

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
 Data File : BF079720.D
 Acq On : 5 Jun 2015 1:20
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Jun 06 03:31:22 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 05 19:33:57 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	105053	20.00	ng	0.00
21) Naphthalene-d8	8.80	136	411415	20.00	ng	0.00
38) Acenaphthene-d10	10.58	164	213635	20.00	ng	0.00
63) Phenanthrene-d10	12.08	188	426935	20.00	ng	-0.02
75) Chrysene-d12	14.98	240	464181	20.00	ng	-0.18
86) Perylene-d12	16.91	264	446065	20.00	ng	-0.25

System Monitoring Compounds						
5) 2-Fluorophenol	6.12	112	485672	79.27	ng	0.00
7) Phenol-d6	7.10	99	616639	77.32	ng	0.00
23) Nitrobenzene-d5	8.07	82	579066	85.70	ng	0.00
41) 2,4,6-Tribromophenol	11.37	330	153254	79.88	ng	0.00
44) 2-Fluorobiphenyl	9.87	172	1132098	76.72	ng	0.00
78) Terphenyl-d14	13.74	244	1505267	76.40	ng	-0.09

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.59	88	105806	38.23	ng	98
3) Pyridine	4.30	79	311753	40.23	ng	96
4) n-Nitrosodimethylamine	4.22	42	141832	39.11	ng	88
6) Aniline	7.16	93	461711	41.58	ng	99
8) 2-Chlorophenol	7.29	128	280377	39.91	ng	99
9) Benzaldehyde	7.05	77	206338	41.14	ng	98
10) Phenol	7.12	94	320704	38.79	ng	100
11) bis(2-Chloroethyl)ether	7.22	93	288654	42.05	ng	100
12) 1,3-Dichlorobenzene	7.45	146	317875	39.79	ng	99
13) 1,4-Dichlorobenzene	7.53	146	339849	39.85	ng	99
14) 1,2-Dichlorobenzene	7.69	146	300335	39.91	ng	99
15) Benzyl Alcohol	7.63	79	243625	40.48	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.77	45	462177	40.68	ng	100
17) 2-Methylphenol	7.72	107	229631	39.98	ng	# 92
18) Hexachloroethane	8.03	117	110405	41.20	ng	94
19) n-Nitroso-di-n-propylamine	7.90	70	231722	42.00	ng	97
20) 3+4-Methylphenols	7.88	107	317345	40.46	ng	100
22) Acetophenone	7.91	105	415573	41.47	ng	# 98
24) Nitrobenzene	8.08	77	281399	40.28	ng	98
25) Isophorone	8.32	82	592766	39.94	ng	100
26) 2-Nitrophenol	8.40	139	105109	40.46	ng	# 66
27) 2,4-Dimethylphenol	8.42	122	268037	39.78	ng	99
28) bis(2-Chloroethoxy)methane	8.51	93	317202	38.33	ng	99
29) 2,4-Dichlorophenol	8.64	162	237972	41.12	ng	98
30) 1,2,4-Trichlorobenzene	8.74	180	276195	39.69	ng	99
31) Naphthalene	8.82	128	884320	39.01	ng	99
32) Benzoic acid	8.49	122	96453	38.65	ng	99
33) 4-Chloroaniline	8.86	127	360204m	40.00	ng	
34) Hexachlorobutadiene	8.94	225	159972	39.56	ng	100
35) Caprolactam	9.20	113	71086	40.52	ng	96
36) 4-Chloro-3-methylphenol	9.31	107	235908	38.79	ng	# 75
37) 2-Methylnaphthalene	9.52	142	616437	40.30	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.69	216	274256	39.75	ng	99
40) Hexachlorocyclopentadiene	9.68	237	135680	41.14	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.79	196	171610	40.28	ng	98
43) 2,4,5-Trichlorophenol	9.83	196	202774	41.54	ng	97
45) 1,1'-Biphenyl	9.98	154	665295	38.90	ng	94
46) 2-Chloronaphthalene	10.01	162	560153	40.10	ng	99
47) 2-Nitroaniline	10.09	65	129750	42.57	ng	95
48) Acenaphthylene	10.43	152	941659	39.40	ng	99
49) Dimethylphthalate	10.26	163	651548	39.59	ng	100
50) 2,6-Dinitrotoluene	10.33	165	122154	39.86	ng	97
51) Acenaphthene	10.62	154	557282	39.62	ng	99
52) 3-Nitroaniline	10.50	138	141470	42.60	ng	99
53) 2,4-Dinitrophenol	10.60	184	26841	41.44	ng	71
54) Dibenzofuran	10.79	168	793546	39.36	ng	98
55) 4-Nitrophenol	10.64	139	120776	43.12	ng	97
56) 2,4-Dinitrotoluene	10.74	165	146844	42.11	ng	97
57) Fluorene	11.13	166	664544	38.58	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.89	232	147606	40.79	ng	100
59) Diethylphthalate	10.97	149	670542	40.48	ng	99
60) 4-Chlorophenyl-phenylether	11.11	204	301073	39.70	ng	98
61) 4-Nitroaniline	11.12	138	139322	40.13	ng	100
62) Azobenzene	11.27	77	631193	40.03	ng	99
64) 4,6-Dinitro-2-methylphenol	11.15	198	50121	43.46	ng	88
65) n-Nitrosodiphenylamine	11.22	169	547690	40.78	ng	99
66) 4-Bromophenyl-phenylether	11.61	248	204608	45.26	ng	# 90
67) Hexachlorobenzene	11.69	284	207358	40.64	ng	# 68
68) Atrazine	11.74	200	175651	38.87	ng	93
69) Pentachlorophenol	11.87	266	93156	38.17	ng	99
70) Phenanthrene	12.11	178	1030146	41.48	ng	100
71) Anthracene	12.16	178	988514	40.22	ng	97
72) Carbazole	12.31	167	977553	42.68	ng	98
73) Di-n-butylphthalate	12.63	149	1077581	39.56	ng	# 80
74) Fluoranthene	13.35	202	1091432	40.51	ng	97
76) Benzidine	13.46	184	611618	43.78	ng	99
77) Pyrene	13.60	202	1117843	38.76	ng	99
79) Butylbenzylphthalate	14.26	149	466397	41.94	ng	# 77
80) Benzo(a)anthracene	14.97	228	1117078	39.37	ng	99
81) 3,3'-Dichlorobenzidine	14.90	252	380255	38.95	ng	# 98
82) Chrysene	15.02	228	1041584	39.22	ng	99
83) Bis(2-ethylhexyl)phthalate	14.93	149	723319	41.45	ng	# 97
84) Di-n-octyl phthalate	15.71	149	1199667	40.97	ng	99
85) Indeno(1,2,3-cd)pyrene	18.93	276	1171358	39.09	ng	99
87) Benzo(h)fluoranthene	16.33	252	1068196m	37.12	ng	
88) Benzo(k)fluoranthene	16.38	252	1077537	39.25	ng	98
89) Benzo(a)pyrene	16.83	252	1005148	38.24	ng	99
90) Dibenzo(a,h)anthracene	18.95	278	941835	38.78	ng	99
91) Benzo(g,h,i)perylene	19.53	276	974454	39.24	ng	99

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

