

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010116\
 Data File : BM003872.D
 Acq On : 30 Dec 2015 10:23
 Operator : SJ/UM
 Sample : SSTD00537
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD00537

Quant Time: Dec 30 13:13:09 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM010116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 30 13:02:01 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	60561	20.00	ng/ul	0.00
18) Naphthalene-d8	10.46	136	260494	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.33	164	148163	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	320292	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	295614	20.00	ng/ul	0.00
86) Perylene-d12	23.51	264	245284	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.13	96	3259	2.90	ng/uL	0.00
5) Phenol-d5	0.00	99	0d	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	0.00	67	0d	0.00	ng/ul	
9) 2-Chlorophenol-d4	7.20	132	19040	4.57	ng/ul	0.00
13) 4-Methylphenol-d8	0.00	113	0d	0.00	ng/ul	
19) Nitrobenzene-d5	8.83	128	8763	4.64	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	9647	4.54	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.09	165	17862	4.60	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
44) Dimethylphthalate-d6	13.73	166	57011	4.66	ng/ul	0.00
47) Acenaphthylene-d8	14.02	160	67597	4.85	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
58) Fluorene-d10	15.32	176	49702	4.91	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul	
71) Anthracene-d10	17.17	188	74397	5.11	ng/ul	0.00
79) Pyrene-d10	19.47	212	77033	5.97	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	59597	4.89	ng/ul	0.00

Target Compounds

					Ovalue	
2) 1,4-Dioxane	3.17	88	3233	2.45	ng/uL#	80
10) 2-Chlorophenol	7.24	128	20200	4.67	ng/ul	97
15) N-Nitroso-di-n-propylamine	8.47	70	14975	4.48	ng/ul	94
17) Hexachloroethane	8.74	117	7653	4.65	ng/ul	97
20) Nitrobenzene	8.87	77	21516	4.93	ng/ul	98
21) Isophorone	9.39	82	40094	4.61	ng/ul	98
23) 2-Nitrophenol	9.58	139	10641	4.51	ng/ul	98
24) 2,4-Dimethylphenol	9.65	107	22416	4.64	ng/ul	99
25) Bis(2-Chloroethoxy)methane	9.87	93	25553	4.87	ng/ul	99
27) 2,4-Dichlorophenol	10.12	162	18288	4.53	ng/ul	99
28) Naphthalene	10.51	128	66630	4.96	ng/ul	100
31) Hexachlorobutadiene	10.79	225	11678	4.70	ng/ul	99
33) 4-Chloro-3-methylphenol	11.76	107	20904	4.42	ng/ul	98
34) 2-Methylnaphthalene	12.13	142	47195	4.71	ng/ul	99
35) 1-Methylnaphthalene	12.35	142	44840	4.71	ng/ul	97
37) 1,2,4,5-Tetrachlorobenzene	12.50	216	22524	5.02	ng/ul	97
39) 2,4,6-Trichlorophenol	12.75	196	14205	4.69	ng/ul	99
40) 2,4,5-Trichlorophenol	12.83	196	15310	4.57	ng/ul	98
41) 1,1'-Biphenyl	13.15	154	60993	5.07	ng/ul	99
42) 2-Chloronaphthalene	13.19	162	44726	4.95	ng/ul	99
43) 2-Nitroaniline	13.40	65	11379	4.55	ng/ul	92
45) Dimethylphthalate	13.78	163	58132	4.64	ng/ul	99
46) 2,6-Dinitrotoluene	13.91	165	10202	4.18	ng/ul	89

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48) Acenaphthylene	14.05	152	75117	4.93	ng/ul	99
50) Acenaphthene	14.39	153	50630	4.97	ng/ul	98
54) Dibenzofuran	14.73	168	71405	4.91	ng/ul	96
55) 2,4-Dinitrotoluene	14.70	165	14950	4.03	ng/ul#	91
56) 2,3,4,6-Tetrachlorophenol	14.96	232	12395	4.30	ng/ul#	98
57) Diethylphthalate	15.15	149	59273	4.47	ng/ul	100
59) Fluorene	15.37	166	58093	4.98	ng/ul	98
60) 4-Chlorophenyl-phenylether	15.37	204	27939	4.94	ng/ul	95
65) N-Nitrosodiphenylamine	15.59	169	49305	5.26	ng/ul	99
66) 4-Bromophenyl-phenylether	16.27	248	16234	5.09	ng/ul	94
67) Hexachlorobenzene	16.39	284	17628	4.97	ng/ul	98
70) Phenanthrene	17.12	178	90627	5.12	ng/ul	99
72) Anthracene	17.20	178	92847	5.14	ng/ul	99
73) 1,2,3,4-Tetrachlorobenzene	13.12	216	24748	5.64	ng/ul	97
74) Pentachlorobenzene	14.65	250	22260	5.38	ng/ul	99
76) Di-n-butylphthalate	18.04	149	99946	4.66	ng/ul	100
80) Pyrene	19.50	202	102566	6.03	ng/ul	98
81) Butylbenzylphthalate	20.41	149	38703	4.81	ng/ul	98
83) Benzo(a)anthracene	21.26	228	86728	4.87	ng/ul	100
84) Bis(2-ethylhexyl)phthalate	21.19	149	49167	4.11	ng/ul	99
85) Chrysene	21.31	228	82672	4.81	ng/ul	100
88) Benzo(b)fluoranthene	22.84	252	75505	5.11	ng/ul	99
89) Benzo(k)fluoranthene	22.88	252	68596	4.90	ng/ul	98
91) Benzo(a)pyrene	23.41	252	68540	4.85	ng/ul	99
92) Indeno(1,2,3-cd)pyrene	25.76	276	74306	4.88	ng/ul	98
93) Dibenzo(a,h)anthracene	25.76	278	62503	4.89	ng/ul	98
94) Benzo(g,h,i)perylene	26.44	276	61966	4.96	ng/ul	99

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

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