

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021219\
 Data File : BM018770.D
 Acq On : 12 Feb 2019 13:28
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Feb 13 01:31:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 11 16:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	306443	20.00	ng	0.00
21) Naphthalene-d8	10.50	136	1289598	20.00	ng	0.00
39) Acenaphthene-d10	14.36	164	767734	20.00	ng	0.00
64) Phenanthrene-d10	17.12	188	1761705	20.00	ng	0.00
76) Chrysene-d12	21.32	240	1945389	20.00	ng	0.00
87) Perylene-d12	23.57	264	1967037	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.30	112	1515299	81.02	ng	0.00
7) Phenol-d6	6.89	99	2165998	82.21	ng	0.00
23) Nitrobenzene-d5	8.89	82	2270692	81.42	ng	0.00
42) 2,4,6-Tribromophenol	15.86	330	916921	82.85	ng	0.00
45) 2-Fluorobiphenyl	12.98	172	4532708	78.37	ng	0.00
79) Terphenyl-d14	19.75	244	7965867	84.47	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.25	88	316863	38.898	ng	100
3) Pyridine	3.65	79	956986	43.069	ng	98
4) n-Nitrosodimethylamine	3.58	42	234202	41.432	ng	97
6) Aniline	7.06	93	1315981	40.764	ng	99
8) 2-Chlorophenol	7.28	128	834900	40.659	ng	99
9) Benzaldehyde	6.87	77	591774	41.511	ng	100
10) Phenol	6.92	94	1135867	40.972	ng	99
11) bis(2-Chloroethyl)ether	7.15	93	871097	41.191	ng	98
12) 1,3-Dichlorobenzene	7.60	146	920689	39.874	ng	99
13) 1,4-Dichlorobenzene	7.75	146	942621	40.101	ng	100
14) 1,2-Dichlorobenzene	8.06	146	900554	39.884	ng	98
15) Benzyl Alcohol	7.96	79	875242	41.806	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.25	45	821985	40.159	ng	99
17) 2-Methylphenol	8.16	107	724191	41.042	ng	98
18) Hexachloroethane	8.78	117	346991	40.438	ng	99
19) n-Nitroso-di-n-propylamine	8.53	70	846298	41.484	ng	100
20) 3+4-Methylphenols	8.49	107	1019146	41.828	ng	98
22) Acetophenone	8.55	105	1437432	40.589	ng	100
24) Nitrobenzene	8.93	77	1164351	40.219	ng	99
25) Isophorone	9.45	82	1848014	41.527	ng	100
26) 2-Nitrophenol	9.63	139	486602	42.207	ng	98
27) 2,4-Dimethylphenol	9.69	122	685228	40.694	ng	98
28) bis(2-Chloroethoxy)methane	9.93	93	1170947	40.690	ng	100
29) 2,4-Dichlorophenol	10.16	162	803560	41.560	ng	98
30) 1,2,4-Trichlorobenzene	10.36	180	889747	39.847	ng	99
31) Naphthalene	10.56	128	2502568	39.790	ng	99
32) Benzoic acid	9.85	122	571564	38.912	ng	98
33) 4-Chloroaniline	10.67	127	1159864	41.254	ng	100
34) Hexachlorobutadiene	10.82	225	514802	39.720	ng	99
35) Caprolactam	11.50	113	319324	40.468	ng	93
36) 4-Chloro-3-methylphenol	11.80	107	972793	41.730	ng	99
37) 2-Methylnaphthalene	12.17	142	1780166	40.201	ng	100
38) 1-Methylnaphthalene	12.39	142	1730491	39.963	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.54	216	962592	39.195	ng	100

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021219\
 Data File : BM018770.D
 Acq On : 12 Feb 2019 13:28
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Feb 13 01:31:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 11 16:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.50	237	493523	39.305	ng	98
43) 2,4,6-Trichlorophenol	12.79	196	665925	40.960	ng	99
44) 2,4,5-Trichlorophenol	12.86	196	703289	41.271	ng	98
46) 1,1'-Biphenyl	13.19	154	2589232	39.180	ng	100
47) 2-Chloronaphthalene	13.24	162	2006142	39.303	ng	100
48) 2-Nitroaniline	13.46	65	715822	42.225	ng	100
49) Acenaphthylene	14.09	152	3051755	40.153	ng	99
50) Dimethylphthalate	13.83	163	2554057	39.918	ng	99
51) 2,6-Dinitrotoluene	13.96	165	570793	41.183	ng	99
52) Acenaphthene	14.43	154	1824959	39.159	ng	100
53) 3-Nitroaniline	14.29	138	631681	40.336	ng	99
54) 2,4-Dinitrophenol	14.50	184	296818	37.081	ng	98
55) Dibenzofuran	14.77	168	3015773	39.370	ng	99
56) 4-Nitrophenol	14.60	139	499563	39.961	ng	97
57) 2,4-Dinitrotoluene	14.74	165	805881	42.200	ng	99
58) Fluorene	15.42	166	2321738	39.619	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.99	232	607411	40.326	ng	99
60) Diethylphthalate	15.20	149	2635557	39.932	ng	99
61) 4-Chlorophenyl-phenylether	15.42	204	1296999	39.474	ng	97
62) 4-Nitroaniline	15.46	138	621912	40.508	ng	99
63) Azobenzene	15.70	77	2910960	41.134	ng	99
65) 4,6-Dinitro-2-methylphenol	15.51	198	482896	39.041	ng	94
66) n-Nitrosodiphenylamine	15.63	169	2237651	40.203	ng	100
67) 4-Bromophenyl-phenylether	16.31	248	786365	39.878	ng	100
68) Hexachlorobenzene	16.41	284	905665	39.513	ng	99
69) Atrazine	16.59	200	718706	40.053	ng	100
70) Pentachlorophenol	16.76	266	615162	41.683	ng	99
71) Phenanthrene	17.16	178	3765585	39.773	ng	100
72) Anthracene	17.25	178	3709873	40.258	ng	100
73) Carbazole	17.53	167	3906518	40.576	ng	99
74) Di-n-butylphthalate	18.08	149	4643049	41.933	ng	100
75) Fluoranthene	19.18	202	4420028	40.416	ng	99
77) Benzidine	19.38	184	2137960	48.797	ng	100
78) Pyrene	19.54	202	4648542	41.143	ng	99
80) Butylbenzylphthalate	20.44	149	2254160	43.777	ng	98
81) Benzo(a)anthracene	21.30	228	4609184	40.645	ng	99
82) 3,3'-Dichlorobenzidine	21.24	252	1879570	41.938	ng	99
83) Chrysene	21.35	228	4446667	40.910	ng	100
84) Bis(2-ethylhexyl)phthalate	21.22	149	3311699	43.942	ng	100
85) Di-n-octyl phthalate	22.10	149	5553548	43.844	ng	100
86) Indeno(1,2,3-cd)pyrene	25.87	276	4733779	37.722	ng	100
88) Benzo(b)fluoranthene	22.89	252	4610847	40.970	ng	99
89) Benzo(k)fluoranthene	22.94	252	4546571	40.579	ng	99
90) Benzo(a)pyrene	23.48	252	4285310	40.195	ng	100
91) Dibenzo(a,h)anthracene	25.88	278	4068216	38.515	ng	99
92) Benzo(g,h,i)perylene	26.57	276	3675476	37.377	ng	99

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

