

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM101918\
 Data File : BM017254.D
 Acq On : 20 Oct 2018 07:13
 Operator : MJ/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Oct 22 02:36:33 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	311310	20.00	ng	-0.01
21) Naphthalene-d8	10.44	136	1461378	20.00	ng	-0.01
39) Acenaphthene-d10	14.30	164	879823	20.00	ng	-0.01
64) Phenanthrene-d10	17.06	188	1904945	20.00	ng	-0.01
76) Chrysene-d12	21.25	240	1679042	20.00	ng	-0.01
87) Perylene-d12	23.46	264	1511848	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.28	112	1454354	79.01	ng	-0.01
7) Phenol-d6	6.85	99	2069623	82.56	ng	-0.01
23) Nitrobenzene-d5	8.82	82	2117627	84.75	ng	-0.01
42) 2,4,6-Tribromophenol	15.80	330	991241	83.31	ng	-0.01
45) 2-Fluorobiphenyl	12.92	172	5222157	78.97	ng	-0.01
79) Terphenyl-d14	19.70	244	7003817	94.02	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.26	88	326275	34.714	ng	# 86
3) Pyridine	3.64	79	848934	39.255	ng	# 85
4) n-Nitrosodimethylamine	3.56	42	330492	38.068	ng	# 69
6) Aniline	7.00	93	1286886	41.042	ng	97
8) 2-Chlorophenol	7.23	128	889124	41.755	ng	97
9) Benzaldehyde	6.82	77	626516	39.197	ng	96
10) Phenol	6.87	94	1098324	40.710	ng	98
11) bis(2-Chloroethyl)ether	7.10	93	876610	40.759	ng	99
12) 1,3-Dichlorobenzene	7.55	146	982513	39.581	ng	98
13) 1,4-Dichlorobenzene	7.70	146	998564	39.374	ng	96
14) 1,2-Dichlorobenzene	8.01	146	977007	39.985	ng	99
15) Benzyl Alcohol	7.91	79	836857	45.671	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.19	45	1196420	39.818	ng	99
17) 2-Methylphenol	8.10	107	743876	42.956	ng	98
18) Hexachloroethane	8.73	117	359747	41.450	ng	98
19) n-Nitroso-di-n-propylamine	8.47	70	758213	45.934	ng	98
20) 3+4-Methylphenols	8.43	107	1039876	43.905	ng	98
22) Acetophenone	8.49	105	1450258	39.149	ng	# 99
24) Nitrobenzene	8.86	77	1075970	40.709	ng	94
25) Isophorone	9.39	82	1745876	42.313	ng	98
26) 2-Nitrophenol	9.56	139	536060	43.890	ng	99
27) 2,4-Dimethylphenol	9.63	122	762805	41.810	ng	99
28) bis(2-Chloroethoxy)methane	9.87	93	1178852	40.593	ng	99
29) 2,4-Dichlorophenol	10.09	162	904847	42.715	ng	99
30) 1,2,4-Trichlorobenzene	10.30	180	999044	39.314	ng	99
31) Naphthalene	10.49	128	2795847	39.039	ng	99
32) Benzoic acid	9.79	122	652980	47.967	ng	94
33) 4-Chloroaniline	10.60	127	1291368	41.751	ng	97
34) Hexachlorobutadiene	10.77	225	636229	39.514	ng	99
35) Caprolactam	11.41	113	316604	41.011	ng	95
36) 4-Chloro-3-methylphenol	11.74	107	1022451	42.820	ng	98
37) 2-Methylnaphthalene	12.11	142	2031736	41.458	ng	98
38) 1-Methylnaphthalene	12.33	142	1973450	41.374	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.48	216	1236007	39.891	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.46	237	683304	45.618	ng	99
43) 2,4,6-Trichlorophenol	12.73	196	785808	43.689	ng	99
44) 2,4,5-Trichlorophenol	12.80	196	824528	42.749	ng	98
46) 1,1'-Biphenyl	13.13	154	3119589	39.401	ng	100
47) 2-Chloronaphthalene	13.17	162	2311997	39.693	ng	99
48) 2-Nitroaniline	13.39	65	682308	42.238	ng	99
49) Acenaphthylene	14.03	152	3447690	40.836	ng	100
50) Dimethylphthalate	13.77	163	2755446	40.594	ng	99
51) 2,6-Dinitrotoluene	13.89	165	620630	41.542	ng	97
52) Acenaphthene	14.37	154	2103153	39.674	ng	98
53) 3-Nitroaniline	14.23	138	658369	40.689	ng	98
54) 2,4-Dinitrophenol	14.43	184	331342	42.144	ng	96
55) Dibenzofuran	14.71	168	3397700	38.543	ng	99
56) 4-Nitrophenol	14.53	139	527372	38.370	ng	95
57) 2,4-Dinitrotoluene	14.69	165	846771	42.170	ng	99
58) Fluorene	15.36	166	2580607	39.550	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.94	232	755297	42.960	ng	99
60) Diethylphthalate	15.15	149	2819081	41.878	ng	100
61) 4-Chlorophenyl-phenylether	15.36	204	1482791	39.861	ng	96
62) 4-Nitroaniline	15.39	138	680529	38.709	ng	96
63) Azobenzene	15.65	77	2642413	40.385	ng	98
65) 4,6-Dinitro-2-methylphenol	15.45	198	536884	44.489	ng	98
66) n-Nitrosodiphenylamine	15.58	169	2485149	43.120	ng	99
67) 4-Bromophenyl-phenylether	16.26	248	924530	44.201	ng	96
68) Hexachlorobenzene	16.36	284	1014581	41.758	ng	99
69) Atrazine	16.54	200	884802	42.895	ng	98
70) Pentachlorophenol	16.71	266	712520	42.810	ng	99
71) Phenanthrene	17.10	178	4058616	39.489	ng	100
72) Anthracene	17.19	178	4031033	41.277	ng	100
73) Carbazole	17.47	167	3939454	39.490	ng	100
74) Di-n-butylphthalate	18.03	149	4518795	42.948	ng	100
75) Fluoranthene	19.12	202	4332834	37.464	ng	99
77) Benzidine	19.32	184	2152017	50.547	ng	99
78) Pyrene	19.49	202	4490126	47.848	ng	100
80) Butylbenzylphthalate	20.40	149	1751914	48.901	ng	97
81) Benzo(a)anthracene	21.23	228	3797980	40.196	ng	100
82) 3,3'-Dichlorobenzidine	21.17	252	1459849	41.455	ng	99
83) Chrysene	21.29	228	3676841	39.315	ng	99
84) Bis(2-ethylhexyl)phthalate	21.17	149	2371063	43.906	ng	99
85) Di-n-octyl phthalate	22.05	149	3606365	40.011	ng	100
86) Indeno(1,2,3-cd)pyrene	25.67	276	3565947	34.090	ng	99
88) Benzo(b)fluoranthene	22.80	252	3252596	39.350	ng	100
89) Benzo(k)fluoranthene	22.84	252	3407251	41.080	ng	99
90) Benzo(a)pyrene	23.37	252	3108707	39.611	ng	99
91) Dibenzo(a,h)anthracene	25.69	278	3014170	38.678	ng	99
92) Benzo(g,h,i)perylene	26.35	276	2959447	38.316	ng	99

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

