

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071415\
 Data File : BG017880.D
 Acq On : 15 Jul 2015 5:33
 Operator : TP/UM
 Sample : MDL-04-W
 Misc : MDL-WATER-SVOC-4PPM
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleID :
 MDL-04-W

Manual Integrations
 APPROVED

apatel
 7/16/2015 8:19:26 AM

Quant Time: Jul 15 08:48:28 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG071415.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 15 07:22:18 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	66671	20.00	ng/ul	0.00
18) Naphthalene-d8	10.37	136	308192	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.24	164	174555	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.99	188	357315	20.00	ng/ul	0.00
77) Chrysene-d12	21.20	240	350936	20.00	ng/ul	0.00
85) Perylene-d12	23.42	264	328240	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.12	96	9130	4.90	ng/uL	0.00
5) Phenol-d5	6.77	99	166508	28.95	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.93	67	117335	29.85	ng/ul	0.00
9) 2-Chlorophenol-d4	7.12	132	132620	29.70	ng/ul	0.00
13) 4-Methylphenol-d8	8.29	113	128259	29.10	ng/ul	0.00
19) Nitrobenzene-d5	8.74	128	74396	30.94	ng/ul	0.00
22) 2-Nitrophenol-d4	9.46	143	53294	29.52	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.99	165	102498	27.67	ng/ul	0.00
29) 4-Chloroaniline-d4	10.51	131	198076	37.14	ng/ul	0.00
43) Dimethylphthalate-d6	13.66	166	404820	33.26	ng/ul	0.00
46) Acenaphthylene-d8	13.93	160	513790	32.49	ng/ul	0.00
51) 4-Nitrophenol-d4	14.45	143	64320	28.01	ng/ul	0.00
57) Fluorene-d10	15.24	176	379433	33.52	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.36	200	30027	17.96	ng/ul	0.00
70) Anthracene-d10	17.09	188	588145	36.32	ng/ul	0.00
78) Pyrene-d10	19.40	212	639138	38.67	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.28	264	553950	34.43	ng/ul	0.00

Target Compounds

					Qvalue
2) 1,4-Dioxane	3.16	88	1492	0.75 ng/uL#	85
4) Benzaldehyde	6.74	77	13810	4.10 ng/ul	96
6) Phenol	6.79	94	17223	2.73 ng/ul	98
8) Bis(2-Chloroethyl)ether	7.02	93	15481	3.01 ng/ul	93
10) 2-Chlorophenol	7.15	128	11160	2.47 ng/ul	92
11) 2-Methylphenol	8.03	108	10994	2.40 ng/ul	89
12) 2,2'-oxybis(1-Chloropropan	8.13	45	21080	2.82 ng/ul#	94
14) Acetophenone	8.40	105	21607	2.95 ng/ul	100
15) N-Nitroso-di-n-propylamine	8.39	70	12723	2.81 ng/ul	89
16) 4-Methylphenol	8.36	108	12518	2.51 ng/ul	98
17) Hexachloroethane	8.66	117	5710	2.65 ng/ul	94
20) Nitrobenzene	8.78	77	17355	2.74 ng/ul#	88
21) Isophorone	9.30	82	31649	2.67 ng/ul#	96
23) 2-Nitrophenol	9.49	139	5371	2.50 ng/ul	93
24) 2,4-Dimethylphenol	9.56	107	13058	2.38 ng/ul	98
25) Bis(2-Chloroethoxy)methane	9.80	93	17430	2.61 ng/ul	96
27) 2,4-Dichlorophenol	10.01	162	9494	2.54 ng/ul#	80
28) Naphthalene	10.42	128	45418	2.85 ng/ul	98
30) 4-Chloroaniline	10.53	127	15487	2.80 ng/ul	89
31) Hexachlorobutadiene	10.71	225	6830	2.90 ng/ul#	85
32) Caprolactam	11.31	113	3383m	1.24 ng/ul	
33) 4-Chloro-3-methylphenol	11.67	107	10405	2.27 ng/ul#	95
34) 2-Methylnaphthalene	12.04	142	29040	2.68 ng/ul	92

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 1,2,4,5-Tetrachlorobenzene	12.42	216	13094	2.85	ng/ul	90
37) Hexachlorocyclopentadiene	12.41	237	4508	1.89	ng/ul#	96
38) 2,4,6-Trichlorophenol	12.66	196	5850	2.35	ng/ul	94
39) 2,4,5-Trichlorophenol	12.72	196	5779m	2.13	ng/ul	
40) 1,1'-Biphenyl	13.07	154	37759	2.79	ng/ul	97
41) 2-Chloronaphthalene	13.11	162	29841	2.88	ng/ul	90
42) 2-Nitroaniline	13.32	65	5872	1.93	ng/ul#	78
44) Dimethylphthalate	13.71	163	40083	3.14	ng/ul	96
45) 2,6-Dinitrotoluene	13.82	165	4911	2.11	ng/ul#	84
47) Acenaphthylene	13.96	152	53175	3.04	ng/ul	97
48) 3-Nitroaniline	14.15	138	5721	2.15	ng/ul	90
49) Acenaphthene	14.31	153	33125	2.82	ng/ul	98
50) 2,4-Dinitrophenol	14.36	184	663	0.68	ng/ul#	42
52) 4-Nitrophenol	14.46	109	4665	2.40	ng/ul	81
53) Dibenzofuran	14.64	168	45063	2.91	ng/ul	99
54) 2,4-Dinitrotoluene	14.61	165	8255	2.47	ng/ul	100
55) 2,3,4,6-Tetrachlorophenol	14.87	232	4544	2.17	ng/ul#	94
56) Diethylphthalate	15.08	149	37440	2.85	ng/ul	98
58) Fluorene	15.30	166	39428	2.93	ng/ul	88
59) 4-Chlorophenyl-phenylether	15.30	204	17895	2.86	ng/ul	93
60) 4-Nitroaniline	15.32	138	7909	2.50	ng/ul	92
63) 4,6-Dinitro-2-methylphenol	15.38	198	2788	1.50	ng/ul#	67
64) N-Nitrosodiphenylamine	15.51	169	31195	2.87	ng/ul	99
65) 4-Bromophenyl-phenylether	16.20	248	10350	3.02	ng/ul	83
66) Hexachlorobenzene	16.30	284	11863	3.19	ng/ul	90
67) Atrazine	16.47	200	9769	2.62	ng/ul	95
68) Pentachlorophenol	16.65	266	2739	1.49	ng/ul#	57
69) Phenanthrene	17.04	178	62781	3.18	ng/ul	97
71) Anthracene	17.13	178	65454	3.24	ng/ul	97
72) 1,2,3,4-Tetrachlorobenzene	13.03	216	11898	2.62	ng/uL	92
73) Pentachlorobenzene	14.57	250	13316	3.05	ng/uL	95
74) Carbazole	17.41	167	56774	3.22	ng/ul	97
75) Di-n-butylphthalate	17.98	149	56066	2.54	ng/ul	97
76) Fluoranthene	19.06	202	67071	3.35	ng/ul#	97
79) Pyrene	19.43	202	77274	3.58	ng/ul#	97
80) Butylbenzylphthalate	20.35	149	16396	1.88	ng/ul	96
81) 3,3'-Dichlorobenzidine	21.12	252	9091	1.73	ng/ul#	88
82) Benzo(a)anthracene	21.18	228	64039	3.31	ng/ul	100
83) Bis(2-ethylhexyl)phthalate	21.13	149	21746	1.80	ng/ul	98
84) Chrysene	21.24	228	63971	3.44	ng/ul	97
86) Di-n-octyl phthalate	22.00	149	29348	1.53	ng/ul	100
87) Benzo(b)fluoranthene	22.75	252	54025	2.98	ng/ul	95
88) Benzo(k)fluoranthene	22.79	252	54718	2.93	ng/ul	93
90) Benzo(a)pyrene	23.32	252	58914	3.24	ng/ul	96
91) Indeno(1,2,3-cd)pyrene	25.65	276	53843m	2.77	ng/ul	
92) Dibenzo(a,h)anthracene	25.67	278	45587	2.75	ng/ul	93
93) Benzo(g,h,i)perylene	26.34	276	44758	2.75	ng/ul	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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