

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030116\
 Data File : BG021191.D
 Acq On : 1 Mar 2016 19:45
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Mar 02 00:03:44 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 29 17:40:25 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	8634	20.00	ng	-0.01
21) Naphthalene-d8	10.96	136	45424	20.00	ng	0.00
38) Acenaphthene-d10	14.75	164	29431	20.00	ng	0.00
63) Phenanthrene-d10	17.49	188	69560	20.00	ng	0.00
75) Chrysene-d12	21.75	240	69936	20.00	ng	-0.01
86) Perylene-d12	25.02	264	65089	20.00	ng	-0.03

System Monitoring Compounds						
5) 2-Fluorophenol	5.69	112	45378	83.55	ng	-0.01
7) Phenol-d6	7.29	99	73784	82.62	ng	0.00
23) Nitrobenzene-d5	9.31	82	87456	81.47	ng	-0.01
41) 2,4,6-Tribromophenol	16.23	330	25224	82.38	ng	0.00
44) 2-Fluorobiphenyl	13.39	172	166181	80.18	ng	0.00
78) Terphenyl-d14	20.08	244	239194	82.86	ng	-0.01

Target Compounds						Qvalue
2) 1,4-Dioxane	3.58	88	10234	41.96	ng	# 89
3) Pyridine	3.98	79	33446	43.38	ng	94
4) n-Nitrosodimethylamine	3.90	42	15370	39.57	ng	98
6) Aniline	7.47	93	48121	40.65	ng	# 87
8) 2-Chlorophenol	7.71	128	27666	41.55	ng	84
9) Benzaldehyde	7.28	77	21857	39.18	ng	# 80
10) Phenol	7.32	94	39650	41.65	ng	82
11) bis(2-Chloroethyl)ether	7.56	93	29819	40.64	ng	93
12) 1,3-Dichlorobenzene	8.03	146	27915	40.89	ng	# 89
13) 1,4-Dichlorobenzene	8.18	146	28712	40.85	ng	93
14) 1,2-Dichlorobenzene	8.50	146	27591	41.16	ng	96
15) Benzyl Alcohol	8.37	79	34372	42.15	ng	# 85
16) 2,2'-oxybis(1-Chloropropan	8.67	45	51575	40.71	ng	98
17) 2-Methylphenol	8.57	107	27012	41.01	ng	# 89
18) Hexachloroethane	9.23	117	11963	40.64	ng	97
19) n-Nitroso-di-n-propylamine	8.94	70	34199	41.45	ng	# 92
20) 3+4-Methylphenols	8.90	107	38323	41.08	ng	93
22) Acetophenone	8.97	105	54946	40.65	ng	# 92
24) Nitrobenzene	9.35	77	47368	40.43	ng	# 87
25) Isophorone	9.87	82	97944	42.10	ng	94
26) 2-Nitrophenol	10.07	139	18338	42.97	ng	# 71
27) 2,4-Dimethylphenol	10.11	122	31036	40.54	ng	88
28) bis(2-Chloroethoxy)methane	10.35	93	43012	39.82	ng	99
29) 2,4-Dichlorophenol	10.59	162	27253	39.51	ng	96
30) 1,2,4-Trichlorobenzene	10.82	180	27120	38.98	ng	95
31) Naphthalene	11.01	128	94654	39.97	ng	99
32) Benzoic acid	10.22	122	18988	31.44	ng	87
33) 4-Chloroaniline	11.11	127	43704	40.01	ng	# 88
34) Hexachlorobutadiene	11.28	225	17061	40.29	ng	97
35) Caprolactam	11.88	113	14502	43.12	ng	# 80
36) 4-Chloro-3-methylphenol	12.21	107	42925	41.02	ng	85
37) 2-Methylnaphthalene	12.60	142	66975	39.72	ng	# 95
39) 1,2,4,5-Tetrachlorobenzene	12.96	216	35984	40.72	ng	# 98
40) Hexachlorocyclopentadiene	12.94	237	21070	43.48	ng	96

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42) 2,4,6-Trichlorophenol	13.19	196	27272	41.09	nq	94
43) 2,4,5-Trichlorophenol	13.26	196	28803	40.58	nq #	89
45) 1,1'-Biphenyl	13.59	154	98339	40.63	nq	96
46) 2-Chloronaphthalene	13.64	162	74101	40.78	nq	98
47) 2-Nitroaniline	13.83	65	36858	41.62	nq #	67
48) Acenaphthylene	14.48	152	125874	40.92	nq	98
49) Dimethylphthalate	14.20	163	106603	40.64	nq	97
50) 2,6-Dinitrotoluene	14.33	165	22792	41.53	nq #	72
51) Acenaphthene	14.82	154	79484	40.70	nq	99
52) 3-Nitroaniline	14.65	138	25336	41.61	nq #	66
53) 2,4-Dinitrophenol	14.85	184	9183	34.04	nq #	70
54) Dibenzofuran	15.15	168	104331	40.01	nq	91
55) 4-Nitrophenol	14.94	139	21260	42.10	nq #	53
56) 2,4-Dinitrotoluene	15.11	165	33203	42.00	nq #	77
57) Fluorene	15.80	166	98948	39.82	nq	98
58) 2,3,4,6-Tetrachlorophenol	15.37	232	23307	41.05	nq #	93
59) Diethylphthalate	15.56	149	117749	40.62	nq	98
60) 4-Chlorophenyl-phenylether	15.78	204	50033	40.33	nq	96
61) 4-Nitroaniline	15.81	138	26785	41.19	nq #	59
62) Azobenzene	16.08	77	125832	40.41	nq	93
64) 4,6-Dinitro-2-methylphenol	15.86	198	16380	35.39	nq	90
65) n-Nitrosodiphenylamine	16.00	169	87913	41.10	nq	93
66) 4-Bromophenyl-phenylether	16.68	248	29954	41.20	nq #	83
67) Hexachlorobenzene	16.79	284	29038	41.15	nq #	85
68) Atrazine	16.94	200	31155	42.54	nq	97
69) Pentachlorophenol	17.13	266	17630	38.25	nq	89
70) Phenanthrene	17.53	178	153274	40.83	nq	99
71) Anthracene	17.62	178	155449	40.47	nq	98
72) Carbazole	17.89	167	138942	40.79	nq	97
73) Di-n-butylphthalate	18.43	149	193386	40.38	nq #	98
74) Fluoranthene	19.53	202	173407	39.32	nq	100
76) Benzidine	19.70	184	90321	46.16	nq	99
77) Pyrene	19.89	202	174034	41.71	nq	98
79) Butylbenzylphthalate	20.76	149	88140	41.40	nq #	85
80) Benzo(a)anthracene	21.73	228	167028	40.34	nq	98
81) 3,3'-Dichlorobenzidine	21.64	252	65902	42.06	nq	94
82) Chrysene	21.80	228	158086	40.29	nq	97
83) Bis(2-ethylhexyl)phthalate	21.62	149	128157	41.07	nq	98
84) Di-n-octyl phthalate	22.85	149	216181	41.59	nq	98
85) Indeno(1,2,3-cd)pyrene	28.76	276	181485	40.34	nq #	100
87) Benzo(b)fluoranthene	23.98	252	165306	40.38	nq	99
88) Benzo(k)fluoranthene	24.04	252	159321	38.86	nq	98
89) Benzo(a)pyrene	24.87	252	155863	39.38	nq	96
90) Dibenzo(a,h)anthracene	28.82	278	153459	39.19	nq #	97
91) Benzo(g,h,i)perylene	29.92	276	146286	38.67	nq	93

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

