

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062119\
 Data File : BF115078.D
 Acq On : 21 Jun 2019 14:38
 Operator : HP/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICCC040

Quant Time: Jun 21 16:35:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 21 16:30:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.62	152	198977	20.00	ng	0.00
21) Naphthalene-d8	7.90	136	816147	20.00	ng	0.00
39) Acenaphthene-d10	9.65	164	399645	20.00	ng	0.00
64) Phenanthrene-d10	11.12	188	784869	20.00	ng	0.00
76) Chrysene-d12	13.75	240	581758	20.00	ng	0.00
87) Perylene-d12	15.11	264	675206	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.21	112	1067563	89.59	ng	0.00
7) Phenol-d6	6.25	99	1298213	90.31	ng	0.00
23) Nitrobenzene-d5	7.18	82	1096020	80.83	ng	0.00
42) 2,4,6-Tribromophenol	10.43	330	353603	82.62	ng	0.00
45) 2-Fluorobiphenyl	8.98	172	2007097	85.12	ng	0.00
79) Terphenyl-d14	12.71	244	2249942	72.74	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.20	88	249252	44.686	ng	100
3) Pyridine	2.85	79	698118	44.455	ng	100
4) n-Nitrosodimethylamine	2.80	42	272914	48.950	ng	100
6) Aniline	6.28	93	895975	46.543	ng	100
8) 2-Chlorophenol	6.40	128	603065	44.632	ng	100
9) Benzaldehyde	6.16	77	434341	52.559	ng	100
10) Phenol	6.27	94	757133	47.078	ng	100
11) bis(2-Chloroethyl)ether	6.36	93	554656	44.434	ng	100
12) 1,3-Dichlorobenzene	6.56	146	663633	42.049	ng	100
13) 1,4-Dichlorobenzene	6.64	146	671991	42.352	ng	100
14) 1,2-Dichlorobenzene	6.79	146	626218	41.921	ng	100
15) Benzyl Alcohol	6.77	79	481896	44.745	ng	100
16) 2,2'-oxybis(1-Chloropropan	6.91	45	804722	48.167	ng	100
17) 2-Methylphenol	6.88	107	489463	43.429	ng	100
18) Hexachloroethane	7.14	117	240346	41.945	ng	100
19) n-Nitroso-di-n-propylamine	7.05	70	408196	46.739	ng	100
20) 3+4-Methylphenols	7.04	107	571159	42.581	ng	100
22) Acetophenone	7.03	105	769945	41.675	ng	100
24) Nitrobenzene	7.20	77	604728	43.835	ng	100
25) Isophorone	7.45	82	1103217	43.442	ng	100
26) 2-Nitrophenol	7.52	139	300642	39.098	ng	100
27) 2,4-Dimethylphenol	7.57	122	488503	44.038	ng	100
28) bis(2-Chloroethoxy)methane	7.66	93	712431	43.726	ng	100
29) 2,4-Dichlorophenol	7.76	162	506721	43.241	ng	100
30) 1,2,4-Trichlorobenzene	7.84	180	563839	41.641	ng	100
31) Naphthalene	7.92	128	1679740	43.499	ng	100
32) Benzoic acid	7.68	122	296982	37.002	ng	100
33) 4-Chloroaniline	7.97	127	759480	42.608	ng	100
34) Hexachlorobutadiene	8.05	225	308415	40.978	ng	100
35) Caprolactam	8.35	113	165302	42.352	ng	100
36) 4-Chloro-3-methylphenol	8.45	107	512058	43.360	ng	100
37) 2-Methylnaphthalene	8.61	142	1127550	43.779	ng	100
38) 1-Methylnaphthalene	8.71	142	1083832	44.524	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.78	216	478727	40.546	ng	100

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062119\
 Data File : BF115078.D
 Acq On : 21 Jun 2019 14:38
 Operator : HP/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICCC040

Quant Time: Jun 21 16:35:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 21 16:30:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.78	237	279279	48.002	ng	100
43) 2,4,6-Trichlorophenol	8.89	196	366586	41.569	ng	100
44) 2,4,5-Trichlorophenol	8.92	196	367302	41.214	ng	100
46) 1,1'-Biphenyl	9.08	154	1316677	41.113	ng	100
47) 2-Chloronaphthalene	9.10	162	1095390	42.298	ng	100
48) 2-Nitroaniline	9.19	65	317378	41.463	ng	100
49) Acenaphthylene	9.51	152	1709390	42.931	ng	100
50) Dimethylphthalate	9.38	163	1240488	41.274	ng	100
51) 2,6-Dinitrotoluene	9.44	165	282393	41.421	ng	100
52) Acenaphthene	9.68	154	1068037	46.884	ng	100
53) 3-Nitroaniline	9.60	138	346390	41.469	ng	100
54) 2,4-Dinitrophenol	9.70	184	121423	36.719	ng	100
55) Dibenzofuran	9.85	168	1493640	43.504	ng	100
56) 4-Nitrophenol	9.75	139	282683	49.075	ng	100
57) 2,4-Dinitrotoluene	9.84	165	370249	42.668	ng	100
58) Fluorene	10.20	166	1046638	41.680	ng	100
59) 2,3,4,6-Tetrachlorophenol	9.97	232	315572	43.836	ng	100
60) Diethylphthalate	10.08	149	1230006	42.801	ng	100
61) 4-Chlorophenyl-phenylether	10.20	204	501142	40.431	ng	100
62) 4-Nitroaniline	10.21	138	346796	41.293	ng	100
63) Azobenzene	10.35	77	1157301	46.033	ng	100
65) 4,6-Dinitro-2-methylphenol	10.24	198	162698	36.985	ng	100
66) n-Nitrosodiphenylamine	10.31	169	1025298	43.702	ng	100
67) 4-Bromophenyl-phenylether	10.68	248	354495	42.066	ng	100
68) Hexachlorobenzene	10.74	284	382548	41.232	ng	100
69) Atrazine	10.84	200	329385	43.545	ng	100
70) Pentachlorophenol	10.93	266	250032	49.624	ng	100
71) Phenanthrene	11.15	178	1704265	42.056	ng	100
72) Anthracene	11.20	178	1746284	42.715	ng	100
73) Carbazole	11.35	167	1599581	44.827	ng	100
74) Di-n-butylphthalate	11.71	149	1853700	41.207	ng	100
75) Fluoranthene	12.32	202	1775062	42.837	ng	100
77) Benzidine	12.45	184	1070022	45.473	ng	100
78) Pyrene	12.55	202	1825713	34.449	ng	100
80) Butylbenzylphthalate	13.19	149	805440	32.538	ng	100
81) Benzo(a)anthracene	13.74	228	1717809	39.459	ng	100
82) 3,3'-Dichlorobenzidine	13.71	252	635232	38.060	ng	100
83) Chrysene	13.77	228	1569900	36.264	ng	100
84) Bis(2-ethylhexyl)phthalate	13.76	149	1055685	36.084	ng	100
85) Di-n-octyl phthalate	14.37	149	1791008	35.550	ng	100
86) Indeno(1,2,3-cd)pyrene	16.39	276	1900327	49.973	ng	100
88) Benzo(b)fluoranthene	14.74	252	1792457	39.934	ng	100
89) Benzo(k)fluoranthene	14.77	252	1454807	36.926	ng	100
90) Benzo(a)pyrene	15.06	252	1587198	40.980	ng	100
91) Dibenzo(a,h)anthracene	16.42	278	1530661	45.890	ng	100
92) Benzo(g,h,i)perylene	16.78	276	1563928	45.975	ng	100

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

