

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022817\
 Data File : BG026019.D
 Acq On : 1 Mar 2017 10:42
 Operator : SJ/MA
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Mar 01 12:33:18 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG021717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 28 15:36:56 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	129597	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	630603	20.00	ng	-0.01
38) Acenaphthene-d10	14.68	164	463933	20.00	ng	-0.01
63) Phenanthrene-d10	17.41	188	950387	20.00	ng	0.00
75) Chrysene-d12	21.66	240	940625	20.00	ng	-0.02
86) Perylene-d12	24.86	264	960553	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	609142	81.83	ng	-0.01
7) Phenol-d6	7.23	99	899304	81.64	ng	0.00
23) Nitrobenzene-d5	9.23	82	805623	80.62	ng	-0.01
41) 2,4,6-Tribromophenol	16.16	330	381866	89.09	ng	0.00
44) 2-Fluorobiphenyl	13.31	172	2057884	77.51	ng	-0.01
78) Terphenyl-d14	20.01	244	2884017	74.57	ng	-0.01

Target Compounds						Qvalue
2) 1,4-Dioxane	3.54	88	126320	37.13	ng	# 80
3) Pyridine	3.94	79	378892	37.37	ng	# 69
4) n-Nitrosodimethylamine	3.85	42	125234	34.35	ng	# 19
6) Aniline	7.40	93	593201	38.73	ng	# 90
8) 2-Chlorophenol	7.64	128	362851	41.01	ng	# 85
9) Benzaldehyde	7.21	77	228659	35.03	ng	# 73
10) Phenol	7.26	94	481928	40.26	ng	# 74
11) bis(2-Chloroethyl)ether	7.49	93	370650	37.67	ng	# 83
12) 1,3-Dichlorobenzene	7.97	146	404321	39.54	ng	# 86
13) 1,4-Dichlorobenzene	8.12	146	419942	40.54	ng	# 93
14) 1,2-Dichlorobenzene	8.43	146	400725	39.40	ng	# 92
15) Benzyl Alcohol	8.31	79	315102	39.06	ng	# 78
16) 2,2'-oxybis(1-Chloropropan	8.60	45	443496	31.06	ng	# 88
17) 2-Methylphenol	8.51	107	346782	40.22	ng	# 81
18) Hexachloroethane	9.17	117	136972	39.89	ng	# 89
19) n-Nitroso-di-n-propylamine	8.87	70	294361	36.59	ng	# 76
20) 3+4-Methylphenols	8.84	107	501775	40.84	ng	# 88
22) Acetophenone	8.89	105	591118	39.06	ng	# 78
24) Nitrobenzene	9.27	77	425894	39.41	ng	# 77
25) Isophorone	9.80	82	851576	38.73	ng	# 91
26) 2-Nitrophenol	9.98	139	232577	46.03	ng	# 68
27) 2,4-Dimethylphenol	10.04	122	392193	41.59	ng	# 87
28) bis(2-Chloroethoxy)methane	10.28	93	529682	38.01	ng	# 96
29) 2,4-Dichlorophenol	10.52	162	401994	44.66	ng	# 94
30) 1,2,4-Trichlorobenzene	10.74	180	408944	41.77	ng	# 99
31) Naphthalene	10.93	128	1299430	39.86	ng	# 100
32) Benzoic acid	10.17	122	333849	46.36	ng	# 77
33) 4-Chloroaniline	11.02	127	585397	41.34	ng	# 84
34) Hexachlorobutadiene	11.22	225	227155	40.78	ng	# 98
35) Caprolactam	11.79	113	160942	44.28	ng	# 67
36) 4-Chloro-3-methylphenol	12.14	107	461524	42.86	ng	# 77
37) 2-Methylnaphthalene	12.52	142	1007505	40.57	ng	# 93
39) 1,2,4,5-Tetrachlorobenzene	12.89	216	477809	41.09	ng	# 98
40) Hexachlorocyclopentadiene	12.88	237	243737	38.31	ng	# 96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.12	196	331661	44.16	ng	98
43) 2,4,5-Trichlorophenol	13.19	196	385696	45.24	ng #	94
45) 1,1'-Biphenyl	13.52	154	1321364	39.30	ng	99
46) 2-Chloronaphthalene	13.56	162	976940	40.20	ng	97
47) 2-Nitroaniline	13.76	65	318014	41.25	ng #	65
48) Acenaphthylene	14.40	152	1731856	40.02	ng	99
49) Dimethylphthalate	14.13	163	1242273	39.03	ng	99
50) 2,6-Dinitrotoluene	14.25	165	311021	43.50	ng #	61
51) Acenaphthene	14.74	154	1153840	40.37	ng	100
52) 3-Nitroaniline	14.57	138	340589	42.59	ng #	72
53) 2,4-Dinitrophenol	14.78	184	160899	43.01	ng #	82
54) Dibenzofuran	15.08	168	1516795	39.81	ng #	89
55) 4-Nitrophenol	14.87	139	266421	40.93	ng #	44
56) 2,4-Dinitrotoluene	15.03	165	425331	44.22	ng #	74
57) Fluorene	15.72	166	1349048	39.62	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.30	232	324453	43.14	ng	99
59) Diethylphthalate	15.48	149	1278928	38.85	ng	99
60) 4-Chlorophenyl-phenylether	15.71	204	640260	40.58	ng	93
61) 4-Nitroaniline	15.73	138	345311	41.65	ng #	53
62) Azobenzene	16.00	77	1059848	37.45	ng	84
64) 4,6-Dinitro-2-methylphenol	15.80	198	243885	43.85	ng	88
65) n-Nitrosodiphenylamine	15.92	169	1176776	40.83	ng	98
66) 4-Bromophenyl-phenylether	16.60	248	421492	42.31	ng	97
67) Hexachlorobenzene	16.72	284	431064	41.99	ng	94
68) Atrazine	16.86	200	409920	40.18	ng	98
69) Pentachlorophenol	17.06	266	271572	42.88	ng	97
70) Phenanthrene	17.45	178	2012133	39.00	ng	98
71) Anthracene	17.55	178	2075592	39.46	ng	99
72) Carbazole	17.80	167	2040577	38.62	ng	97
73) Di-n-butylphthalate	18.36	149	2104255	40.57	ng #	95
74) Fluoranthene	19.45	202	2193731	38.62	ng	95
76) Benzidine	19.62	184	978756	32.96	ng	98
77) Pyrene	19.81	202	2275302	38.25	ng	99
79) Butylbenzylphthalate	20.69	149	962135	42.62	ng	86
80) Benzo(a)anthracene	21.64	228	2180045	39.70	ng	99
81) 3,3'-Dichlorobenzidine	21.55	252	837062	42.76	ng #	98
82) Chrysene	21.71	228	2093833	39.65	ng	100
83) Bis(2-ethylhexyl)phthalate	21.55	149	1368079	41.97	ng	97
84) Di-n-octyl phthalate	22.75	149	2340656	43.48	ng #	84
85) Indeno(1,2,3-cd)pyrene	28.52	276	2645739	42.53	ng #	87
87) Benzo(b)fluoranthene	23.84	252	2276772	39.58	ng #	96
88) Benzo(k)fluoranthene	23.92	252	2214197	39.45	ng #	94
89) Benzo(a)pyrene	24.71	252	2198798	40.36	ng #	95
90) Dibenzo(a,h)anthracene	28.58	278	2238588	40.90	ng #	94
91) Benzo(g,h,i)perylene	29.66	276	2245669	40.93	ng #	87

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

