

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100617\
 Data File : BF099138.D
 Acq On : 6 Oct 2017 10:12
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 06 14:30:30 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 05 17:20:48 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	171366	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	717810	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	319332	20.00	ng	0.00
63) Phenanthrene-d10	11.39	188	550177	20.00	ng	0.00
75) Chrysene-d12	14.03	240	389589	20.00	ng	0.00
86) Perylene-d12	15.50	264	312575	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	829777	77.08	ng	0.00
7) Phenol-d6	6.50	99	1054348	79.40	ng	0.00
23) Nitrobenzene-d5	7.44	82	878105	78.45	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	209810	80.29	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1466385	76.66	ng	0.00
78) Terphenyl-d14	12.97	244	1355543	79.13	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.63	88	190914	37.13	ng	96
3) Pyridine	3.40	79	523188	37.61	ng	94
4) n-Nitrosodimethylamine	3.38	42	208590	35.36	ng	83
6) Aniline	6.54	93	696659	38.87	ng	98
8) 2-Chlorophenol	6.66	128	478174	39.22	ng	95
9) Benzaldehyde	6.42	77	269546	34.99	ng	93
10) Phenol	6.52	94	610351	41.10	ng	98
11) bis(2-Chloroethyl)ether	6.61	93	455541	39.45	ng	95
12) 1,3-Dichlorobenzene	6.81	146	505712	39.61	ng	98
13) 1,4-Dichlorobenzene	6.89	146	506273	38.97	ng	98
14) 1,2-Dichlorobenzene	7.04	146	470026	39.16	ng	99
15) Benzyl Alcohol	7.01	79	358903	36.84	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.15	45	645730	35.90	ng	97
17) 2-Methylphenol	7.12	107	393284	39.43	ng	97
18) Hexachloroethane	7.38	117	177564	38.03	ng	96
19) n-Nitroso-di-n-propylamine	7.29	70	300347	35.42	ng	96
20) 3+4-Methylphenols	7.27	107	451419	35.77	ng	# 85
22) Acetophenone	7.29	105	620262	35.67	ng	# 93
24) Nitrobenzene	7.46	77	482372	37.99	ng	95
25) Isophorone	7.69	82	882742	38.15	ng	99
26) 2-Nitrophenol	7.77	139	274513	44.96	ng	88
27) 2,4-Dimethylphenol	7.80	122	445429	38.63	ng	97
28) bis(2-Chloroethoxy)methane	7.90	93	563085	39.04	ng	99
29) 2,4-Dichlorophenol	8.01	162	396851	40.09	ng	97
30) 1,2,4-Trichlorobenzene	8.09	180	393013	39.69	ng	100
31) Naphthalene	8.17	128	1283282	38.07	ng	99
32) Benzoic acid	7.95	122	343517	36.27	ng	96
33) 4-Chloroaniline	8.22	127	548585	38.97	ng	98
34) Hexachlorobutadiene	8.29	225	197603	38.75	ng	99
35) Caprolactam	8.62	113	126030	39.69	ng	# 85
36) 4-Chloro-3-methylphenol	8.70	107	429438	38.59	ng	97
37) 2-Methylnaphthalene	8.86	142	845200	38.57	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.03	216	380251	39.93	ng	99
40) Hexachlorocyclopentadiene	9.01	237	146829	33.74	ng	98

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42) 2,4,6-Trichlorophenol	9.14	196	263246	39.02	ng	100
43) 2,4,5-Trichlorophenol	9.18	196	267447	39.71	ng	97
45) 1,1'-Biphenyl	9.33	154	1058911	38.58	ng	100
46) 2-Chloronaphthalene	9.36	162	815491	39.58	ng	96
47) 2-Nitroaniline	9.45	65	245007	38.83	ng	94
48) Acenaphthylene	9.77	152	1223990	38.80	ng	100
49) Dimethylphthalate	9.63	163	957620	39.73	ng	99
50) 2,6-Dinitrotoluene	9.69	165	219864	41.90	ng	93
51) Acenaphthene	9.94	154	721933	36.13	ng	99
52) 3-Nitroaniline	9.87	138	258024	42.17	ng	88
53) 2,4-Dinitrophenol	9.97	184	95906	38.38	ng	94
54) Dibenzofuran	10.11	168	1025997	38.56	ng	99
55) 4-Nitrophenol	10.02	139	191102	36.94	ng	91
56) 2,4-Dinitrotoluene	10.10	165	282890	41.54	ng	94
57) Fluorene	10.46	166	819983	38.35	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.23	232	179655	38.62	ng	99
59) Diethylphthalate	10.33	149	891815	37.87	ng	97
60) 4-Chlorophenyl-phenylether	10.45	204	368469	38.36	ng	94
61) 4-Nitroaniline	10.49	138	257099	43.56	ng	89
62) Azobenzene	10.60	77	786720	36.33	ng	94
64) 4,6-Dinitro-2-methylphenol	10.51	198	152590	45.58	ng	80
65) n-Nitrosodiphenylamine	10.57	169	753165	39.52	ng	99
66) 4-Bromophenyl-phenylether	10.93	248	215048	39.93	ng	96
67) Hexachlorobenzene	11.00	284	233413	40.58	ng	92
68) Atrazine	11.09	200	226223	39.45	ng	100
69) Pentachlorophenol	11.19	266	127528	35.63	ng	98
70) Phenanthrene	11.42	178	1152951	39.15	ng	100
71) Anthracene	11.47	178	1184559	39.31	ng	100
72) Carbazole	11.63	167	1166000	39.51	ng	99
73) Di-n-butylphthalate	11.95	149	1352087	38.07	ng	99
74) Fluoranthene	12.60	202	1165681	39.09	ng	99
76) Benzidine	12.72	184	535835	37.19	ng	99
77) Pyrene	12.84	202	1213294	40.84	ng	100
79) Butylbenzylphthalate	13.45	149	567031	40.61	ng	92
80) Benzo(a)anthracene	14.02	228	948389	40.20	ng	99
81) 3,3'-Dichlorobenzidine	13.99	252	352936	40.81	ng	98
82) Chrysene	14.06	228	890311	39.64	ng	99
83) Bis(2-ethylhexyl)phthalate	14.00	149	734282	38.19	ng	99
84) Di-n-octyl phthalate	14.61	149	1165840	37.66	ng	96
85) Indeno(1,2,3-cd)pyrene	16.99	276	704414	39.21	ng	96
87) Benzo(b)fluoranthene	15.07	252	749890	39.02	ng	97
88) Benzo(k)fluoranthene	15.10	252	748073	39.41	ng	99
89) Benzo(a)pyrene	15.44	252	701975	39.79	ng	# 97
90) Dibenzo(a,h)anthracene	17.00	278	563467	39.91	ng	96
91) Benzo(g,h,i)perylene	17.44	276	591604	42.39	ng	95

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

