

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
 Data File : BG028064.D
 Acq On : 28 Jul 2017 20:05
 Operator : SJ/JU
 Sample : MDL-S-ML-07
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MDL-S-ML-07

Manual Integrations
 APPROVED

Sohil
 7/31/2017 11:54:02 AM

Quant Time: Jul 29 00:10:24 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 28 19:32:36 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	31381	20.00	ng/ul	0.00
18) Naphthalene-d8	11.19	136	135752	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.97	164	106397	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.72	188	311465	20.00	ng/ul	0.00
78) Chrysene-d12	22.06	240	422456	20.00	ng/ul	0.00
86) Perylene-d12	25.58	264	425093	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.85	96	3403	7.03	ng/uL	0.00
5) Phenol-d5	7.51	99	61583	25.98	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.68	67	33228	28.04	ng/ul	0.00
9) 2-Chlorophenol-d4	7.89	132	58275	29.16	ng/ul	0.00
13) 4-Methylphenol-d8	9.06	113	56585	26.74	ng/ul	0.00
19) Nitrobenzene-d5	9.53	128	28724	30.27	ng/ul	0.00
22) 2-Nitrophenol-d4	10.26	143	37004	30.62	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.80	165	69502	29.22	ng/ul	0.00
29) 4-Chloroaniline-d4	11.32	131	80839	33.70	ng/ul	0.00
44) Dimethylphthalate-d6	14.36	166	308544	32.04	ng/ul	0.00
47) Acenaphthylene-d8	14.67	160	319646	30.69	ng/ul	0.00
52) 4-Nitrophenol-d4	15.15	143	39122	27.05	ng/ul	0.00
58) Fluorene-d10	15.96	176	273328	30.82	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	16.07	200	60933	27.47	ng/ul	0.00
71) Anthracene-d10	17.82	188	470350	31.53	ng/ul	0.00
79) Pyrene-d10	20.10	212	590224	31.45	ng/ul	0.00
90) Benzo(a)pyrene-d12	25.33	264	599377	30.65	ng/ul	0.00

Target Compounds

					Qvalue
2) 1,4-Dioxane	3.89	88	725	1.397 ng/uL#	82
4) Benzaldehyde	7.51	77	1080	0.921 ng/ul	99
6) Phenol	7.54	94	7433m	3.233 ng/ul	
8) Bis(2-Chloroethyl)ether	7.77	93	6159	3.785 ng/ul	88
10) 2-Chlorophenol	7.93	128	7026	3.817 ng/ul#	80
11) 2-Methylphenol	8.79	108	6253	3.489 ng/ul#	65
12) 2,2'-oxybis(1-Chloropropan	8.89	45	6814	3.190 ng/ul#	92
14) Acetophenone	9.19	105	11659	3.793 ng/ul	87
15) N-Nitroso-di-n-propylamine	9.16	70	3422	2.371 ng/ul#	89
16) 4-Methylphenol	9.12	108	6339	3.188 ng/ul	96
17) Hexachloroethane	9.46	117	2813	3.820 ng/ul#	70
20) Nitrobenzene	9.58	77	8048	3.950 ng/ul#	93
21) Isophorone	10.10	82	14121	3.627 ng/ul#	95
23) 2-Nitrophenol	10.29	139	4721	4.054 ng/ul	95
24) 2,4-Dimethylphenol	10.33	107	8312	3.569 ng/ul	90
25) Bis(2-Chloroethoxy)methane	10.57	93	8454	3.738 ng/ul#	91
27) 2,4-Dichlorophenol	10.83	162	8708	3.842 ng/ul#	93
28) Naphthalene	11.24	128	23300	3.803 ng/ul	97
30) 4-Chloroaniline	11.34	127	6334m	2.774 ng/ul	
31) Hexachlorobutadiene	11.51	225	7879	4.018 ng/ul	98
32) Caprolactam	12.13	113	2563m	3.523 ng/ul	
33) 4-Chloro-3-methylphenol	12.44	107	7555	3.421 ng/ul	94
34) 2-Methylnaphthalene	12.81	142	20936	3.959 ng/ul	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
35) 1-Methylnaphthalene	13.03	142	20158	3.931	ng/ul	93
37) 1,2,4,5-Tetrachlorobenzene	13.18	216	16236	3.961	ng/ul	96
38) Hexachlorocyclopentadiene	13.15	237	6811	2.834	ng/ul	91
39) 2,4,6-Trichlorophenol	13.41	196	7911	3.234	ng/ul	90
40) 2,4,5-Trichlorophenol	13.50	196	8579	3.301	ng/ul	85
41) 1,1'-Biphenyl	13.81	154	30210	3.906	ng/ul	99
42) 2-Chloronaphthalene	13.85	162	23826	3.820	ng/ul	95
43) 2-Nitroaniline	14.06	65	4610	3.136	ng/ul	93
45) Dimethylphthalate	14.41	163	34378	3.897	ng/ul#	98
46) 2,6-Dinitrotoluene	14.54	165	6195	3.498	ng/ul	91
48) Acenaphthylene	14.69	152	40041	4.056	ng/ul	97
49) 3-Nitroaniline	14.87	138	4946	3.361	ng/ul#	81
50) Acenaphthene	15.03	153	26651	3.942	ng/ul	99
51) 2,4-Dinitrophenol	15.09	184	2085	1.845	ng/ul#	81
53) 4-Nitrophenol	15.17	109	4470m	3.800	ng/ul	
54) Dibenzofuran	15.37	168	42070	4.093	ng/ul	97
55) 2,4-Dinitrotoluene	15.32	165	10908	4.125	ng/ul#	84
56) 2,3,4,6-Tetrachlorophenol	15.59	232	8898	3.215	ng/ul	97
57) Diethylphthalate	15.76	149	33105	3.575	ng/ul	98
59) Fluorene	16.01	166	35112	3.937	ng/ul	98
60) 4-Chlorophenyl-phenylether	16.00	204	21949	4.121	ng/ul	94
61) 4-Nitroaniline	16.04	138	5769	3.300	ng/ul	90
64) 4,6-Dinitro-2-methylphenol	16.09	198	7216	3.396	ng/ul#	65
65) N-Nitrosodiphenylamine	16.21	169	30070	3.825	ng/ul	98
66) 4-Bromophenyl-phenylether	16.90	248	15671	3.913	ng/ul#	84
67) Hexachlorobenzene	17.02	284	18468	4.055	ng/ul	92
68) Atrazine	17.15	200	13351	3.559	ng/ul	94
69) Pentachlorophenol	17.37	266	5835m	2.454	ng/ul	
70) Phenanthrene	17.77	178	62273	4.021	ng/ul	96
72) Anthracene	17.86	178	63201	3.920	ng/ul	97
73) 1,2,3,4-Tetrachlorobenzene	13.78	216	15350	3.533	ng/uL	96
74) Pentachlorobenzene	15.29	250	24788	3.840	ng/uL	93
75) Carbazole	18.12	167	51747	3.919	ng/ul	98
76) Di-n-butylphthalate	18.65	149	53673	3.661	ng/ul	98
77) Fluoranthene	19.76	202	80979	4.092	ng/ul#	94
80) Pyrene	20.13	202	89836	4.097	ng/ul#	93
81) Butylbenzylphthalate	20.99	149	22104	3.392	ng/ul#	96
82) 3,3'-Dichlorobenzidine	21.94	252	23960	3.100	ng/ul#	93
83) Benzo(a)anthracene	22.04	228	85980	3.832	ng/ul	99
84) Bis(2-ethylhexyl)phthalate	21.90	149	31373	3.291	ng/ul#	91
85) Chrysene	22.11	228	81323	3.962	ng/ul	98
87) Di-n-octyl phthalate	23.21	149	53200	3.175	ng/ul	100
88) Benzo(b)fluoranthene	24.45	252	84100	3.554	ng/ul#	97
89) Benzo(k)fluoranthene	24.52	252	90586	3.926	ng/ul#	95
91) Benzo(a)pyrene	25.41	252	84029	3.690	ng/ul#	96
92) Indeno(1,2,3-cd)pyrene	29.60	276	100835m	3.649	ng/ul	
93) Dibenzo(a,h)anthracene	29.67	278	84745	3.575	ng/ul	97
94) Benzo(g,h,i)perylene	30.84	276	84352m	3.682	ng/ul	

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

