

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110719\
 Data File : BF117705.D
 Acq On : 7 Nov 2019 14:57
 Operator : JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Nov 07 15:37:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 17:27:10 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	386786	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	1447457	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	764387	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	1378691	20.00	ng	0.00
76) Chrysene-d12	14.13	240	964566	20.00	ng	0.00
87) Perylene-d12	15.64	264	935574	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.55	112	1657376	77.56	ng	0.00
7) Phenol-d6	6.56	99	2045678	77.64	ng	0.00
23) Nitrobenzene-d5	7.50	82	1896934	77.37	ng	0.00
42) 2,4,6-Tribromophenol	10.77	330	584174	78.12	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	3284060	84.77	ng	0.00
79) Terphenyl-d14	13.07	244	3525715	75.25	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.76	88	414007	38.886	ng	99
3) Pyridine	3.52	79	1148740	39.226	ng	99
4) n-Nitrosodimethylamine	3.47	42	498280	38.610	ng	99
6) Aniline	6.60	93	1441068	38.973	ng	100
8) 2-Chlorophenol	6.72	128	934636	39.320	ng	100
9) Benzaldehyde	6.49	77	737344	40.222	ng	99
10) Phenol	6.57	94	1120643	39.473	ng	100
11) bis(2-Chloroethyl)ether	6.67	93	910623	38.844	ng	100
12) 1,3-Dichlorobenzene	6.88	146	1080466	39.213	ng	99
13) 1,4-Dichlorobenzene	6.96	146	1092577	39.612	ng	99
14) 1,2-Dichlorobenzene	7.11	146	1014173	39.935	ng	99
15) Benzyl Alcohol	7.07	79	821913	39.125	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.21	45	1427782	38.519	ng	100
17) 2-Methylphenol	7.18	107	786184	38.510	ng	100
18) Hexachloroethane	7.46	117	390665	38.319	ng	99
19) n-Nitroso-di-n-propylamine	7.35	70	658025	37.984	ng	99
20) 3+4-Methylphenols	7.33	107	960857	39.526	ng	99
22) Acetophenone	7.35	105	1283982	38.812	ng	99
24) Nitrobenzene	7.52	77	1005612	39.280	ng	100
25) Isophorone	7.76	82	1762206	37.580	ng	100
26) 2-Nitrophenol	7.83	139	458088	38.543	ng	99
27) 2,4-Dimethylphenol	7.87	122	761668	38.006	ng	98
28) bis(2-Chloroethoxy)methane	7.97	93	1149061	38.174	ng	100
29) 2,4-Dichlorophenol	8.07	162	794616	38.486	ng	99
30) 1,2,4-Trichlorobenzene	8.16	180	928137	39.268	ng	99
31) Naphthalene	8.25	128	2601102	39.315	ng	99
32) Benzoic acid	7.98	122	588155	37.075	ng	98
33) 4-Chloroaniline	8.29	127	1183595	38.501	ng	98
34) Hexachlorobutadiene	8.37	225	525071	38.793	ng	99
35) Caprolactam	8.66	113	253332	37.019	ng	96
36) 4-Chloro-3-methylphenol	8.76	107	801554	38.472	ng	99
37) 2-Methylnaphthalene	8.94	142	1778870	39.495	ng	99
38) 1-Methylnaphthalene	9.04	142	1677256	39.213	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.10	216	826622	39.529	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.10	237	422709	37.469	ng	99
43) 2,4,6-Trichlorophenol	9.21	196	585568	38.448	ng	97
44) 2,4,5-Trichlorophenol	9.25	196	597176	38.683	ng	94
46) 1,1'-Biphenyl	9.40	154	2111294	39.438	ng	99
47) 2-Chloronaphthalene	9.43	162	1737934	39.120	ng	99
48) 2-Nitroaniline	9.52	65	536677	38.548	ng	99
49) Acenaphthylene	9.85	152	2552259	39.448	ng	100
50) Dimethylphthalate	9.70	163	1979019	38.297	ng	100
51) 2,6-Dinitrotoluene	9.76	165	451896	39.053	ng	99
52) Acenaphthene	10.02	154	1624778	38.918	ng	100
53) 3-Nitroaniline	9.93	138	548181	39.268	ng	95
54) 2,4-Dinitrophenol	10.03	184	138975	35.033	ng	92
55) Dibenzofuran	10.19	168	2303012	39.422	ng	100
56) 4-Nitrophenol	10.07	139	400749	38.171	ng	99
57) 2,4-Dinitrotoluene	10.17	165	588163	39.624	ng	98
58) Fluorene	10.54	166	1694545	39.357	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.30	232	506264	39.450	ng	99
60) Diethylphthalate	10.40	149	1964712	38.660	ng	100
61) 4-Chlorophenyl-phenylether	10.53	204	876595	39.111	ng	99
62) 4-Nitroaniline	10.55	138	531441	38.678	ng	99
63) Azobenzene	10.69	77	1872445	38.901	ng	99
65) 4,6-Dinitro-2-methylphenol	10.57	198	200609	36.440	ng	97
66) n-Nitrosodiphenylamine	10.64	169	1591592	39.087	ng	100
67) 4-Bromophenyl-phenylether	11.02	248	588680	39.097	ng	97
68) Hexachlorobenzene	11.09	284	680720	39.136	ng	100
69) Atrazine	11.17	200	516813	38.515	ng	99
70) Pentachlorophenol	11.27	266	374994	38.566	ng	97
71) Phenanthrene	11.50	178	2587794	39.160	ng	99
72) Anthracene	11.56	178	2584744	39.314	ng	100
73) Carbazole	11.70	167	2356769	39.213	ng	100
74) Di-n-butylphthalate	12.04	149	2834116	38.780	ng	100
75) Fluoranthene	12.69	202	2631005	38.991	ng	100
77) Benzidine	12.81	184	1100086	32.804	ng	99
78) Pyrene	12.93	202	2696816	38.300	ng	99
80) Butylbenzylphthalate	13.54	149	1228888	37.785	ng	99
81) Benzo(a)anthracene	14.12	228	2303432	38.365	ng	99
82) 3,3'-Dichlorobenzidine	14.07	252	877623	39.138	ng	98
83) Chrysene	14.16	228	2280370	38.460	ng	99
84) Bis(2-ethylhexyl)phthalate	14.10	149	1612922	38.445	ng	99
85) Di-n-octyl phthalate	14.73	149	2514577	38.089	ng	100
86) Indeno(1,2,3-cd)pyrene	17.18	276	2091162	37.893	ng	99
88) Benzo(b)fluoranthene	15.19	252	2220920	37.691	ng	98
89) Benzo(k)fluoranthene	15.23	252	2062639	39.248	ng	99
90) Benzo(a)pyrene	15.58	252	1937544	37.932	ng	99
91) Dibenzo(a,h)anthracene	17.20	278	1730080	37.765	ng	99
92) Benzo(g,h,i)perylene	17.64	276	1649759	37.159	ng	99

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

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