

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050118\
 Data File : BF105017.D
 Acq On : 2 May 2018 7:47
 Operator : JU/SJ
 Sample : PB108770BSD
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB108770BSD

Quant Time: May 02 09:01:01 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 01 16:13:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	118872	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	507911	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	235882	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	385160	20.00	ng	0.00
75) Chrysene-d12	14.02	240	248143	20.00	ng	0.05
86) Perylene-d12	15.57	264	223111	20.00	ng	0.14

System Monitoring Compounds

5) 2-Fluorophenol	5.41	112	593387	76.33	ng	0.00
7) Phenol-d6	6.43	99	722020	75.59	ng	0.00
23) Nitrobenzene-d5	7.36	82	626067	80.49	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	178807	73.08	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1186330	79.45	ng	0.00
78) Terphenyl-d14	12.92	244	935839	85.98	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.57	88	101367	27.983	ng	# 88
3) Pyridine	3.30	79	285671	28.049	ng	# 83
4) n-Nitrosodimethylamine	3.27	42	110563	31.055	ng	# 77
6) Aniline	6.45	93	139043	10.477	ng	# 1
8) 2-Chlorophenol	6.58	128	307458	37.470	ng	91
9) Benzaldehyde	6.34	77	93374	15.506	ng	95
10) Phenol	6.45	94	358742	35.396	ng	95
11) bis(2-Chloroethyl)ether	6.53	93	277439	34.303	ng	93
12) 1,3-Dichlorobenzene	6.73	146	316092	34.004	ng	99
13) 1,4-Dichlorobenzene	6.80	146	321985	34.748	ng	98
14) 1,2-Dichlorobenzene	6.96	146	298697	34.784	ng	99
15) Benzyl Alcohol	6.94	79	233226	32.147	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.06	45	308782	28.429	ng	53
17) 2-Methylphenol	7.05	107	248301	34.812	ng	# 93
18) Hexachloroethane	7.30	117	109102	33.023	ng	95
19) n-Nitroso-di-n-propylamine	7.21	70	193072	30.932	ng	91
20) 3+4-Methylphenols	7.21	107	310994	35.156	ng	# 83
22) Acetophenone	7.20	105	400572	32.034	ng	99
24) Nitrobenzene	7.38	77	299245	34.845	ng	94
25) Isophorone	7.62	82	548644	33.355	ng	99
26) 2-Nitrophenol	7.69	139	163970	46.901	ng	94
27) 2,4-Dimethylphenol	7.73	122	264636	36.455	ng	98
28) bis(2-Chloroethoxy)methane	7.82	93	355616	35.361	ng	99
29) 2,4-Dichlorophenol	7.94	162	256354	37.011	ng	97
30) 1,2,4-Trichlorobenzene	8.02	180	260693	34.295	ng	99
31) Naphthalene	8.10	128	862706	34.111	ng	100
32) Benzoic acid	7.87	122	113517	19.599	ng	98
33) 4-Chloroaniline	8.15	127	66232	6.113	ng	96
34) Hexachlorobutadiene	8.20	225	142817	34.301	ng	97
35) Caprolactam	8.53	113	79408	32.864	ng	# 75
36) 4-Chloro-3-methylphenol	8.65	107	268862	36.201	ng	99
37) 2-Methylnaphthalene	8.79	142	580796	35.502	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.96	216	252013	37.323	ng	99
40) Hexachlorocyclopentadiene	8.93	237	60588	16.028	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.07	196	168271	35.899	ng	100
43) 2,4,5-Trichlorophenol	9.12	196	181042	34.515	ng	93
45) 1,1'-Biphenyl	9.25	154	687354	34.687	ng	98
46) 2-Chloronaphthalene	9.28	162	536959	34.306	ng	100
47) 2-Nitroaniline	9.39	65	152898	39.501	ng	87
48) Acenaphthylene	9.70	152	787340	32.061	ng	100
49) Dimethylphthalate	9.55	163	655098	35.218	ng	100
50) 2,6-Dinitrotoluene	9.63	165	140483	40.815	ng	# 86
51) Acenaphthene	9.87	154	464500	31.001	ng	100
52) 3-Nitroaniline	9.79	138	53119	13.183	ng	99
53) 2,4-Dinitrophenol	9.91	184	43320	40.841	ng	# 23
54) Dibenzofuran	10.04	168	677362	32.440	ng	97
55) 4-Nitrophenol	9.97	139	166286	52.534	ng	97
56) 2,4-Dinitrotoluene	10.03	165	170780	37.827	ng	94
57) Fluorene	10.39	166	550133	36.196	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.16	232	129761	33.160	ng	93
59) Diethylphthalate	10.25	149	619320	33.431	ng	99
60) 4-Chlorophenyl-phenylether	10.37	204	260278	38.453	ng	95
61) 4-Nitroaniline	10.42	138	122489	31.089	ng	95
62) Azobenzene	10.53	77	546385	30.871	ng	93
64) 4,6-Dinitro-2-methylphenol	10.45	198	39141	25.019	ng	78
65) n-Nitrosodiphenylamine	10.49	169	498547	37.332	ng	99
66) 4-Bromophenyl-phenylether	10.86	248	158628	38.719	ng	93
67) Hexachlorobenzene	10.93	284	168462	36.544	ng	# 89
68) Atrazine	11.03	200	161497	40.064	ng	99
69) Pentachlorophenol	11.13	266	120368	42.207	ng	98
70) Phenanthrene	11.35	178	728415	33.960	ng	99
71) Anthracene	11.40	178	757030	34.721	ng	99
72) Carbazole	11.56	167	676469	31.750	ng	99
73) Di-n-butylphthalate	11.88	149	916942	35.231	ng	99
74) Fluoranthene	12.54	202	671415	29.960	ng	97
76) Benzidine	12.67	184	276722	30.657	ng	99
77) Pyrene	12.77	202	648544	35.260	ng	99
79) Butylbenzylphthalate	13.40	149	330516	38.142	ng	93
80) Benzo(a)anthracene	14.01	228	554445	37.401	ng	98
81) 3,3'-Dichlorobenzidine	13.97	252	172495	31.208	ng	98
82) Chrysene	14.05	228	522828	34.336	ng	98
83) Bis(2-ethylhexyl)phthalate	13.99	149	460388	41.306	ng	100
84) Di-n-octyl phthalate	14.67	149	735575	39.497	ng	# 93
85) Indeno(1,2,3-cd)pyrene	17.06	276	426814	32.997	ng	96
87) Benzo(b)fluoranthene	15.14	252	522880	37.107	ng	98
88) Benzo(k)fluoranthene	15.17	252	479066	36.011	ng	99
89) Benzo(a)pyrene	15.52	252	467056	37.044	ng	100
90) Dibenzo(a,h)anthracene	17.07	278	366083	35.353	ng	98
91) Benzo(g,h,i)perylene	17.50	276	338887	33.779	ng	96

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

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