

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\
 Data File : BP002620.D
 Acq On : 02 Jul 2020 08:47
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Quant Time: Jul 02 10:42:21 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 29 16:51:57 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	201910	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	812417	20.00	ng	0.00
39) Acenaphthene-d10	14.22	164	501244	20.00	ng	0.00
64) Phenanthrene-d10	16.98	188	1055168	20.00	ng	0.00
76) Chrysene-d12	21.09	240	932033	20.00	ng	0.00
86) Perylene-d12	23.29	264	988604	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.19	112	952082	79.84	ng	0.00
7) Phenol-d6	6.77	99	1363066	81.89	ng	0.00
23) Nitrobenzene-d5	8.72	82	1227766	77.87	ng	0.00
42) 2,4,6-Tribromophenol	15.73	330	433515	71.11	ng	0.00
45) 2-Fluorobiphenyl	12.84	172	2511599	73.75	ng	0.00
79) Terphenyl-d14	19.57	244	3265944	75.64	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.15	88	194383	38.031	ng	99
3) Pyridine	3.52	79	607103	40.077	ng	99
4) n-Nitrosodimethylamine	3.44	42	238297	37.059	ng	90
6) Aniline	6.90	93	867294	41.477	ng	99
8) 2-Chlorophenol	7.15	128	529510	39.972	ng	99
9) Benzaldehyde	6.72	77	354372	41.080	ng	99
10) Phenol	6.79	94	695525	41.437	ng	95
11) bis(2-Chloroethyl)ether	7.00	93	569450	41.334	ng	98
12) 1,3-Dichlorobenzene	7.46	146	592537	38.763	ng	99
13) 1,4-Dichlorobenzene	7.60	146	591599	37.988	ng	100
14) 1,2-Dichlorobenzene	7.92	146	574239	38.307	ng	98
15) Benzyl Alcohol	7.82	79	498789	39.881	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.11	45	824354	42.947	ng	99
17) 2-Methylphenol	8.03	107	501553	39.921	ng	99
18) Hexachloroethane	8.64	117	216826	38.577	ng	99
19) n-Nitroso-di-n-propylamine	8.38	70	435334	39.863	ng	98
20) 3+4-Methylphenols	8.36	107	682818	40.312	ng	98
22) Acetophenone	8.39	105	796930	38.756	ng	98
24) Nitrobenzene	8.76	77	657809	39.415	ng	100
25) Isophorone	9.29	82	1268661	40.145	ng	99
26) 2-Nitrophenol	9.47	139	306885	40.319	ng	96
27) 2,4-Dimethylphenol	9.55	122	462625	40.111	ng	97
28) bis(2-Chloroethoxy)methane	9.77	93	746565	41.126	ng	100
29) 2,4-Dichlorophenol	10.01	162	520812	39.182	ng	98
30) 1,2,4-Trichlorobenzene	10.22	180	564156	37.700	ng	98
31) Naphthalene	10.39	128	1660804	38.617	ng	100
32) Benzoic acid	9.71	122	279798	33.989	ng	99
33) 4-Chloroaniline	10.50	127	749367	39.834	ng	100
34) Hexachlorobutadiene	10.69	225	336398	36.727	ng	98
35) Caprolactam	11.28	113	180484	40.393	ng	93
36) 4-Chloro-3-methylphenol	11.66	107	562337	38.599	ng	100
37) 2-Methylnaphthalene	12.02	142	1193507	38.061	ng	99
38) 1-Methylnaphthalene	12.24	142	1127370	38.173	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.40	216	603606	38.566	ng	100

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\
 Data File : BP002620.D
 Acq On : 02 Jul 2020 08:47
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Quant Time: Jul 02 10:42:21 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 29 16:51:57 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.38	237	341880	43.934	ng	97
43) 2,4,6-Trichlorophenol	12.65	196	414631	39.152	ng	98
44) 2,4,5-Trichlorophenol	12.72	196	467896	38.380	ng	99
46) 1,1'-Biphenyl	13.05	154	1441464	38.018	ng	99
47) 2-Chloronaphthalene	13.09	162	1168537	37.861	ng	97
48) 2-Nitroaniline	13.29	65	400025	40.627	ng	98
49) Acenaphthylene	13.94	152	1863260	38.359	ng	99
50) Dimethylphthalate	13.69	163	1483500	37.682	ng	99
51) 2,6-Dinitrotoluene	13.80	165	330652	38.889	ng	96
52) Acenaphthene	14.29	154	1105293	38.568	ng	99
53) 3-Nitroaniline	14.13	138	377475	40.773	ng	99
54) 2,4-Dinitrophenol	14.35	184	174300	36.424	ng	96
55) Dibenzofuran	14.63	168	1760440	37.905	ng	99
56) 4-Nitrophenol	14.47	139	267042	37.863	ng	92
57) 2,4-Dinitrotoluene	14.60	165	451215	39.450	ng	96
58) Fluorene	15.28	166	1305617	36.912	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.86	232	370078	37.579	ng	99
60) Diethylphthalate	15.07	149	1445198	36.917	ng	100
61) 4-Chlorophenyl-phenylether	15.28	204	693412	36.556	ng	96
62) 4-Nitroaniline	15.30	138	371002	40.095	ng	91
63) Azobenzene	15.57	77	1486817	39.279	ng	97
65) 4,6-Dinitro-2-methylphenol	15.37	198	258305	38.597	ng	96
66) n-Nitrosodiphenylamine	15.50	169	1228147	39.871	ng	99
67) 4-Bromophenyl-phenylether	16.18	248	465468	39.255	ng	95
68) Hexachlorobenzene	16.30	284	481215	37.882	ng	97
69) Atrazine	16.46	200	383929	38.844	ng	99
70) Pentachlorophenol	16.65	266	249184	35.844	ng	98
71) Phenanthrene	17.03	178	2181966	38.434	ng	100
72) Anthracene	17.12	178	2184336	38.665	ng	98
73) Carbazole	17.39	167	2140153	39.198	ng	99
74) Di-n-butylphthalate	17.97	149	2530591	39.356	ng	99
75) Fluoranthene	19.01	202	2595106	37.764	ng	98
77) Benzidine	19.19	184	990700	34.488	ng	100
78) Pyrene	19.36	202	2658870	41.911	ng	99
80) Butylbenzylphthalate	20.26	149	1167044	42.137	ng	100
81) Benzo(a)anthracene	21.07	228	2403206	39.440	ng	99
82) 3,3'-Dichlorobenzidine	21.02	252	886448	39.737	ng	94
83) Chrysene	21.13	228	2259911	38.677	ng	99
84) Bis(2-ethylhexyl)phthalate	21.03	149	1582353	39.787	ng	99
85) Di-n-octyl phthalate	21.89	149	2801076	39.876	ng	99
87) Indeno(1,2,3-cd)pyrene	25.51	276	2852612	40.075	ng	98
88) Benzo(b)fluoranthene	22.63	252	2499208	39.618	ng	98
89) Benzo(k)fluoranthene	22.68	252	2346406	39.872	ng	98
90) Benzo(a)pyrene	23.20	252	2314774	39.987	ng	99
91) Dibenzo(a,h)anthracene	25.52	278	2322954	39.938	ng	99
92) Benzo(g,h,i)perylene	26.19	276	2341802	40.257	ng	98

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

