l           85-01-8			ND	0.10	0.015	ug/l	
111-44-4         bis(2-Chloroethyl)ether         ND         0.010         0.0026         ug/l           53-70-3         Dibenzo(a,h)anthracene         ND         0.10         0.017         ug/l           206-44-0         Fluoranthene         ND         0.10         0.013         ug/l           86-73-7         Fluorene         ND         0.10         0.017         ug/l           118-74-1         Hexachlorobenzene         ND         0.020         0.017         ug/l           87-68-3         Hexachlorobutadiene         ND         0.10         0.017         ug/l           77-47-4         Hexachlorocyclopentadiene         ND         0.20         0.0088         ug/l           193-39-5         Indeno(1,2,3-cd)pyrene         ND         0.10         0.014         ug/l           90-12-0         1-Methylnaphthalene         ND         0.10         0.023         ug/l           91-57-6         2-Methylnaphthalene         ND         0.10         0.023         ug/l           91-20-3         Naphthalene         ND         0.10         0.036         ug/l           85-01-8         Phenanthrene         ND         0.10         0.021         ug/l           129-00-0		* *	ND	0.10	0.012	ug/l	
53-70-3         Dibenzo (a,h) anthracene         ND         0.10         0.017         ug/l           206-44-0         Fluoranthene         ND         0.10         0.013         ug/l           86-73-7         Fluorene         ND         0.10         0.017         ug/l           118-74-1         Hexachlorobenzene         ND         0.020         0.017         ug/l           87-68-3         Hexachlorobutadiene         ND         0.10         0.017         ug/l           77-47-4         Hexachlorocyclopentadiene         ND         0.20         0.0088         ug/l           193-39-5         Indeno(1,2,3-cd)pyrene         ND         0.10         0.014         ug/l           90-12-0         1-Methylnaphthalene         ND         0.10         0.023         ug/l           91-57-6         2-Methylnaphthalene         ND         0.10         0.023         ug/l           91-20-3         Naphthalene         ND         0.10         0.036         ug/l           85-01-8         Phenanthrene         ND         0.10         0.021         ug/l           129-00-0         Pyrene         ND         0.10         0.015         ug/l			ND	0.010	0.0026	ug/I	
86-73-7         Fluorene         ND         0.10         0.017         ug/l           118-74-1         Hexachlorobenzene         ND         0.020         0.017         ug/l           87-68-3         Hexachlorobutadiene         ND         0.10         0.017         ug/l           77-47-4         Hexachlorocyclopentadiene         ND         0.20         0.0088         ug/l           193-39-5         Indeno(1,2,3-cd)pyrene         ND         0.10         0.014         ug/l           90-12-0         1-Methylnaphthalene         ND         0.10         0.023         ug/l           91-57-6         2-Methylnaphthalene         ND         0.10         0.023         ug/l           91-20-3         Naphthalene         ND         0.10         0.036         ug/l           85-01-8         Phenanthrene         ND         0.10         0.021         ug/l           129-00-0         Pyrene         ND         0.10         0.015         ug/l	53-70-3		ND	0.10	0.017	ug/I	
118-74-1         Hexachlorobenzene         ND         0.020         0.017         ug/l           87-68-3         Hexachlorobutadiene         ND         0.10         0.017         ug/l           77-47-4         Hexachlorocyclopentadiene         ND         0.20         0.0088         ug/l           193-39-5         Indeno(1,2,3-cd)pyrene         ND         0.10         0.014         ug/l           90-12-0         1-Methylnaphthalene         ND         0.10         0.023         ug/l           91-57-6         2-Methylnaphthalene         ND         0.10         0.023         ug/l           91-20-3         Naphthalene         ND         0.10         0.036         ug/l           85-01-8         Phenanthrene         ND         0.10         0.021         ug/l           129-00-0         Pyrene         ND         0.10         0.015         ug/l	206-44-0	Fluoranthene	ND	0.10	0.013	ug/I	
87-68-3         Hexachlorobutadiene         ND         0.10         0.017         ug/I           77-47-4         Hexachlorocyclopentadiene         ND         0.20         0.0088         ug/I           193-39-5         Indeno(1,2,3-cd)pyrene         ND         0.10         0.014         ug/I           90-12-0         1-Methylnaphthalene         ND         0.10         ug/I           91-57-6         2-Methylnaphthalene         ND         0.10         0.023         ug/I           91-20-3         Naphthalene         ND         0.10         0.036         ug/I           85-01-8         Phenanthrene         ND         0.10         0.021         ug/I           129-00-0         Pyrene         ND         0.10         0.015         ug/I	86-73-7	Fluorene	ND -	0.10	0.017	ug/I	
77-47-4         Hexachlorocyclopentadiene         ND         0.20         0.0088         ug/l           193-39-5         Indeno(1,2,3-cd)pyrene         ND         0.10         0.014         ug/l           90-12-0         1-Methylnaphthalene         ND         0.10         ug/l           91-57-6         2-Methylnaphthalene         ND         0.10         0.023         ug/l           91-20-3         Naphthalene         ND         0.10         0.036         ug/l           85-01-8         Phenanthrene         ND         0.10         0.021         ug/l           129-00-0         Pyrene         ND         0.10         0.015         ug/l	118-74-1	Hexachlorobenzene	ND	0.020	0.017	ug/l	
193-39-5         Indeno(1,2,3-cd)pyrene         ND         0.10         0.014         ug/l           90-12-0         1-Methylnaphthalene         ND         0.10         ug/l           91-57-6         2-Methylnaphthalene         ND         0.10         0.023         ug/l           91-20-3         Naphthalene         ND         0.10         0.036         ug/l           85-01-8         Phenanthrene         ND         0.10         0.021         ug/l           129-00-0         Pyrene         ND         0.10         0.015         ug/l		Hexachlorobutadiene	ND	0.10	0.017	ug/I	
193-39-5         Indeno(1,2,3-cd)pyrene         ND         0.10         0.014         ug/l           90-12-0         1-Methylnaphthalene         ND         0.10         ug/l           91-57-6         2-Methylnaphthalene         ND         0.10         0.023         ug/l           91-20-3         Naphthalene         ND         0.10         0.036         ug/l           85-01-8         Phenanthrene         ND         0.10         0.021         ug/l           129-00-0         Pyrene         ND         0.10         0.015         ug/l	77-47-4	Hexachlorocyclopentadiene	ND	0.20	0.0088	ug/l	
91-57-6         2-Methylnaphthalene         ND         0.10         0.023         ug/l           91-20-3         Naphthalene         ND         0.10         0.036         ug/l           85-01-8         Phenanthrene         ND         0.10         0.021         ug/l           129-00-0         Pyrene         ND         0.10         0.015         ug/l			ND	0.10	0.014	ug/l	
91-57-6       2-Methylnaphthalene       ND       0.10       0.023       ug/l         91-20-3       Naphthalene       ND       0.10       0.036       ug/l         85-01-8       Phenanthrene       ND       0.10       0.021       ug/l         129-00-0       Pyrene       ND       0.10       0.015       ug/l	90-12-0	1-Methylnaphthalene	ND	0.10		ug/l	
91-20-3         Naphthalene         ND         0.10         0.036         ug/l           85-01-8         Phenanthrene         ND         0.10         0.021         ug/l           129-00-0         Pyrene         ND         0.10         0.015         ug/l	91-57-6		ND	0.10	0.023	ug/I	
129-00-0 Pyrene ND 0.10 0.015 ug/l	91-20-3		ND	0.10	0.036	ug/l	
The state of the s	85-01-8	Phenanthrene	ND	0.10	0.021	ug/I	
Toluidines (Total) ND 0.20 ug/l	129-00-0	Pyrene	ND	0.10	0.015	ug/l	
Tournaires (Total)		Toluidines (Total)	ND	0.20		ug/l	
CAS No. Surrogate Recoveries Limits	CAS No.	Surrogate Recoveries		Limits	3		
367-12-4 2-Fluorophenol 28% 10-110%	367-12-4	2-Fluorophenol	MARKET TO A THE PARTY	SCAL			

CAS No.	Surrogate Recoveries		Limits
367-12-4 4165-62-2 118-79-6 4165-60-0 321-60-8 1718-51-0	2-Fluorophenol Phenol-d5 2,4,6-Tribromophenol Nitrobenzene-d5 2-Fluorobiphenyl Terphenyl-d14	28% 20% 72% 58% 63% 87%	10-110% 10-110% 10-157% 23-131% 24-120% 10-125%

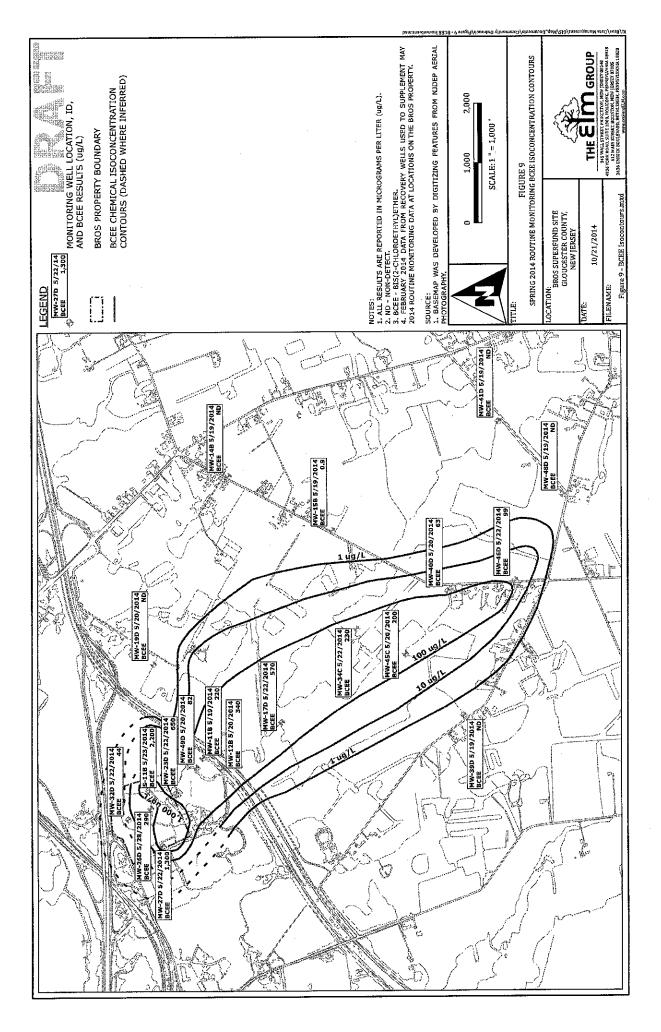
#### Exhibit 7

Figure 9, BCEE Isocontours

#### **Exhibit 8**

**BCEE PRGs, Bridgeport Rental** 

# Figure 9 BCEE Isocontours



### BCEE PRGS Bridgeport Rental

July 2006

Bridgeport Rental and Oil Services (BROS)

Logan Township

Gloucester County, New Jersey

Proposed Plan

U.S. Environmental Protection Agency Superfund Program – Region 2



#### INTRODUCTION

This Proposed Plan identifies the Preferred Alternative for cleanup of groundwater, soil and sediment contamination at the Bridgeport Rental and Oil Services (BROS) Superfund site, and provides the rationale for this preference. The proposed remedial actions are categorized as Groundwater Work and Wetlands Work. The actions contemplated under this Proposed Plan are designated as Operable Unit (OU) 2 work.

The Groundwater Work has been further divided into two major sub-categories to address media or area-specific concerns. The sub-categories are Soils, Light Non-Aqueous Phase Liquids (LNAPLs), Shallow Groundwater, and Deep Groundwater.

The Groundwater Work includes addressing onproperty and off-property shallow and deep
groundwater contamination associated with drinking
water aquifers (water bearing strata) of the PotomacRaritan-Magothy (PRM) system. It also includes
components designed to address residual subsurface
soil and LNAPL contamination which impact the
groundwater system and are themselves media which
represent potential exposure points in the human
health risk model. For example, the contaminated
soils and LNAPL are a potential exposure point/source
to construction workers through a direct contact
pathway and future building occupants through the
vapor intrusion pathway.

The PRM aquifers were impacted by releases from on-site sources including the former 13-acre waste oil lagoon, tank farms, contaminated debris, residually contaminated soil and LNAPLs. The former waste oil lagoon and tank farms were cleaned up as part of the OU-1 remedy for the site. The OU-1 actions were completed in 1997.

The Wetlands Work includes addressing the stream corridor and wetland areas both on-property and offproperty. The wetland areas were impacted by the releases of contaminants from the former waste oil lagoon during a series of overflow/spill events in the late 1960's and 1970's.

#### Preferred Alternative

The Preferred Alternative is a set of alternatives which combines technologies, within an adaptive management approach, to address both impacted media as well as post-lagoon remediation residual contamination. The preferred alternative, which is described in greater detail in the Preferred Alternative section of this Proposed Plan, includes the following Groundwater and Wetlands Work activities:

- Soil "Hot Spot" management through cover and drainage improvements, improved water budget management (using phytoremediation techniques), enhanced biodegradation, and institutional controls (ICs);
- <u>LNAPLs</u> management through cover and drainage improvements, limited property excavation, improved water budget management, enhanced LNAPL recovery via bioslurping with steam injection, where warranted, and ICs;
- <u>Shallow Groundwater</u> management through residual source remediation controls, improved water budget management (using phytoremediation techniques), groundwater extraction concurrent with the LNAPL removal system, monitored natural attenuation and ICs;
  - Deep Groundwater management via in-situe chemical oxidation treatment and enhanced biodegradation in conjunction with source area pumping and treatment (with a contingency for hydraulic containment see below for additional information); and,
    - Wetlands sediment excavation, ex-situ treatment, off-site disposal (via landfilling),

LNAPL and metals are present in the DMZ. In this zone, calculated risks to red fox and Eastern screech owl were higher than the reference areas. However, in LTCS II/III (the areas of highest concern), no ecological risk exceeding an HQ of 1.0 was noted for the red fox (representative of the upper trophic level predatory mammal) or the Eastern screech owl (representative of the upper trophic level predatory avian population). Although not an ideal habitat for mink, HQ values were also calculated for this sensitive receptor. HQ values for mink were well below one.

Similar results were obtained for areas of CS, where great blue heron (receptor representing piscivorous bird population), red fox and Eastern screech owl had HQ values less that one. Risks to white perch (representing water column biota) and mommichugs (benthic feeding forage fish species) were not considered significant.

Swindell Pond sampling indicated no observed results above aquatic benchmarks. It was concluded that Swindell Pond does not contain BROS-related COPECs at concentrations of potential ecological concern. Potential adverse effects to benthic organisms, however, are probable within a limited area surrounding the seep area of Gaventa Pond.

As previously noted, while habitat for threatened and endangered species potentially occurs in Gloucester County, no rare plants or animals have been observed on the site.

#### REMEDIAL ACTION OBJECTIVES

It is EPA's current judgment that the Preferred Alternative identified in this Proposed Plan is necessary to protect public health and the environment from actual and threatened releases of hazardous substances, pollutants or contaminants from the site which may present an imminent and substantial endangerment to public health or welfare.

Remedial action objectives (RAOs) were developed in accordance with the National Contingency Plan and current EPA guidance. RAOs were developed for each media of concern considering on-property use in the commercial/industrial category, and off-property use under a residential/agricultural/recreational setting.

The RAOs for shallow groundwater on- and offproperty include:

· Protection of the public against ingestion and

direct contact with VOCs, SVOCs and metals in the groundwater above preliminary remediation goals (PRGs — including the cleanup of shallow groundwater or reductions in movement to ensure that contaminants do not occur at potential exposure points);

- Protection against vapor intrusion from VOCs, SVOCs and PCBs in groundwater above PRGs; and,
- Protection of the public and utility workers against direct contact with VOCs, SVOCs LNAPL, PCBs and metals in groundwater above PRGs.

From a numerical standpoint, the PRGs identified, to the extent practicable at this stage for shallow groundwater, include the lower of federal or state maximum contaminant levels (MCLs) or the state groundwater quality criteria which are considered ARARs for the site. However, there is the potential for a future waiver of these chemical-specific requirements based on the ability of the proposed remedial alternatives to reduce localized areas to below threshold levels. Any such waiver would be subject to the NCP process including public participation. The following table provides numerical values for key site COCs. Values for all of the groundwater COCs are found in the RI. It is believed that the RAOs for off-property shallow groundwater are achievable in a reasonable period of time.

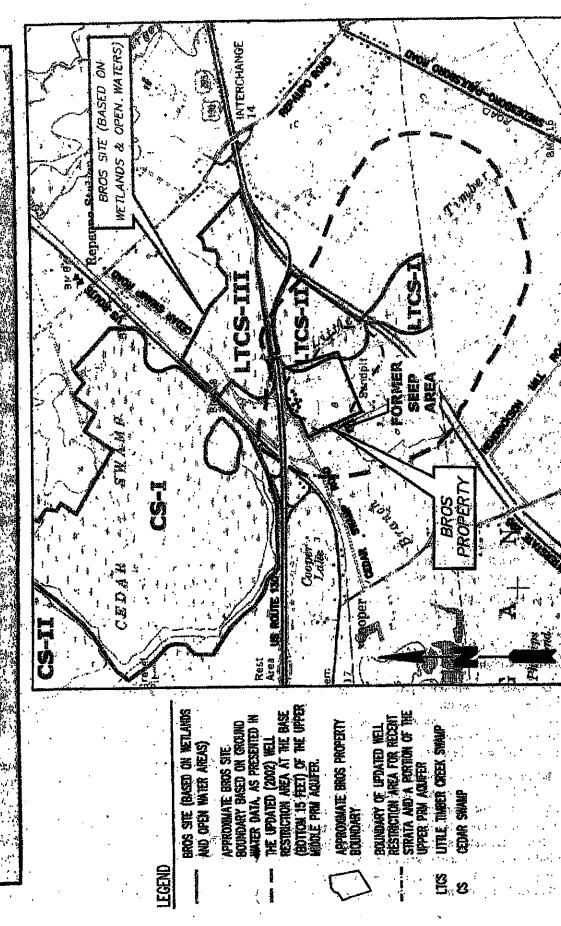
#### **Groundwater PRGs**

Analyte/ Contaminant Group	njdep Gwqc	Federal MCX
Serie Granic Companies as L Benzene		5
TCR Vinyl Chloride	I I	.2.
emi Politi Grenie comunidate DES		
BCEE  **********************************		
Total PCBs	0.5	0.5

TABLE 2
BROS Preferred Remedy Components by Media

Media	Alternative No.	Remedy Components	Estimated Costs
Wetlands	DMZ-5	Excavation/ Ex-Situ Treatment/ Off-Site Disposal/ Application of Sorptive Agent (prior to Capping or incorporated into Backfilly/ Wetlands Restoration Physical Amenities: Includes the excavation of approximately 17,500 cubic yards of contaminated sediment and application of sorptive material over 10.6 acres.	\$11.9M
	IZ-2	Natural Remediation (Monitored)/ Institutional Controls	\$0.6M
	SHS-4	Enhanced Biodegradation Component Only Physical Amenities: Includes the installation of at least 230 chemical injection points.	\$2.3M
Soil, LNAPL, and Shallow Groundwater	LNAPL-5	Institutional Controls/ Cover and Drainage Improvements/ Limited Off-Property Excavation (Gaventa Pond Seep and Green Acres Area)/ Enhanced LNAPL Recovery via Bioslurping and Thermal/Steam Injection (where warranted following Bioslurping)/ Containment-Water Bidget Management via Phytoremediation/Alternate Final Cover Physical Amenities: Includes the installation of approximately 72 hioslurping extraction points. Water budget management may include planting up to 1,000 trees per zero in LNAPL areas.	\$16.5M
	SGW-2	Institutional Controls/ Source Remediation/Control/ Monitored Natural Attenuation	\$1.9
Deep Groundwater	DGW-6	Phased Combination: Source Area (Principal Threat Zone) Pumping and Treatment (Mass Reduction)/ Followed by In-Situ Chemical Oxidation Treatment in Significant Rebound Areas/ Followed by Enhanced Biodegradation—Lower Threat Zone Pumping and Treatment (Mass Reduction)/ Followed by Enhanced Biodegradation in Significant Rebound Areas  Downgradient Area Enhanced Biodegradation  Physical Amenities: includes the installation of over 50 extraction wells in the PTZ/LTZ and 300 Chemical Oxidant injection wells. Will include the inocalation of groundwater with an estimated 600,000 pounds of oxidant and the extraction of over 100 million gallons of contaminated groundwater over the first two years of operation. Long-term operations could realize the extraction of over 500 million gallons of groundwater.	\$57.7M
Deep Groundwater Contingency	DGW-4	Source Area Containment Pumping and Treatment/ Downgradient Area Enhanced Aerobic Biodegradation  Physical Amenities: Includes the groundwater treatment plant constructed for DGW-6 with additional wells to capture the plume.	\$48.5M. (1)  (1) If implemented as a contingency, cost would be reduced by capital expendity for treatment plant construction under DGW-6

Estimated total cost: \$90.9M (Wetlands \$12.5M; Soils, LNAPLs, Shallow GW \$20.7; Deep GW \$57.7M)



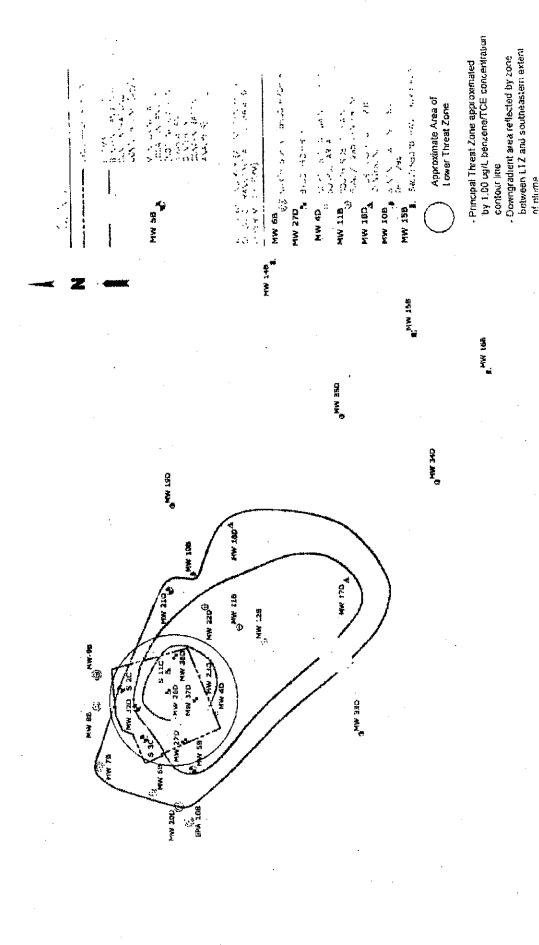
# **EPA Superfund Record of Decision:**

BRIDGEPORT RENTAL & OIL SERVICES EPA ID: NJD053292652 OU 02 BRIDGEPORT, NJ 09/27/2006

Table 9-1: Groundwater Preliminary Remediation Goals

Analyte/ Contaminant Group	NJDEP GWQC	Federal MCL
Volatile Organic Compounds (ug/L)	NOTE OF COLUMN TO SERVICE OF THE OWNER OWNER OF THE OWNER OWN	
Benzene	1	5
TCE	1	5
Vinyl Chloride	I I	2
Semi-Volatile Organic Compounds (ug/L)		
BCEE	7	
Polychlorinated Biphenyls (ug/L)	· •	·
Total PCBs	0.5	0.5

# Figure 6-3: Deep Groundwater Extent of Contamination (As Represented by Benzene and TCE Distribution)



#### Exhibit 9

Brosha Excerpt

brosha excerpt

# PUBLIC HEALTH ASSESSMENT BRIDGEPORT RENTAL AND OIL SERVICE LOGAN TOWNSHIP, GLOUCESTER COUNTY, NEW JERSEY CERCLIS NO. NJD053292652 JANUARY 23, 1995

Prepared by:

New Jersey Department of Health Environmental Health Service

Under Cooperative Agreement with the Agency for Toxic Substances and Disease Registry

#### **SUMMARY**

The 30 acre Bridgeport Rental and Oil Services, Inc. (BROS) site is located in southwest New Jersey, Logan Township, approximately 1 mile east of the town of Bridgeport and about 2 miles south of the Delaware River.

Groundwater and surface soils at the BROS site have been contaminated as a result of the mismanagement of large quantities of waste oil and other fluids. A second superfund site, Chemical Leaman Tank Lines, Inc., is located near the BROS site and may be contributing to groundwater contamination within the BROS study area.

Groundwater in the top two aquifers has been impacted by significant levels of contaminants. The contaminants include, among others: volatile organic compounds (VOC's), such as methylene chloride, trichloroethene, benzene, and vinyl chloride; semi-volatile organic compounds, bis (2-chloroethyl) ether; metals, lead and chromium; and polychlorinated biphenyls (PCBs).

Past exposure of residents living near the BROS lagoon to VOC's in residential well water is likely to have occurred from before 1983 through 1987. In the past, VOC's had migrated from the BROS lagoon through the shallow groundwater system. Before connections to the public water supply were completed, all of the known residents in the area were supplied with bottled water for drinking and other domestic purposes. A permanent source of water to 33 of these residents was provided by extension of the public water supply pipeline in April 1987.

The full expanse of the contamination plume at the BROS site has not been totally characterized. Groundwater contamination has been documented in off site monitoring wells and has been estimated to have moved as much as a mile from the site towards the southeast. It is not known the extent to which residents may have used this water in the past, how it is presently used, or how it may be used in the future.

It is possible that there are residents who are downgradient from the BROS lagoon who drank water from contaminated private wells and were chronically exposed to contaminants in those wells.

Based upon information reviewed, the BROS Site is considered a public health hazard because of past exposures, and is currently considered an indeterminate public health hazard. Human exposure to site related contaminants has occurred in the past. The indeterminate nature of the current health hazard is based on the fact that site related contaminants are migrating off-site and may have contaminated groundwater in an area where residential wells are in current use. The contaminants detected in this groundwater are of sufficient concentrations as to constitute a public health hazard. ATSDR's Health Activities Recommendation Panel has evaluated the BROS site to determine appropriate follow-up health actions. The panel determined that community health education is needed; however, the level of educational activity will be determined by the NJDOH following an evaluation of, among other factors, comments received during the public comment period for this



#### BACKGROUND

#### A. Site Description and History

Groundwater and surface soils at the Bridgeport Rental and Oil Services, Inc. (BROS) site have been contaminated as a result of the mismanagement of large quantities of waste oil and other fluids. Groundwater in the top two aquifers has been impacted by significant levels of contaminants. The contaminants include, among others: volatile organic compounds (VOC's), such as methylene chloride, trichloroethene, benzene, and vinyl chloride; semi-volatile organic compounds, bis (2-chloroethyl) ether; metals, lead and chromium; and polychlorinated biphenyls (PCBs). A groundwater contaminant plume is suspected to have moved as much as a mile southeast of the site.

Due to the contamination of drinking water wells near the BROS site, first identified in the early 1980's, the U.S. Environmental Protection Agency (USEPA) supplied an alternate water supply to 33 affected residents. A permanent source of water to these residents was provided by extension of the public water supply pipeline in April 1987.

In the early 1970's, the eastern dike of the BROS lagoon was breached and an adjacent 3 acre area was covered with a superficial layer of PCB contaminated oil.

In the spring of 1981, the level of the lagoon again threatened to overflow the dike. In response to the threat, the U.S. Coast Guard increased the height of the dike by 5 feet. When the dike again threatened to overflow in 1982 and 1983, the EPA took emergency action to lower the level of the lagoon by pumping and treating the aqueous phase layer and discharging the treated material into the Little Timber Creek.

The BROS site was included on the EPA's National Priorities List (NPL) in 1983. A Phase 1 Remedial Investigation/Feasibility Study (RI/FS) was conducted in 1983 and 1984 (1), and a Record of Decision (ROD) was signed in 1984 approving remedial activity at the site. EPA is currently performing a detailed Phase 2 RI/FS to determine the nature of the extent of the groundwater contamination and to evaluate remedial alternatives.

Remedial activities have included the demolition and removal of approximately 100 tanks and process vessels used to store hazardous wastes; off-site disposal of about 400,000 gallons of oils and sludges taken from the tanks; demolition and off-site disposal of buildings, drums, and other site debris; and the construction and operation of an aqueous wastewater treatment system (AWTS). In addition, an on-site transportable incinerator has been setup on the site. The incinerator is being used for the thermal destruction of the approximately 5000 cubic yards of PCB contaminated lagoon surface oil , and the 80,000 to 100,000 cubic yards of contaminated sludges, sediments, soils, and levee material.

The listed activities were implemented through an interagency agreement between USEPA and the U.S. Army Corps of Engineers (COE). The New Jersey Department of Environmental Protection and Energy (NJDEPE) provides support to the USEPA.

The Agency for Toxic Substances and Disease Registry (ATSDR) has been involved with this site in the past. A Health Assessment was preformed by the agency in March 1986 (2) and an Addendum to the Health Assessment was completed on August 9, 1991 (10). This addendum evaluated data collected in and adjacent to Swindell Pond. The New Jersey Department of Health (NJDOH), in cooperation with ATSDR, completed a Site Review and Update (SRU) on September 2, 1992 that recommended the site be revaluated through the public health assessment process.

The 30 acre BROS site is located in southwest New Jersey, Logan Township, approximately 1 mile east of the town of Bridgeport and about 2 miles south of the Delaware River (Figure # 1). The site borders an active peach orchard on its western edge. Located on Cedar Swamp Road, it's northern border, BROS is situated at the point where Route 130 and Interstate 295 diverge. The eastern edge of the site is a swampy area leading to the Little Timber Creek. Two "man-made" ponds are found adjacent to the lagoon, south and south west of the site (Figure # 2).

The site was originally a sand and gravel pit in the 1930's, and there is evidence of waste disposal activities beginning in the 1950's. When the current owners acquired the site in the late 1960's the site was already a waste oil storage and recovery operation. Today, waste handling activities are prohibited at the site by court order. Before remediation began, the site consisted of a tank Farm, drums, tank trucks and an approximately 13 acre waste oil lagoon. The lagoon was up to 25 feet deep in certain locations, with the bottom 13 feet of the lagoon in contact with the ground water. The liquid in the lagoon is divided into three phases including: an oily layer with drums, trash, and other debris floating in it; an aqueous layer; and sludge/sediment deposits on the bottom. At one time it was estimated that there were about 2.5 million gallons of oil in the lagoon. The depth of the sludge layer was estimated to be between 2 and 4 feet. Sampling of the oil and the sludge layers revealed average PCB concentrations in excess of 500 ppm. Beneath the sludge layer are contaminated groundwater and soils. Most of the oil has been removed from the lagoon and the water level has been lowered by pumping from 8.5 to 3.0 feet below mean sea level. The water level during remediation is held at or slightly below the natural water level in the superficial aquifer (Upper Potomac-Raritan-Magothy [PRM] Aquifer). The treated water is discharged to the Little Timber Creek and is monitored to ensure acceptable effluent standards. Contaminated soils excavated from the site are incinerated in the on-site thermal destruction facility.

Access to the BROS site is strictly limited. There is tall chain-link fence surrounding the entire site. The entrance to the on-site areas is controlled by a security guard.

#### B. Actions Implemented during the Health Assessment Process

Since the completion of the data review for this site, March 31, 1993, there have been no additional

actions implemented. The remedial work is continuing with a few minor incidents reported. These include: the discovery of additional contamination and short term release of organic vapor while workers were digging in the western levee of the lagoon; and the discovery of an apparent human hip bone in the same area.

#### C. Site Hydrogeology

The BROS lagoon is underlain by three aquifers and three confining units as follows: the Upper Potomac-Raritan-Magothy (PRM) Aquifer; an intermediate confining unit; the Middle PRM Aquifer; a continuous clay confining unit; the lower PRM Aquifer; and a basal confining unit defined by the Wissahikon Formation (Figure # 3). Only the Upper and Middle PRM aquifers have been investigated as part of the Phase II RI/FS (4). In order to accurately address the vertical and horizontal migration of contaminants from the lagoon, the middle PRM Aquifer has been separated into the Upper Middle PRM and the Lower Middle PRM Aquifers.

#### D. Site Visit

NJDOH (J. Pasqualo, N. P. Singh, and J. Winegar), ATSDR (A. Block and S. Jones; Region II), USEPA (Region II) and Gloucester County Health Department personnel visited the BROS site on December 1, 1992. The site was well secured with a chain-link fence, guarded, and posted with warning signs. The on-site area actually consists of two distinct areas: an administrative area consisting of numerous large office trailers; and an inner exclusion zone surrounding the lagoon itself where remediation is taking place. Both of these regions were examined during our visit. During our visit the following observations were made:

- The site is undergoing active remediation;
- The large storage lagoon was obviously heavily contaminated and stained with oil, with the level of the fluid approximately 10-15 feet below the lip of the levee;
- In the eastern portion of the lagoon numerous steel drums were observed in and protruding from the oil. Each drum appears to have been punctured. In addition, there is a tanker trailer, trash, and a partially submerged boat in this section;
- There is an aqueous wastewater treatment system (AWTS) and a large mobile thermal destruction facility located near the north east corner of the lagoon. The AWTS facility pumps and treats the aqueous layer of the lagoon to keep the level at or below the water table in the surface aquifer and to remove volatile organic and inorganic contaminants. The treated water is discharged into the Little Timber Creek. The incinerator is used to burn oil, contaminated

soil, and sediments from the lagoon. At the time of our site visit the incinerator was not operating due to a dispute between two of the subcontractors;

- On the north side of the lagoon were two large piles of incinerated soil and ash. This material is stockpiled and being used as part of the back fill material for the lagoon;
- Backfilling with a mixture of incinerated soil, clean fill, and cement mix, has begun in a small area in the south west corner of the lagoon;
- The area contaminated by a previous levee breach northeast of the lagoon has just been cleared of trees and brush as part of the initial remediation of this area. The perimeter fence has been extended around this region;
- There is an active peach orchard directly adjacent to the western border of the lagoon;
- South and south west of the BROS lagoon, within 100 feet, are two "manmade" ponds, known as Swindell Pond and Gaventa Pond, respectively. According to the local health officials, the ponds are used for swimming and fishing;
- Several residences were noticed about 800 feet north of the site. A single residence was observed about 150 feet to the west.
- At the north east corner of Gaventa Pond, the point closest to the BROS Lagoon, a oil boom was noticed in the corner. There was an obvious oil slick on the lagoon side of the boom. There were also several drums of unknown origin noticed in the pond; and
- The land area just south of the lagoon, and between the two ponds, showed evidence of human activity. Off-road bike tracks were visible throughout this area. Closer to Swindell Pond there was an area noticed where persons have been building campfires and drinking. This area was littered with beer cans.

#### E. Demographics, Land Use, and Natural Resources Use

#### **Demographics**

The population of Logan Township was listed as 5,147 in the 1990 U.S. census. The Township



consists of approximately 15,369 acres in northwest Gloucester County, New Jersey and consists of the following towns: Bridgeport; Gibbstown; Paulsboro; Swedesboro Repaupo; and Beckett. A large percentage of the residents are employed by the local industries.

There is at least one daycare center known to be near the BROS site. The facility is located approximately 1/4 mile south and west of the site (Figure # 4, Residential Well #44).

#### Land Use

The area surrounding the BROS site is predominately rural and agricultural. There is an active peach orchard adjacent to the western border of the site. A truck repair garage is located about 300 feet north east of the site and there are three homes about 800 feet north.

Approximately 1/2 of Logan Township is undeveloped, including marshes, vegetated areas, and woodlands. About 5,200 acres of the Township are covered by privately owned farms that cultivate vegetable crops and fruit. About 500 acres of land is devoted to residential use.

There is another EPA Superfund site located approximately 1200 feet west of the BROS Site (Figure #5). This site is known as the Chemical Learnan Tank Lines Site (CLTL), and it occupies 31.4 acres in Logan Township. The site includes a 14.1 acre tanker washing terminal, which has been operating since 1960. Contamination of the soil and the groundwater as a result of CLTL's activities has been confirmed. Contaminants largely consist of volatile organic compounds (VOC's).

#### **Natural Resources Use**

Residents near the BROS site use groundwater from either private wells or public water supplies for drinking and other household uses. As noted previously, the U.S. Environmental Protection Agency (USEPA) supplied an alternate water supply to 33 residents who reside just north of the site after it was first identified that their water was being contaminated by BROS. A permanent source of water was provided by extension of the public water supply pipeline by April 1987. Most of the residents to the south of the site, including those to the southeast in the projected path of the contamination plume moving in the upper middle aquifer, are utilizing private wells. Both the Upper PRM and Upper Middle PRM aquifers are used for this water source (private water). It is believed that most of these wells draw from the upper PRM aquifer (5).

Seven municipal supply wells lie within 4 miles of the BROS site. Five of the wells are screened within the Middle PRM; however, none of the wells appear to lie within the path of the BROS contaminant plume. Penns Grove Water Company wells in Bridgeport are hydraulically upgradient of the site. The Purulent Water Company wells, which lie southwest of BROS along Raccoon Creek, are hydraulically cross-gradient and are more than 3 miles away.



Only four private wells proximal to the BROS site are screened within the Middle PRM and appear to lie hydraulically downgradient and in the path of the contaminant plume. Wells 40, 43, 44, and 61 (Figure #4) are screened within the Middle PRM and lie within 1 mile of the southwest boundary of the site. The remaining domestic wells lying to the south-southeast and proximal to the site are screened in the Upper PRM aquifer and surficial Cape May Formation.

There is an active peach orchard directly adjacent to the western border of the lagoon, which according to the local health department, sells peaches for local consumption. It seems likely, although not documented, that water from Gaventa Pond is used as irrigation water for the peach orchard.

According to the Gloucester County Health Department (personal communication) local residents use the two "man-made" ponds, known as Swindell Pond and Gaventa Pond respectively, for swimming and fishing. No information exists, however, on the number of persons swimming or fishing in the area.

The Little Timber Creek flows north along the eastern edge of the BROS site and discharges into a small drainage ditch through a tidal gate north of Route 44. It eventually drains into an unnamed tidal channel that runs along the western edge of Cedar Swamp. Water in the tidal channel finally flows into the Delaware River.

#### F. Health Outcome Data

There are multiple sources of health outcome data in New Jersey. State and local data for heath outcome information include the New Jersey State Cancer Registry, Birth Defects Registry, Vital Statistics Records, Renal Dialysis Network, and Hospital Discharge Reports. Federal databases such as those maintained by the agencies within the US Department of Health and Human Services (i.e., National Cancer Institute, National Institute for Occupational Safety and Health, and ATSDR) are not site-specific, but may be used for comparison or evaluation purposes.

#### COMMUNITY HEALTH CONCERNS

Initially, the public was most concerned about drinking water contamination in the direct vicinity of the BROS lagoon, particularly north and west of the site. They were also upset with the amount of time the study of the site was taking before actual remediation of the site was to take place. They felt action should be taken immediately.

A recent well inventory (October 1992) of domestic/residential, commercial, and municipal supply wells that lie towards the south and southeast of the site has generated and renewed local community concern regarding contaminated drinking water. This survey was conducted by the USEPA.



In general, the public was opposed to the idea of on-site incineration. This was due, in part, to negative experience that they had with a nearby hazardous waste incinerator, and a perception that incineration is a poor and inefficient technology.

At least one local resident who lives about 800 feet north of the BROS lagoon had complained about odors emanating from the site.

The owner of Swindell Pond has, in the past, expressed concern about possible contamination of his pond by the adjacent BROS lagoon.

#### ENVIRONMENTAL CONTAMINATION AND OTHER HAZARDS

The tables in this section list the contaminants of concern. NJDOH evaluates these contaminants in the subsequent sections of the Health Assessment to determine whether exposure to them has public health significance. NJDOH selects and discusses these contaminants based upon the following factors:

- 1. Concentrations of contaminants on and off the site.
- 2. Field data quality, laboratory data quality, and sample design.
- 3. Comparison of on-site and off-site concentrations with health assessment comparison values for (1) noncarcinogenic endpoints and (2) carcinogenic endpoints.
- 4. Community health concerns.

In the data tables that follow the On-site Contamination subsection and the Off-site Contamination subsection (appendices), the listed contaminant does not mean that it will cause adverse health effects from exposures. When selected as a contaminant of concern in one medium, that contaminant will be reported in all media.

The Data tables include the following acronyms:

- CREG = ATSDR Cancer Risk Evaluation Guide
- EMEG = ATSDR Environmental Media Evaluation Guide
- RMEG = Reference Dose Media Evaluation Guide, calculated from EPA's reference dose (RfD).

■ LTHA = USEPA's Lifetime Health Advisory

■ NJ MCL =NJ Maximum Contaminant Level

■ PPB = Parts Per Billion

■ ND = Not Detected

Comparison values for public health assessments are contaminant concentrations in specific media that are used to select contaminants for further evaluation. These values include Environmental Media Evaluation Guides (EMEGs), Cancer Risk Evaluation Guides (CREGs), and other relevant guidelines. CREGs are estimated contaminant concentrations based on a one excess cancer in a million persons exposed over a lifetime. CREGs are calculated from EPA's cancer slope factors. Maximum contaminant levels (MCLs) represent contaminant concentrations that New Jersey or a Federal regulatory agency, e.g., EPA, deems protective of public health (considering the availability and economics of water treatment technology) over a lifetime (70 years) at an exposure rate of 2 liters of water per day. MCLs are regulatory concentrations. EPA's Reference Dose (RfD) is an estimate of the daily exposure to a contaminant that is unlikely to cause health effects.

The environmental contamination section includes sampling data from a variety of media sources including: groundwater (monitoring wells and residential wells); surface water; surface soil; subsurface soil; and sediments.

Local residential water supplies in the area of the BROS lagoon are derived primarily from the Upper PRM Aquifer and the Upper Middle PRM Aquifer. Due to this fact, the groundwater section of this public health assessment will focus primarily on these two aquifers.

The Upper PRM Aquifer is the water table aquifer. Originally the movement of groundwater in this aquifer was primarily to the north and west, however, due to the ongoing site remediation the local hydraulic gradient has reversed so that the groundwater currently moves back towards the lagoon (personal communication).

Groundwater flow in the Upper Middle PRM Aquifer is consistently toward the southeast, regardless of tidal

stage or season. Contaminant concentration in this aquifer is strongest (most contaminants, highest concentration) towards the southeast, along the direction of the groundwater flow.

As of the writing of this report the extent of the contaminant plume in the Upper Middle PRM Aquifer has not been delineated (6). The conservative estimate of the distance of contaminant transport from the lagoon in the Upper Middle PRM Aquifer was calculated to be 5,000 feet, based on 40 years of travel time and an average of the hydraulic conductivity of this aquifer. Swedesboro-Paulsboro Road is about 5,000 feet from the BROS lagoon.

TABLE # 3. CONTAMINANTS OF CONCERN; ON-SITE GROUNDWATER MONITORING WELLS; BROS SITE, LOGAN TWP., NJ., CONTAMINANTS WITH COMPARISON VALUES.

	CONCENTRATION  RANGE PPB				elowiparison	
CONDAMINANT	STATELON	De Gr		REF		ikye dike
1,2-Dibromo-3-chloro-propane	ND-10	. ND	9/90	5	20	EMEG
Benzene	ND-810	ND-95	9/90	5	1,2	CREG
Acetone	ND-77	17-33,000	9/90	5	1000	RMEG
1,1 Dichloroethene	ND	ND	9/90	5	.06	CREG
Cis-1,2 Dichloro- ethene	ND-3	ND-100	9/90	5	70	LTHA
1,2 Dichloroethane	ND	ND	9/90	5	0.4	CREG
Methylene Chloride	ND-70	ND-579	9/90	5	5	CREG
Trichloroetheue	ND	ND-1000	9/90	5	3	CREG
Tetrachloroethene	ND-2	ND-25	9/90	5	0.7	CREG
1,1,2,2-Tetrachloroethane	ND	ND	9/90	5	0.2	CREG
Vinyl Chloride	ND	ND	9/90	5	0.2	EMEG
Benzo (A) Pyrene	ND-15	ND	9/90	5_	.005	CREG
ВЕНР	ND-120	ND	9/90	5	3	CREG
BCEE	ND-11	ND-780	9/90	5	.03	CREG
Isophorone	ND-35	ND-1300	9/90	5	9	CREG
Naphthalene	ND-180	ND-25	9/90	5	20	LTHA
Manganese	292-6040	79.1- 6240	9/90	5	50	RMEC
Barium	ND-1200	ND	9/90	5	700	RMEC
Zinc	ND	ND- 567,000	9/90	5	3000	RMEC
Chromium	120-1730	ND	9/90	5	10	NJ MC

<sup>\* =</sup> Concentration range given for both the shallow and the deep aquifers

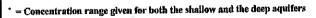
ND = Not Detected

Source: Phase II RI/FS

35

TABLE #4. CONTAMINANTS OF CONCERN; OFF-SITE GROUNDWATER; BROS SITE, LOGAN TWP.,NJ. CONTAMINANTS WITH COMPARISON VALUES.

		рудоў — году			COMPA	RISON VALUE
CONTAMINANT			DATE	REF	200	Source
1,2-Dibromo-3- chloro-propane	ND	ND	9/90	5	20	EMEG
Benzene	ND	ND-30	9/90	5	1.2	CREG
Acetone	ND-630	ND-3500	9/90	5	1000	RMEG
1,1 Dichloroethene	ND	ND9	9/90	5	.06	CREG
Cis-1,2 Dichloro- ethene	ND	ND-28	9/90	5	70	LTHA
1,2 Dichloroethane	ND	ND-17	9/90	5	0.4	CREG
Methylene Chloride	ND-16	ND	9/90	5	5	CREG
Trichloroethene	ND	ND-35	9/90	5	3	CREG
Tetrachloroethenc	ND	ND-3	9/90	5	0.7	CREG
1,1,2,2-Tetrachloroethane	ND	ND-2	9/90	5	0.2	CREG
Vinyl Chloride	ND ·	ND-37	9/90	5	0.2	EMEG
Benzo (A) Pyrene	ND	ND	9/90	5	.005	CREG
ВЕНР	1-5	ND-2	9/90	5	3	CREG
BCEE	ND	ND-380	9/90	5	.03	CREG
Isophorone	ND	ND-1	9/90	5	9	CREG
Naphthalene	ND	ND-5	9/90	5	20	LTHA
Manganese	249-2040	ND-293	9/90	5	50	RMEG
Barium	ND	ND	9/90	5	700	RMEG
Zinc	ND	ND- 16,800	9/90	5	3000	RMEG
Chromium	ND-234	ND	9/90	5_	10	NJMCI



ND = Not Detected

Source: Phase II RI/FS

TABLE # 5. RANGE OF CONTAMINANTS IN ON-SITE GROUNDWATER, CONTAMINANTS WITHOUT ATSDR COMPARISON VALUES.

CONTAMINANT	GONG:NI		
CONTAMINANT	SENE CW		REPERCISE
1,2,4 Trimethylbenzene	ND-360	ND	
1,3,5 Trimethylbenzene	ND-250	ND	5
1,2 Dibromoethane	ND-36	ND	5
2-Butanone	ND-27	ND-2	5
2-Hexanone	ND-93	ND	5
N-Propylbenzene	ND-59	ND	5
N-Butylbenzene	ND-28	. ND	5
P-Isopropyltoluene	ND-15	ND	5
Sec-Butylbenzene	ND-5	ND	5
2-Methylnapthalene	ND-74	ND	5
4-Methylphenol	ND-130	ND-160	5
Benzo (B) Fluoranthrene	ND-13	ND	5
Chrysene	ND-32	ND	5
Di-N-Octylphthalate	ND-12	ND	5
Dibenzofuran	ND-16	ND	5
Phenanthrene	ND-130	ND	5
Benzylalcohol	ND-2,800	ND	5
4-methyl-2-Pentanone	ND	ND-2,000	5
Lead	ND-1380	ND-2920	5

<sup>\* =</sup> Concentration range given for both the shallow and the deep aquifers

ND = Not Detected

Source: Phase II RI/FS

TABLE #6. RANGE OF CONTAMINANTS IN OFF-SITE GROUNDWATER, CONTAMINANTS WITHOUT ATSDR COMPARISON VALUES.

	CONGENI THE RANGE		
CONTAMINANT			REFERENCE
1,2,4 Trimethylbenzene	ND-11	ND-3	5
1,3,5 Trimethylbenzene	ND	ND7	5
1,2 Dibromoethane	ND	ND	
2-Butanone	ND	ND5	5
2-Hexanone	ND	ND	5
N-Propylbenzene	מא	ND	5
N-Butylbouzene	ND_	ND	5
P-Isopropyltoluene	ND .	ND	5
Sec-Butylbenzene	ND	ND	5
2-Methylnapthalene	ND	ND	
4-Methylphenol	ND	ND	5
Benze (B) Fluoranthrene	ND	ND	5
Chrysene	ND	ND	5
Di-N-Octylphthalate	ND	ND	5
Dibenzofuran	ND	ND	5
Phonanthrene	ND	ND	5
Benzylaicohol .	ND	ND	5
4-methyl-2-Pentanone	ND	ND	5
Lead	ND-140	ND-55.6	5

\* = Concentration range given for both the shallow and the deep aquifers

ND = Not Detected

Source: Phase II RI/FS

TABLE #7. CONTAMINANTS IN 11 OFF-SITE RESIDENTIAL GROUNDWATER (1983 TO 1985).

		of Malatylet use bit bloomil and were the second	Autoroxy ales, and
CONCENTRATIONERS		SOURCE	
20	200	RMEG	NO
1.9	0.4	CREG	YES
7.1	.005	MCL	YES
. 1.9			NA
170	0.2	EMEG	YES
	20 1.9 7.1	20 200 1.9 0.4 7.1 .005 1.9 1.70 0.2	PRB   SOURCE

Table #8. RANGE OF CONTAMINATION IN Off-SITE GROUNDWATER, RESIDENTIAL WELLS, SEPTEMBER 1992.\*

		111 1250 4 1890 8		2 2	LTUME! FOMPA	
Wolsen	GONPAMINAN	PDT	+ DVIIIC	REF		(SOJ)(KG)()
RW-40	Cis-1,2-dichloro- ethane	9	10/92	15	70	LTHA
RW-40	Trichloroethene	120	10/92	15	3	CREG
RW-65	Dieldrin	.025	10/92	15	.002	CREG
RW-44	Aroclor 1260	0.8	10/92	15	.002	MCL

<sup>\*</sup> Resampling of contaminated wells by the USEPA in 1993 was not able to confirm the presence of the contaminants shown in table.

#### Exhibit 10

**PRGs Delaware Sand & Gravel** 

#### Exhibit 11

**Artesian November 2000 BCEE Incident** 

#### Exhibit 12

Artesian Llangollen 2003 PWSSWA

EPA/ROD/R03-88/048 1988

# **EPA Superfund Record of Decision:**

DELAWARE SAND & GRAVEL LANDFILL EPA ID: DED000605972 OU 01 NEW CASTLE, DE 04/22/1988

COMPOUND	ACCEPTABLE WATER EXPOSURE (UG/1)	SOURCE	KOC.
TOLUENE (9)	2000	MCLG	300
METHYLENE CHLORIDE	0.7	CAGUCR	8.8
ACETONE (9)	3500	SPHEM	2,2
4-METHYL-2PENTANONE (	9) 1750	IRIS (AIC)	4.5**
ETHYLBENZENE (9)	680	MCLG	1100
1,2-DICHLOROETHANE	5	MCL	14
XYLENE (9)	2	MCL	240
PHENOL (9)	3500	IRIS (ORAL) (RFD)	14.2
BIS (2-CHLOROETHYL) E	THER 0.03	(10-6) RISK	13.9
NAPHTHALENE (8)	. 400	ENVIRON RED	1072 A
4-METHYLPHENOL (9)	1750	SPHEM	246 A.
2-METHYLPHENOL (9)	1750	SPHEM	* A
2-BUTANONE	1750	IRIS	35

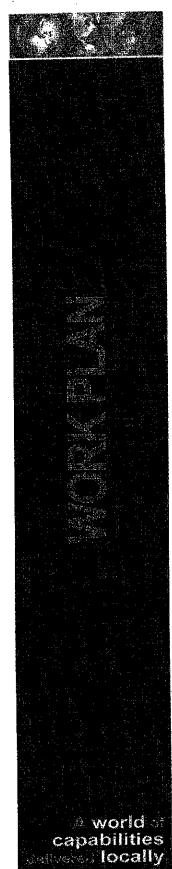
<sup>1</sup> ACCEPTABLE SOIL CONCENTRATIONS CALCULATED FROM PRODUCT OF ACCEPTABLE WATER EXPOSURE AND KOC VALUE, FOLLOWED BY TEN-FOLD ALLOWANCE FOR DILUTION, I.E. (ACCEPTABLE WATER EXPOSURE ) X (KOC) X (10)

EQUATION FOR NAPTHALENE: LOG KOC = 1.00 LOG KOW - 0.21 EQUATION FOR 4-METHYLPHENOL: LOG KOC = 0.544 LOG KOW + 1.377

A = KOC CALCULATED FROM KOC/KOW RELATIONSHIPS REPORTED IN LYMAN ETAL. (1982)

<sup>\*</sup> = NO KOW DATA AVAILABLE FOR 2-METHYLPHENOL; ASSUME VALUE SIMILAR TO 4-METHYLPHENOL

<sup>\*\* =</sup> NO KOW DATA AVAILABLE - ASSUME VALUE SIMILAR TO MIBK PEROXIDE



#### FEASIBILITY STUDY WORK PLAN – REVISION 2

Delaware Sand & Gravel Superfund Site New Castle, Delaware

Submitted To: United States Environmental Protection Agency

Region III

1650 Arch Street

Philadelphia, Pennsylvania 19103-2029 USA

Submitted By: Delaware Sand & Gravel Trust

100 East Market Street

Suite 1

Newport, Delaware 19804 USA

Prepared By: Golder Associates Inc.

670 Commercial Street, Suite 103

Manchester, New Hampshire 03101 USA

#### Distribution:

2 copies and 1 CD - Debra Rossi - USEPA Region III

1 copy and 1 CD - John Cargill - DNREC

1 copy and 1 CD - Aaron Frantz - CDM

1 copy and 1 CD - Michael Barbara

1 copy and 15 CDs - DS&G Trust

4 copies and 5 CDs - Golder Associates

October 2011 Revision 2 Project No.: 013-6052-006





Project No.: 013-6052-006

Table 7-1: Preliminary Remediation Goals for Groundwater Delaware Sand & Gravel Superfund Site New Castle, Delaware

Contaminant of Concern	Preliminary Remediation Goal (μg/L)	Maximum Contaminant Level (µg/L)
Benzene	5	5
bis(2-Chloroethyl)ether	0.1	0.1*
bis(2-Ethylhexyl)phthalate	6	6
1,2-Dichloroethane	5	5
cis-1,2-Dichloroethene	11	70
Ethylbenzene	30	700
Toluene	460	1,000
Tetrachloroethene	0.8	5
Trichloroethene	0.18	5
Vinyl chloride	2	2
Iron	10,900	300**
Manganese	1,900	50**

#### Notes:

- 1) Preliminary Remediation Goals for Groundwater for Delaware Sand & Gravel Landfill Operable Unit 6 as provided by USEPA on June 29, 2011
- 2) PRG = preliminary remediation goal
- 3) MCL = maximum contaminant level
- 4) \* indicates Delaware's Interim Advisory Level for BCEE in drinking water
- 5) \*\* indicates Secondary MCL
- 6) µg/L = micrograms per liter

Prepared by: RWB
Checked by: TAM





Artesian Resources Comporation services P.O. Box 15004 • Wilmington, DE 19850 • (302) 453-6953 • Fax: (302) 453-6957

FOR IMMEDIATE RELEASE Contact: Joseph A. DiNunzio (302) 453-6900

# <u>ARTESIAN'S WATER NEVER UNSAFE TO DRINK BASED ON EXISTING EPA STANDARDS</u>

NOVEMBER 11, 2000—Artesian Water Company today disputed the characterization in a November 11, 2000 News Journal report that the established by the Environmental Protection Agency (EPA). Dian C. Taylor, CEO & President of Artesian, stated that "We are dedicated to available and prepared, upon request, to meet with any civic association, maintenance corporation and/or community group to answer any providing quality water, and we will continue to do what is necessary to ensure a safe, reliable water supply for all our customers. We are water Artesian has provided to the public was unsafe. Artesian supplied water to its customers within the parameters for safe water as further questions.

Division of Public Health (DPH) that levels of bis (2-chloroethyl) ether (BCEE) exceeded the EPA's action level of 0.96 parts per billion, On November 1, 2000, Artesian voluntarily removed its Llangollen well field from service immediately upon notice from the Delaware which has been the level set for the protection of public health and the environment.

in the human body. It metabolizes rapidly and passes out of the body in urine. The DPH has issued an Information Sheet on BCEE which may According to DPH, exposure to BCEE at this low concentration causes no short-term health effects. This compound also does not accumulate be obtained by contacting them.

The possible effects of long-term exposure to BCEE are based upon limited studies on mice that indicated a possible increase in risk for liver



consistency or statistical power to permit a conclusion regarding the presence or absence of a causal association between exposure and cancer, performed to date. The IARC classification of "inadequate evidence" is because, per IARC: "The available studies are of insufficient quality, tumors. Based on these limited studies, the EPA has concluded that a person drinking two liters of water per day for a period of 30 years carcinogen" if consumed at the rate of the two liter of water per day for a period of 30 years. According to the International Agency for Cancer Research (IARC), there is "inadequate evidence for carcinogenicity in humans" from exposure to BCEE based on the studies might have a 1 in 10,000 increased risk for such cancer. Based on these studies, the EPA has classified BCEE as a "probable human or no data on cancer in humans are available."

EPA's action level of 0.96 ppb in order to assure protection of public health under standards set in an appropriately conservative manner by Artesian acted to remove its Llangollen well field from service immediately upon notification of levels of BCEE of 1.0 ppb, just above the EPA. Artesian is proceeding with the installation of an activated carbon treatment system, which will eliminate BCEE from any water supplied to its customers.

standard does not mean that water was unsafe at the standard of 0.96 ppb applied by EPA and followed by Artesian. Artesian expects that this Region 3, DPH, Artesian and the Delaware Department of Natural Resources and Environmental Control on October 11, 2000. The proposed proposed standard will apply to all drinking water in Delaware once approved in the course of the normal regulatory standard setting process. standard for Delaware of 0.096 ppb for BCEE. DPH did not propose this more conservative standard when first notified of the presence of BCEE in Artesian's drinking water in April 2000, nor at meetings held with EPA Region 3, DPH and Artesian in June 2000 and with EPA After Artesian's voluntary removal of its well field from service, DPH has proposed a new and much more conservative drinking water

Sanitation Foundation. The DPH has committed to expedite their review and approval required for the installation of this treatment system. The use of carbon treatment to remove chemicals such as BCEE is the only approved industry practice and is approved by the National

Artesian expects that installation of the treatment equipment at the Llangollen site will begin the week of November 27 and that the well field will be back in service by the end of the year after test results confirm the absence of BCEE in water delivered to customers.

published in the spring and includes all testing results from the previous year. A copy of our latest Water Quality Report may be found on our Artesian regularly tests its water for more than 130 regulated and unregulated substances. As required by the Federal Safe Drinking Water Act, Artesian reports any detection of these substances in a "Water Quality Report" that is distributed to all customers. This report is web site, or you may call our Customer Service Department at (302) 453-6930.

##

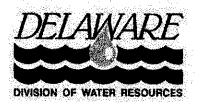
Public Water Supply Source Water Assessment for

Artesian Water Co. (Llangollen)

PWS ID: DE0000552

New Castle County, Delaware





Final Report: December 31, 2003

State of Delaware
Department of Natural Resources and Environmental Control
Division of Water Resources
Source Water Assessment and Protection Program
89 Kings Highway
Dover, Delaware 19901

Phone: (302) 739-4793

fax: (302) 739-2296

#### **Artesian Water Company**

#### Llangollen Wellfield

Welf Number	Sample Date	Contaminant Class	Analyte	Sample Result µg/i*	(P/S)MCL** "If applicable µg/i	(P/S)MCL. Violation?	Greater than 50% of MCL	Synthetic Substance Detection?
	8/22/2001	Other Organics	BCEE (Bis(2-chloroethyl) Ether)	0.044	0.096 <sup>8</sup>			Yes
Well 2	8/22/2001	Other Organics	Chloroform	0.2J	80 <sup>A</sup>	200	4.00	Yes
(35081)	8/22/2001	Pesticides	Dacthal	0,75	37 <sup>B</sup>			Yes
l ' '	8/22/2001	Petroleum Hydrocarbons	MTBE	1.0	10		900	Yes
	8/22/2001	Other Inorganics	pН	6.04	6.5-8.5 units			

Well Number	Sample Date	Contaminant Class	Analyte	Sample Result µg/l*	(P/S)MCL** 'If applicable µg/l	(P/S)MCL Violation?	Greater than 50% of MCL	Synthetic Substance Detection?
	8/22/2001	Other Organics	Cis-1,2-DCE (Cis-1,2-Dichloroethylane)	0.2J	70			Yes
	8/22/2001	Other Organics	1,2-DCA (1,2-Dichloroethane)	6.0	5			
	8/22/2001	Other Organics	Trichlorofluoromethane	2.4	80 <sup>A</sup>		1200	Yes
	8/22/2001	Other Organics	Bromoform	0.2J	80 <sup>A</sup>	and the second		Yes
Well 7	8/22/2001	Other Organics	Chloroform	D.2J	80 <sup>A</sup>			Yes
(10049)	8/22/2001	Other Organics	BCEE (Bis(2-chloroethyl) Ether)	0.6	0.096 <sup>B</sup>			The Agreement
	8/22/2001	Other Organics	PCE (Tetrachloroethylene)	5.8	5			
	8/22/2001	Pesticides	Dacthal	1.1	37 <sup>8</sup>			Yes
	8/22/2001	Petroleum Hydrocarbons	MTBE	0.1J	10			Yes
	8/22/2001	Other Inorganics	pH	5,76	6,5-8,5 units			

Well Number	Sample Date	Contaminant Class	Analyte	Sample Result µg/l*	(P/S)MCL** "If applicable µg/I	(P/S)MCL Violation?	Greater than 56% of MCL	Synthetic Substance Detection?
	8/22/2001	Other Organics	Cls-1,2-DCE (Cis-1,2-Dichloroethylene)	0.1J	70			Yes
		Other Organics	1,1-DCA (1,1-Dichloroethane)	0.1J	5			Yes
	8/22/2001	Other Organics	1,2-DCA (1,2-Dichloroethane)	0.4J	5			Yes
		Other Organics	Chlorobenzene	0.4J	100	1000	37 77 8 20 20 75	Yes
		Other Organics	Diethyl Ether	0.6J	N/A			Yes
Well G-3	8/22/2001	Other Organics	Trichloroffuoromethane	2.4	80 <sup>A</sup>	<b>经</b> 有基础		Yes
(10050)	8/22/2001	Other Organics	Bromoform	1.7	80 <sup>A</sup>		100	Yes
		Other Organics	Dibromochloromethane	0.6	80 <sup>A</sup>	100	1.00	Yes
	8/22/2001	Other Organics	Bromodichloromethane	0.2J	80 <sup>A</sup>		Service Control	Yes
	8/22/2001	Other Organics	BCEE (Bis(2-chloroethyl) Ether)	3.05	0.096 <sup>9</sup>			MARINE.
	8/22/2001	Other Organics	TCE (Trichloroethylene)	1.7	5		4.5	Yes
	8/22/2001	Pesticides	Dacthal	0.17	37 <sup>8</sup>	100		Yes
	8/22/2001	Other Inorganics	pH	5.76	6.5-8.5 units		200	100

Weli Number	Sample Date	Contaminant Class	Analyte	Sample Result µg/l*	(P/S)MCL** 'If applicable µg/I	(P/S)MCL Violation?	Greater than 50% of MCL	Synthetic Substance Detection?
Well K-1	8/22/2001	Other Organics	BCEE (Bis(2-chloroethyl) Ether)	0.010	0,096 <sup>B</sup>			Yes
	8/22/2001	Other Organics	TCE (Trichloroethylene)	2.4	5		4	Yes
(10052)	8/22/2001	Other Organics	Carbon Tetrachloride	0.1J	80 <sup>A</sup>		Salah Salah Salah	Yes
	8/22/2001	Other inorganics	рH	5.51	6.5-8.5 units		Tree Land	200

Well Number	Sample Date	Contaminant Class	Analyte	Sample Result µg/i*	(P/S)MCL** 'Wappilcable µg/I	(P/S)MCL Violation?	Greater than 50% of MCL	Synthetic Substance Detection?
	8/22/2001	Other Organics	Dibromochloromethane	3.4	80 <sup>A</sup>	1.7		Yes
	8/21/2001	Other Organics	Bromodichloromethane	1.1	80 <sup>A</sup>	Service Page		100
Well ASR	8/22/2001	Other Organics	Bromoform	7.1	80 <sup>A</sup>		7. 7. 1. 1.	Yes
(156408)	8/22/2001	Other Organics	Chloroform	0.3J	80 <sup>A</sup>		2 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	Yes
(100400)	8/22/2001	Other Organics	BCEE (Bis(2-chloroethyl) Ether)	0.0296	0.096 <sup>B</sup>			Yes
	8/22/2001	Other Inorganics	рH	6.28	6.5-8.5 units		10.75	



<sup>\*:</sup> Results are in micrograms/liter unless listed otherwise
\*\*: PMCL = Primary Maximum Contaminant Level, SMCL = Secondary Maximum Contaminant Level
A: Bromoform, Chloroform, Bromodichloromethana, Dibromochloromethana, and Trichlorofluoromethana are constituents of the Trihalomethana Group which has a total PMCL of 80 µg/l.
B: Detaware Uniform Risk-Based Standard or Proposed State-instituted MCL
C: pH is a widespread problem in Delaware that will not be used to adjust susceptibility
J: Laboratory Qualifier indicating that substance was present though the quantity is at levels just above the detection limits of the equiment.