

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121319\
 Data File : BF118198.D
 Acq On : 13 Dec 2019 20:52
 Operator : JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Dec 13 22:49:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 06 17:34:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	103113	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	394445	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	209248	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	463821	20.00	ng	0.00
76) Chrysene-d12	14.06	240	292015	20.00	ng	0.00
87) Perylene-d12	15.55	264	316127	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.48	112	465171	79.01	ng	0.00
7) Phenol-d6	6.50	99	554295	77.84	ng	0.00
23) Nitrobenzene-d5	7.43	82	530719	75.69	ng	0.00
42) 2,4,6-Tribromophenol	10.71	330	225524	88.72	ng	0.00
45) 2-Fluorobiphenyl	9.23	172	1094469	79.59	ng	0.00
79) Terphenyl-d14	13.00	244	1465496	86.30	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.59	88	93666	37.732	ng	98
3) Pyridine	3.35	79	250852	39.168	ng	99
4) n-Nitrosodimethylamine	3.30	42	103468	37.554	ng	95
6) Aniline	6.53	93	341020	37.439	ng	96
8) 2-Chlorophenol	6.66	128	261395	39.376	ng	97
9) Benzaldehyde	6.42	77	148893	40.368	ng	97
10) Phenol	6.52	94	311350	38.340	ng	98
11) bis(2-Chloroethyl)ether	6.60	93	228676	38.935	ng	94
12) 1,3-Dichlorobenzene	6.81	146	307220	39.617	ng	98
13) 1,4-Dichlorobenzene	6.89	146	310667	39.785	ng	99
14) 1,2-Dichlorobenzene	7.04	146	293856	40.072	ng	98
15) Benzyl Alcohol	7.01	79	216778	38.999	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.15	45	276797	39.196	ng	98
17) 2-Methylphenol	7.12	107	202768	38.404	ng	98
18) Hexachloroethane	7.38	117	114273	39.715	ng	93
19) n-Nitroso-di-n-propylamine	7.28	70	170152	39.323	ng	98
20) 3+4-Methylphenols	7.27	107	258805	39.024	ng	95
22) Acetophenone	7.28	105	365244	38.542	ng	# 94
24) Nitrobenzene	7.45	77	258689	37.553	ng	98
25) Isophorone	7.69	82	450180	37.875	ng	99
26) 2-Nitrophenol	7.77	139	139060	37.769	ng	94
27) 2,4-Dimethylphenol	7.80	122	184034	37.472	ng	100
28) bis(2-Chloroethoxy)methane	7.90	93	296521	38.179	ng	99
29) 2,4-Dichlorophenol	8.02	162	221175	38.631	ng	97
30) 1,2,4-Trichlorobenzene	8.10	180	259256	38.747	ng	98
31) Naphthalene	8.17	128	771595	38.950	ng	100
32) Benzoic acid	7.93	122	165985	37.431	ng	99
33) 4-Chloroaniline	8.23	127	312403	36.523	ng	97
34) Hexachlorobutadiene	8.29	225	158018	39.384	ng	99
35) Caprolactam	8.60	113	80607	45.653	ng	99
36) 4-Chloro-3-methylphenol	8.71	107	236250	39.211	ng	99
37) 2-Methylnaphthalene	8.87	142	519043	38.702	ng	98
38) 1-Methylnaphthalene	8.97	142	485475	38.552	ng	98
40) 1,2,4,5-Tetrachlorobenzene	9.04	216	244408	38.557	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.02	237	69191	24.362	ng	99
43) 2,4,6-Trichlorophenol	9.15	196	162073	38.294	ng	98
44) 2,4,5-Trichlorophenol	9.19	196	175352	39.990	ng	98
46) 1,1'-Biphenyl	9.33	154	637529	38.631	ng	99
47) 2-Chloronaphthalene	9.36	162	512048	38.572	ng	98
48) 2-Nitroaniline	9.46	65	159435	42.115	ng	95
49) Acenaphthylene	9.78	152	766525	39.148	ng	99
50) Dimethylphthalate	9.63	163	693754	44.886	ng	98
51) 2,6-Dinitrotoluene	9.70	165	146297	43.530	ng	91
52) Acenaphthene	9.95	154	462553	38.704	ng	97
53) 3-Nitroaniline	9.87	138	166275	42.040	ng	98
54) 2,4-Dinitrophenol	9.98	184	52714	31.527	ng	99
55) Dibenzofuran	10.12	168	709586	40.521	ng	99
56) 4-Nitrophenol	10.04	139	111545	40.638	ng	94
57) 2,4-Dinitrotoluene	10.11	165	208347	46.446	ng	93
58) Fluorene	10.47	166	594180	42.696	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.25	232	161575	44.664	ng	98
60) Diethylphthalate	10.33	149	729608	47.912	ng	100
61) 4-Chlorophenyl-phenylether	10.46	204	295092	42.143	ng	97
62) 4-Nitroaniline	10.49	138	175959	42.656	ng	98
63) Azobenzene	10.62	77	562509	43.271	ng	99
65) 4,6-Dinitro-2-methylphenol	10.52	198	82406	32.171	ng	89
66) n-Nitrosodiphenylamine	10.57	169	560511	38.698	ng	100
67) 4-Bromophenyl-phenylether	10.95	248	199412	37.663	ng	97
68) Hexachlorobenzene	11.02	284	240354	38.569	ng	91
69) Atrazine	11.10	200	130489	30.336	ng	97
70) Pentachlorophenol	11.22	266	110064	37.556	ng	99
71) Phenanthrene	11.43	178	954852	38.727	ng	100
72) Anthracene	11.49	178	965951	39.272	ng	100
73) Carbazole	11.65	167	845417	38.309	ng	99
74) Di-n-butylphthalate	11.97	149	1159883	41.996	ng	99
75) Fluoranthene	12.63	202	945672	37.514	ng	98
77) Benzidine	12.75	184	93305	12.137	ng	99
78) Pyrene	12.86	202	948366	41.212	ng	99
80) Butylbenzylphthalate	13.47	149	438825	42.107	ng	98
81) Benzo(a)anthracene	14.05	228	768322	38.051	ng	99
82) 3,3'-Dichlorobenzidine	14.01	252	240784	36.309	ng	96
83) Chrysene	14.09	228	745732	38.663	ng	99
84) Bis(2-ethylhexyl)phthalate	14.03	149	609179	44.329	ng	# 99
85) Di-n-octyl phthalate	14.64	149	904343	43.442	ng	100
86) Indeno(1,2,3-cd)pyrene	17.07	276	841925	51.878	ng	100
88) Benzo(b)fluoranthene	15.12	252	718791	34.379	ng	99
89) Benzo(k)fluoranthene	15.14	252	724815	36.087	ng	99
90) Benzo(a)pyrene	15.49	252	690326	37.385	ng	99
91) Dibenzo(a,h)anthracene	17.09	278	692298	43.711	ng	99
92) Benzo(g,h,i)perylene	17.53	276	674883	44.344	ng	99

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

