

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102218\
 Data File : BM017283.D
 Acq On : 22 Oct 2018 17:40
 Operator : MJ/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Oct 23 07:21:05 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM101718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 17 14:23:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	350414	20.00	ng	-0.02
21) Naphthalene-d8	10.43	136	1637682	20.00	ng	-0.02
39) Acenaphthene-d10	14.30	164	987842	20.00	ng	-0.01
64) Phenanthrene-d10	17.06	188	2227358	20.00	ng	-0.01
76) Chrysene-d12	21.25	240	2050527	20.00	ng	-0.01
87) Perylene-d12	23.46	264	1705313	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.27	112	1770533	85.46	ng	-0.01
7) Phenol-d6	6.85	99	2539650	90.00	ng	-0.01
23) Nitrobenzene-d5	8.82	82	2511497	89.69	ng	-0.02
42) 2,4,6-Tribromophenol	15.80	330	1210267	90.59	ng	-0.01
45) 2-Fluorobiphenyl	12.92	172	6179796	83.24	ng	-0.01
79) Terphenyl-d14	19.69	244	8980365	98.71	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.26	88	387503	36.628	ng	88
3) Pyridine	3.64	79	1009275	41.461	ng	87
4) n-Nitrosodimethylamine	3.56	42	386236	39.524	ng	# 73
6) Aniline	7.00	93	1543761	43.740	ng	98
8) 2-Chlorophenol	7.23	128	1067561	44.540	ng	97
9) Benzaldehyde	6.82	77	729888	40.568	ng	97
10) Phenol	6.87	94	1323145	43.570	ng	98
11) bis(2-Chloroethyl)ether	7.10	93	1037095	42.840	ng	97
12) 1,3-Dichlorobenzene	7.55	146	1167669	41.791	ng	97
13) 1,4-Dichlorobenzene	7.70	146	1196508	41.915	ng	98
14) 1,2-Dichlorobenzene	8.01	146	1162983	42.284	ng	100
15) Benzyl Alcohol	7.90	79	1012093	49.071	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.19	45	1374980	40.654	ng	99
17) 2-Methylphenol	8.10	107	897290	46.033	ng	99
18) Hexachloroethane	8.72	117	423012	43.300	ng	98
19) n-Nitroso-di-n-propylamine	8.47	70	892282	48.024	ng	97
20) 3+4-Methylphenols	8.44	107	1277614	47.922	ng	98
22) Acetophenone	8.48	105	1708979	41.167	ng	# 99
24) Nitrobenzene	8.86	77	1263211	42.648	ng	95
25) Isophorone	9.39	82	2091823	45.240	ng	99
26) 2-Nitrophenol	9.56	139	667969	48.164	ng	98
27) 2,4-Dimethylphenol	9.63	122	933424	45.654	ng	98
28) bis(2-Chloroethoxy)methane	9.86	93	1407128	43.237	ng	99
29) 2,4-Dichlorophenol	10.09	162	1120852	47.216	ng	99
30) 1,2,4-Trichlorobenzene	10.30	180	1206268	42.359	ng	99
31) Naphthalene	10.49	128	3347987	41.716	ng	99
32) Benzoic acid	9.80	122	743261	48.721	ng	92
33) 4-Chloroaniline	10.60	127	1558112	44.952	ng	98
34) Hexachlorobutadiene	10.77	225	744533	41.262	ng	98
35) Caprolactam	11.42	113	397757	45.976	ng	92
36) 4-Chloro-3-methylphenol	11.74	107	1256770	46.967	ng	98
37) 2-Methylnaphthalene	12.10	142	2453703	44.678	ng	99
38) 1-Methylnaphthalene	12.33	142	2366136	44.266	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.47	216	1464782	42.105	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.45	237	673873	40.069	ng	99
43) 2,4,6-Trichlorophenol	12.73	196	966281	47.848	ng	99
44) 2,4,5-Trichlorophenol	12.80	196	1014457	46.845	ng	99
46) 1,1'-Biphenyl	13.13	154	3694549	41.561	ng	99
47) 2-Chloronaphthalene	13.17	162	2730680	41.755	ng	98
48) 2-Nitroaniline	13.39	65	817809	44.706	ng	95
49) Acenaphthylene	14.02	152	4148389	43.762	ng	99
50) Dimethylphthalate	13.77	163	3311530	43.451	ng	99
51) 2,6-Dinitrotoluene	13.89	165	761714	45.410	ng	99
52) Acenaphthene	14.37	154	2516581	42.282	ng	98
53) 3-Nitroaniline	14.22	138	814704	44.845	ng	93
54) 2,4-Dinitrophenol	14.43	184	251988	31.200	ng	98
55) Dibenzofuran	14.71	168	4058910	41.009	ng	99
56) 4-Nitrophenol	14.54	139	616719	39.768	ng	100
57) 2,4-Dinitrotoluene	14.69	165	1043916	46.304	ng	# 99
58) Fluorene	15.36	166	3092340	42.210	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.93	232	908069	46.002	ng	100
60) Diethylphthalate	15.14	149	3399153	44.974	ng	99
61) 4-Chlorophenyl-phenylether	15.36	204	1754270	42.003	ng	97
62) 4-Nitroaniline	15.39	138	843646	42.244	ng	99
63) Azobenzene	15.65	77	3146020	42.824	ng	99
65) 4,6-Dinitro-2-methylphenol	15.44	198	425229	32.395	ng	97
66) n-Nitrosodiphenylamine	15.57	169	2976270	44.166	ng	99
67) 4-Bromophenyl-phenylether	16.25	248	1120371	45.811	ng	96
68) Hexachlorobenzene	16.36	284	1234781	43.465	ng	98
69) Atrazine	16.54	200	1096094	45.446	ng	97
70) Pentachlorophenol	16.70	266	799091	41.061	ng	99
71) Phenanthrene	17.10	178	4945661	41.154	ng	100
72) Anthracene	17.19	178	4940923	43.271	ng	100
73) Carbazole	17.46	167	4960209	42.525	ng	99
74) Di-n-butylphthalate	18.03	149	5831247	47.399	ng	99
75) Fluoranthene	19.12	202	5603502	41.438	ng	99
77) Benzidine	19.32	184	2926793	56.291	ng	99
78) Pyrene	19.48	202	5753579	50.204	ng	99
80) Butylbenzylphthalate	20.40	149	2554052	56.985	ng	97
81) Benzo(a)anthracene	21.23	228	4992094	43.262	ng	100
82) 3,3'-Dichlorobenzidine	21.17	252	1976047	45.947	ng	99
83) Chrysene	21.29	228	4747891	41.570	ng	100
84) Bis(2-ethylhexyl)phthalate	21.17	149	3537708	52.464	ng	98
85) Di-n-octyl phthalate	22.04	149	5427940	48.016	ng	100
86) Indeno(1,2,3-cd)pyrene	25.67	276	4052947	31.726	ng	99
88) Benzo(b)fluoranthene	22.80	252	4114489	44.130	ng	99
89) Benzo(k)fluoranthene	22.84	252	4025785	43.031	ng	99
90) Benzo(a)pyrene	23.36	252	3730727	42.143	ng	99
91) Dibenzo(a,h)anthracene	25.69	278	3416462	38.867	ng	98
92) Benzo(g,h,i)perylene	26.34	276	3361474	38.584	ng	99

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

