

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050218\
 Data File : BF105040.D
 Acq On : 2 May 2018 19:23
 Operator : JU/SJ
 Sample : PB108708BS
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB108708BS

Quant Time: May 03 04:24:55 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 02 17:25:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	83076	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	347335	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	153980	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	255413	20.00	ng	0.00
75) Chrysene-d12	14.05	240	183843	20.00	ng	0.02
86) Perylene-d12	15.64	264	148157	20.00	ng	0.05

System Monitoring Compounds

5) 2-Fluorophenol	5.41	112	593126	109.17	ng	0.00
7) Phenol-d6	6.43	99	711081	106.52	ng	0.00
23) Nitrobenzene-d5	7.36	82	430387	80.91	ng	-0.01
41) 2,4,6-Tribromophenol	10.63	330	170362	106.66	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	831269	85.28	ng	0.00
78) Terphenyl-d14	12.92	244	700425	86.86	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.55	88	62194	24.567	ng	87
3) Pyridine	3.29	79	176069	24.736	ng	86
4) n-Nitrosodimethylamine	3.24	42	67599	27.169	ng	# 82
6) Aniline	6.46	93	89090	9.606	ng	# 88
8) 2-Chlorophenol	6.58	128	190280	33.181	ng	91
9) Benzaldehyde	6.34	77	59877	13.960	ng	99
10) Phenol	6.45	94	219018	30.921	ng	95
11) bis(2-Chloroethyl)ether	6.53	93	167074	29.559	ng	93
12) 1,3-Dichlorobenzene	6.73	146	198312	30.526	ng	98
13) 1,4-Dichlorobenzene	6.80	146	202614	31.287	ng	99
14) 1,2-Dichlorobenzene	6.96	146	190117	31.679	ng	99
15) Benzyl Alcohol	6.94	79	146809	28.955	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.06	45	187516	24.703	ng	55
17) 2-Methylphenol	7.05	107	152073	30.508	ng	95
18) Hexachloroethane	7.30	117	60628	26.258	ng	94
19) n-Nitroso-di-n-propylamine	7.20	70	116349	26.672	ng	92
20) 3+4-Methylphenols	7.21	107	201752	32.634	ng	# 88
22) Acetophenone	7.20	105	253713	29.670	ng	# 97
24) Nitrobenzene	7.38	77	185781	31.634	ng	93
25) Isophorone	7.62	82	331333	29.456	ng	99
26) 2-Nitrophenol	7.69	139	90965	38.048	ng	95
27) 2,4-Dimethylphenol	7.73	122	165225	33.283	ng	99
28) bis(2-Chloroethoxy)methane	7.82	93	216296	31.451	ng	99
29) 2,4-Dichlorophenol	7.94	162	156967	33.139	ng	98
30) 1,2,4-Trichlorobenzene	8.02	180	161165	31.004	ng	99
31) Naphthalene	8.10	128	536121	30.998	ng	99
32) Benzoic acid	7.86	122	68880	17.390	ng	96
33) 4-Chloroaniline	8.15	127	46704	6.303	ng	96
34) Hexachlorobutadiene	8.20	225	87832	30.848	ng	99
35) Caprolactam	8.52	113	48239	29.194	ng	# 73
36) 4-Chloro-3-methylphenol	8.64	107	162947	32.083	ng	99
37) 2-Methylnaphthalene	8.79	142	356267	31.845	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.96	216	152414	34.578	ng	98
40) Hexachlorocyclopentadiene	8.93	237	9186	3.723	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.07	196	99035	32.366	ng	98
43) 2,4,5-Trichlorophenol	9.12	196	107716	31.459	ng	93
45) 1,1'-Biphenyl	9.25	154	422075	32.629	ng	99
46) 2-Chloronaphthalene	9.28	162	324054	31.716	ng	99
47) 2-Nitroaniline	9.38	65	92774	36.717	ng	95
48) Acenaphthylene	9.69	152	484846	30.245	ng	99
49) Dimethylphthalate	9.55	163	389488	32.076	ng	100
50) 2,6-Dinitrotoluene	9.62	165	81185	36.133	ng	96
51) Acenaphthene	9.86	154	284098	29.046	ng	99
52) 3-Nitroaniline	9.79	138	43837	16.666	ng	96
53) 2,4-Dinitrophenol	9.92	184	10942	20.992	ng	# 23
54) Dibenzofuran	10.04	168	424134	31.116	ng	99
55) 4-Nitrophenol	9.97	139	100171	48.480	ng	96
56) 2,4-Dinitrotoluene	10.03	165	99922	33.904	ng	98
57) Fluorene	10.38	166	334845	33.750	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.16	232	76546	29.965	ng	94
59) Diethylphthalate	10.25	149	378408	31.291	ng	99
60) 4-Chlorophenyl-phenylether	10.37	204	161554	36.563	ng	99
61) 4-Nitroaniline	10.41	138	72880	28.337	ng	97
62) Azobenzene	10.53	77	320422	27.734	ng	94
64) 4,6-Dinitro-2-methylphenol	10.45	198	10678	14.617	ng	91
65) n-Nitrosodiphenylamine	10.49	169	298944	33.757	ng	99
66) 4-Bromophenyl-phenylether	10.86	248	93634	34.465	ng	90
67) Hexachlorobenzene	10.93	284	98049	32.074	ng	# 90
68) Atrazine	11.02	200	94292	35.274	ng	99
69) Pentachlorophenol	11.13	266	74636	39.465	ng	98
70) Phenanthrene	11.35	178	439436	30.894	ng	97
71) Anthracene	11.40	178	463184	32.035	ng	98
72) Carbazole	11.56	167	415131	29.382	ng	99
73) Di-n-butylphthalate	11.88	149	554243	32.113	ng	99
74) Fluoranthene	12.55	202	428067	28.805	ng	97
76) Benzidine	12.67	184	242253	38.446	ng	97
77) Pyrene	12.77	202	423128	31.051	ng	99
79) Butylbenzylphthalate	13.41	149	207437	32.311	ng	97
80) Benzo(a)anthracene	14.04	228	371885	33.860	ng	99
81) 3,3'-Dichlorobenzidine	14.00	252	125244	30.584	ng	96
82) Chrysene	14.08	228	352547	31.251	ng	98
83) Bis(2-ethylhexyl)phthalate	14.02	149	292330	35.401	ng	100
84) Di-n-octyl phthalate	14.72	149	486646	35.270	ng	# 93
85) Indeno(1,2,3-cd)pyrene	17.13	276	262562	27.398	ng	95
87) Benzo(b)fluoranthene	15.20	252	294705	31.495	ng	100
88) Benzo(k)fluoranthene	15.23	252	308404	34.910	ng	98
89) Benzo(a)pyrene	15.58	252	276228	32.992	ng	99
90) Dibenzo(a,h)anthracene	17.14	278	223006	32.431	ng	98
91) Benzo(g,h,i)perylene	17.58	276	208375	31.278	ng	95

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

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