

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
 Data File : BF109951.D
 Acq On : 18 Oct 2018 10:42
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 19 02:14:46 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 16 14:58:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	227244	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	936420	20.00	ng	0.00
39) Acenaphthene-d10	10.02	164	435946	20.00	ng	0.00
64) Phenanthrene-d10	11.51	188	805129	20.00	ng	0.00
76) Chrysene-d12	14.16	240	535318	20.00	ng	0.00
87) Perylene-d12	15.68	264	415922	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.59	112	1103471	76.97	ng	0.00
7) Phenol-d6	6.60	99	1389983	78.08	ng	0.00
23) Nitrobenzene-d5	7.55	82	1221948	87.90	ng	0.00
42) 2,4,6-Tribromophenol	10.82	330	335642	88.93	ng	0.00
45) 2-Fluorobiphenyl	9.34	172	2101929	76.94	ng	0.00
79) Terphenyl-d14	13.10	244	2035113	79.83	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.85	88	300509	36.907	ng	93
3) Pyridine	3.62	79	694500	38.258	ng	91
4) n-Nitrosodimethylamine	3.59	42	273980	36.999	ng	98
6) Aniline	6.64	93	925009	38.559	ng	# 53
8) 2-Chlorophenol	6.76	128	635232	39.558	ng	99
9) Benzaldehyde	6.53	77	434958	37.599	ng	95
10) Phenol	6.61	94	748538	38.947	ng	# 61
11) bis(2-Chloroethyl)ether	6.71	93	605756	37.695	ng	93
12) 1,3-Dichlorobenzene	6.92	146	686922	39.011	ng	96
13) 1,4-Dichlorobenzene	7.00	146	689951	38.906	ng	99
14) 1,2-Dichlorobenzene	7.15	146	640242	39.206	ng	98
15) Benzyl Alcohol	7.12	79	537541	39.250	ng	89
16) 2,2'-oxybis(1-Chloropropan	7.25	45	948650	35.966	ng	68
17) 2-Methylphenol	7.22	107	500182	38.627	ng	# 81
18) Hexachloroethane	7.49	117	251017	40.054	ng	97
19) n-Nitroso-di-n-propylamine	7.39	70	423086	37.016	ng	96
20) 3+4-Methylphenols	7.37	107	639241	38.902	ng	94
22) Acetophenone	7.39	105	820354	37.776	ng	98
24) Nitrobenzene	7.56	77	632300	41.871	ng	91
25) Isophorone	7.80	82	1023987	37.490	ng	96
26) 2-Nitrophenol	7.87	139	323310	51.637	ng	92
27) 2,4-Dimethylphenol	7.90	122	516390	39.250	ng	95
28) bis(2-Chloroethoxy)methane	8.00	93	693224	36.981	ng	98
29) 2,4-Dichlorophenol	8.11	162	518863	40.655	ng	98
30) 1,2,4-Trichlorobenzene	8.20	180	565346	39.562	ng	96
31) Naphthalene	8.29	128	1650550	38.438	ng	99
32) Benzoic acid	8.00	122	241693	29.485	ng	93
33) 4-Chloroaniline	8.33	127	752376	38.977	ng	99
34) Hexachlorobutadiene	8.40	225	311365	39.513	ng	98
35) Caprolactam	8.70	113	171841	37.912	ng	# 90
36) 4-Chloro-3-methylphenol	8.80	107	532768	39.570	ng	90
37) 2-Methylnaphthalene	8.97	142	1082914	38.946	ng	99
38) 1-Methylnaphthalene	9.07	142	1042990	38.710	ng	98
40) 1,2,4,5-Tetrachlorobenzene	9.14	216	531756	39.602	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.13	237	275050	42.440	ng	97
43) 2,4,6-Trichlorophenol	9.25	196	352552	41.990	ng	96
44) 2,4,5-Trichlorophenol	9.29	196	364657	41.846	ng	94
46) 1,1'-Biphenyl	9.44	154	1491622	38.403	ng	98
47) 2-Chloronaphthalene	9.47	162	1106788	38.685	ng	99
48) 2-Nitroaniline	9.56	65	343396	46.770	ng	96
49) Acenaphthylene	9.89	152	1608967	38.712	ng	97
50) Dimethylphthalate	9.74	163	1191008	38.583	ng	99
51) 2,6-Dinitrotoluene	9.80	165	277867	47.393	ng #	82
52) Acenaphthene	10.06	154	1051673	39.744	ng	98
53) 3-Nitroaniline	9.97	138	327323	45.895	ng	87
54) 2,4-Dinitrophenol	10.07	184	101644	57.269	ng #	82
55) Dibenzofuran	10.23	168	1462588	38.757	ng	97
56) 4-Nitrophenol	10.12	139	244244	44.882	ng	89
57) 2,4-Dinitrotoluene	10.20	165	362880	48.901	ng #	87
58) Fluorene	10.57	166	1064718	38.758	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.34	232	298028	43.154	ng	97
60) Diethylphthalate	10.43	149	1184421	38.848	ng	98
61) 4-Chlorophenyl-phenylether	10.56	204	567459	39.655	ng	99
62) 4-Nitroaniline	10.59	138	316060	46.815	ng	90
63) Azobenzene	10.72	77	1170272	36.742	ng	95
65) 4,6-Dinitro-2-methylphenol	10.61	198	143654	49.045	ng	86
66) n-Nitrosodiphenylamine	10.68	169	1028687	38.361	ng	97
67) 4-Bromophenyl-phenylether	11.05	248	341838	39.177	ng	94
68) Hexachlorobenzene	11.12	284	367238	39.468	ng	99
69) Atrazine	11.20	200	334909	39.983	ng	99
70) Pentachlorophenol	11.31	266	227525	42.967	ng	96
71) Phenanthrene	11.54	178	1581791	37.949	ng	99
72) Anthracene	11.59	178	1609037	38.397	ng	100
73) Carbazole	11.74	167	1569317	38.105	ng	99
74) Di-n-butylphthalate	12.06	149	1732579	37.695	ng	99
75) Fluoranthene	12.73	202	1580206	37.949	ng	97
77) Benzidine	12.84	184	819342	39.920	ng	98
78) Pyrene	12.96	202	1615699	41.094	ng	100
80) Butylbenzylphthalate	13.57	149	663168	41.720	ng	94
81) Benzo(a)anthracene	14.14	228	1216747	38.588	ng	99
82) 3,3'-Dichlorobenzidine	14.10	252	435267	39.375	ng	97
83) Chrysene	14.19	228	1165297	38.826	ng	100
84) Bis(2-ethylhexyl)phthalate	14.12	149	831967	39.777	ng #	94
85) Di-n-octyl phthalate	14.74	149	1193300	40.467	ng	99
86) Indeno(1,2,3-cd)pyrene	17.25	276	1011858	41.872	ng #	94
88) Benzo(b)fluoranthene	15.23	252	974354	40.634	ng	99
89) Benzo(k)fluoranthene	15.26	252	868297	36.738	ng	99
90) Benzo(a)pyrene	15.62	252	858239	38.916	ng	98
91) Dibenzo(a,h)anthracene	17.27	278	844718	43.220	ng	97
92) Benzo(g,h,i)perylene	17.73	276	863090	43.707	ng #	96

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

