

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061417\
 Data File : BG027421.D
 Acq On : 14 Jun 2017 17:29
 Operator : SJ/MA
 Sample : SSTDICC060
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC060

Quant Time: Jun 14 18:23:46 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG061417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 14 16:47:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.53	152	42266	20.00	ng	0.00
21) Naphthalene-d8	11.35	136	223060	20.00	ng	0.00
38) Acenaphthene-d10	15.11	164	149512	20.00	ng	0.00
63) Phenanthrene-d10	17.86	188	336522	20.00	ng	0.00
75) Chrysene-d12	22.24	240	414213	20.00	ng	0.00
86) Perylene-d12	25.90	264	381008	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	6.11	112	372889	134.63	ng	0.00
7) Phenol-d6	7.67	99	584785	143.83	ng	0.00
23) Nitrobenzene-d5	9.69	82	545990	153.94	ng	0.00
41) 2,4,6-Tribromophenol	16.60	330	287866	146.21	ng	0.00
44) 2-Fluorobiphenyl	13.73	172	1234881	115.55	ng	0.00
78) Terphenyl-d14	20.44	244	1907640	117.49	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	4.12	88	75831	66.03	ng	# 86
3) Pyridine	4.50	79	242405	68.47	ng	# 81
4) n-Nitrosodimethylamine	4.41	42	97702	74.53	ng	# 54
6) Aniline	7.85	93	355945	69.52	ng	97
8) 2-Chlorophenol	8.09	128	195849	66.96	ng	97
9) Benzaldehyde	7.67	77	175943	62.59	ng	88
10) Phenol	7.69	94	295894	71.17	ng	79
11) bis(2-Chloroethyl)ether	7.94	93	229717	67.35	ng	91
12) 1,3-Dichlorobenzene	8.42	146	203151	61.58	ng	94
13) 1,4-Dichlorobenzene	8.56	146	212409	62.74	ng	96
14) 1,2-Dichlorobenzene	8.89	146	206129	62.42	ng	96
15) Benzyl Alcohol	8.75	79	212755	74.27	ng	# 87
16) 2,2'-oxybis(1-Chloropropan	9.03	45	391501	80.22	ng	95
17) 2-Methylphenol	8.94	107	202898	69.03	ng	# 89
18) Hexachloroethane	9.62	117	79239	69.48	ng	86
19) n-Nitroso-di-n-propylamine	9.32	70	220977	77.69	ng	# 86
20) 3+4-Methylphenols	9.27	107	294821	72.42	ng	89
22) Acetophenone	9.35	105	365700	64.11	ng	# 86
24) Nitrobenzene	9.73	77	300596	74.39	ng	# 89
25) Isophorone	10.26	82	595970	71.46	ng	# 97
26) 2-Nitrophenol	10.45	139	119831	79.68	ng	# 82
27) 2,4-Dimethylphenol	10.49	122	234111	65.29	ng	93
28) bis(2-Chloroethoxy)methane	10.72	93	352290	65.24	ng	99
29) 2,4-Dichlorophenol	10.98	162	209154	63.89	ng	97
30) 1,2,4-Trichlorobenzene	11.20	180	213774	58.97	ng	99
31) Naphthalene	11.40	128	735006	60.29	ng	99
32) Benzoic acid	10.61	122	161395	73.72	ng	# 87
33) 4-Chloroaniline	11.50	127	320939	63.18	ng	92
34) Hexachlorobutadiene	11.67	225	133283	59.83	ng	98
35) Caprolactam	12.28	113	94114	83.82	ng	90
36) 4-Chloro-3-methylphenol	12.57	107	292885	71.38	ng	96
37) 2-Methylnaphthalene	12.97	142	538365	60.54	ng	94
39) 1,2,4,5-Tetrachlorobenzene	13.32	216	271900	58.38	ng	99
40) Hexachlorocyclopentadiene	13.30	237	156138	62.98	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.55	196	188121	66.59	nq	97
43) 2,4,5-Trichlorophenol	13.62	196	212188	67.12	nq	93
45) 1,1'-Biphenyl	13.95	154	719459	58.95	nq	97
46) 2-Chloronaphthalene	14.00	162	544025	59.46	nq	99
47) 2-Nitroaniline	14.19	65	227284	96.55	nq	88
48) Acenaphthylene	14.84	152	955819	61.73	nq	98
49) Dimethylphthalate	14.55	163	742236	62.58	nq	98
50) 2,6-Dinitrotoluene	14.67	165	165899	75.09	nq	# 83
51) Acenaphthene	15.17	154	640350	63.58	nq	99
52) 3-Nitroaniline	15.01	138	173442	74.48	nq	96
53) 2,4-Dinitrophenol	15.21	184	91691	105.91	nq	96
54) Dibenzofuran	15.51	168	849199	60.33	nq	94
55) 4-Nitrophenol	15.29	139	145039	75.00	nq	# 62
56) 2,4-Dinitrotoluene	15.46	165	225944	81.09	nq	# 93
57) Fluorene	16.16	166	765252	61.21	nq	98
58) 2,3,4,6-Tetrachlorophenol	15.73	232	187612	66.48	nq	98
59) Diethylphthalate	15.90	149	758591	64.12	nq	98
60) 4-Chlorophenyl-phenylether	16.13	204	390796	62.06	nq	97
61) 4-Nitroaniline	16.17	138	184090	78.18	nq	# 71
62) Azobenzene	16.43	77	779425	71.09	nq	89
64) 4,6-Dinitro-2-methylphenol	16.22	198	130195	94.73	nq	82
65) n-Nitrosodiphenylamine	16.35	169	657322	62.30	nq	99
66) 4-Bromophenyl-phenylether	17.03	248	253621	63.68	nq	95
67) Hexachlorobenzene	17.16	284	280024	65.36	nq	91
68) Atrazine	17.29	200	230611	65.37	nq	96
69) Pentachlorophenol	17.50	266	166214	66.19	nq	97
70) Phenanthrene	17.91	178	1138130	60.11	nq	97
71) Anthracene	17.99	178	1162429	61.28	nq	99
72) Carbazole	18.26	167	1123792	62.81	nq	97
73) Di-n-butylphthalate	18.78	149	1236519	71.28	nq	# 95
74) Fluoranthene	19.89	202	1337380	66.12	nq	93
76) Benzidine	20.06	184	596391	47.80	nq	98
77) Pyrene	20.26	202	1372628	54.85	nq	98
79) Butylbenzylphthalate	21.14	149	601365	68.00	nq	95
80) Benzo(a)anthracene	22.22	228	1449419	61.36	nq	96
81) 3,3'-Dichlorobenzidine	22.11	252	574344	62.50	nq	# 99
82) Chrysene	22.30	228	1370151	60.42	nq	97
83) Bis(2-ethylhexyl)phthalate	22.07	149	868494	64.99	nq	99
84) Di-n-octyl phthalate	23.44	149	1452589	65.63	nq	# 92
85) Indeno(1,2,3-cd)pyrene	30.12	276	1756850	63.96	nq	# 93
87) Benzo(b)fluoranthene	24.73	252	1533813	64.91	nq	# 97
88) Benzo(k)fluoranthene	24.80	252	1449105	62.51	nq	# 93
89) Benzo(a)pyrene	25.73	252	1423492	63.99	nq	# 94
90) Dibenzo(a,h)anthracene	30.20	278	1534937	65.92	nq	# 93
91) Benzo(g,h,i)perylene	31.46	276	1453350	64.45	nq	# 90

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

