

Data Path : Z:\HPCHEM1\BNA G\DATA\BG071916\
 Data File : BG023172.D
 Acq On : 19 Jul 2016 12:07
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Jul 19 15:46:50 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG070816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 19 15:46:37 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	104557	20.00	ng	0.00
21) Naphthalene-d8	10.98	136	498938	20.00	ng	0.00
38) Acenaphthene-d10	14.80	164	365074	20.00	ng	0.00
63) Phenanthrene-d10	17.56	188	872820	20.00	ng	0.00
75) Chrysene-d12	21.88	240	843612	20.00	ng	0.00
86) Perylene-d12	25.27	264	826811	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.68	112	495400	77.49	ng	0.00
7) Phenol-d6	7.30	99	810856	80.98	ng	0.00
23) Nitrobenzene-d5	9.32	82	941518	80.80	ng	0.00
41) 2,4,6-Tribromophenol	16.30	330	407053	62.04	ng	0.00
44) 2-Fluorobiphenyl	13.41	172	1841423	79.45	ng	0.00
78) Terphenyl-d14	20.16	244	2416782	95.03	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.54	88	91355	36.76	ng	91
3) Pyridine	3.95	79	344124	40.47	ng	96
4) n-Nitrosodimethylamine	3.86	42	157717	43.23	ng	# 92
6) Aniline	7.46	93	577908	41.65	ng	95
8) 2-Chlorophenol	7.70	128	283898	41.73	ng	99
9) Benzaldehyde	7.27	77	270161	40.38	ng	94
10) Phenol	7.32	94	423262	41.08	ng	92
11) bis(2-Chloroethyl)ether	7.55	93	335731	41.12	ng	92
12) 1,3-Dichlorobenzene	8.03	146	333179	40.01	ng	98
13) 1,4-Dichlorobenzene	8.18	146	333743	40.03	ng	94
14) 1,2-Dichlorobenzene	8.50	146	330033	39.82	ng	92
15) Benzyl Alcohol	8.38	79	380859	41.53	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.67	45	472691	47.39	ng	93
17) 2-Methylphenol	8.58	107	272603	40.65	ng	96
18) Hexachloroethane	9.23	117	141148	40.11	ng	94
19) n-Nitroso-di-n-propylamine	8.95	70	383953	42.54	ng	96
20) 3+4-Methylphenols	8.92	107	395067	42.24	ng	96
22) Acetophenone	8.97	105	596575	39.84	ng	# 95
24) Nitrobenzene	9.36	77	497804	40.20	ng	95
25) Isophorone	9.88	82	938566	40.05	ng	# 94
26) 2-Nitrophenol	10.08	139	203096	41.80	ng	91
27) 2,4-Dimethylphenol	10.13	122	332249	39.56	ng	94
28) bis(2-Chloroethoxy)methane	10.36	93	540662	41.37	ng	99
29) 2,4-Dichlorophenol	10.62	162	340486	38.01	ng	96
30) 1,2,4-Trichlorobenzene	10.83	180	385774	37.59	ng	97
31) Naphthalene	11.02	128	1016613	40.03	ng	99
32) Benzoic acid	10.28	122	275134	36.01	ng	94
33) 4-Chloroaniline	11.13	127	497702	40.07	ng	97
34) Hexachlorobutadiene	11.30	225	290618	34.93	ng	98
35) Caprolactam	11.91	113	134524	36.50	ng	95
36) 4-Chloro-3-methylphenol	12.26	107	459731	37.53	ng	95
37) 2-Methylnaphthalene	12.63	142	798449	39.77	ng	96
39) 1,2,4,5-Tetrachlorobenzene	12.99	216	574037	36.49	ng	98
40) Hexachlorocyclopentadiene	12.97	237	431406	47.63	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.23	196	336861	37.25	nq	95
43) 2,4,5-Trichlorophenol	13.31	196	342727	36.91	nq	96
45) 1,1'-Biphenyl	13.63	154	1125325	41.08	nq	97
46) 2-Chloronaphthalene	13.68	162	911269	41.01	nq	96
47) 2-Nitroaniline	13.88	65	402578	42.46	nq	97
48) Acenaphthylene	14.52	152	1346059	40.38	nq	98
49) Dimethylphthalate	14.25	163	1147213	38.47	nq	100
50) 2,6-Dinitrotoluene	14.37	165	263586	39.33	nq	88
51) Acenaphthene	14.87	154	870777	39.98	nq	98
52) 3-Nitroaniline	14.71	138	277174	41.47	nq	95
53) 2,4-Dinitrophenol	14.91	184	184100	40.03	nq	93
54) Dibenzofuran	15.20	168	1254058	38.46	nq	96
55) 4-Nitrophenol	15.02	139	251412	44.88	nq	93
56) 2,4-Dinitrotoluene	15.16	165	379988	38.89	nq	93
57) Fluorene	15.85	166	1092391	38.93	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.43	232	312619	33.65	nq	93
59) Diethylphthalate	15.61	149	1173421	38.67	nq	98
60) 4-Chlorophenyl-phenylether	15.83	204	618361	36.18	nq	99
61) 4-Nitroaniline	15.88	138	297455	41.13	nq	# 79
62) Azobenzene	16.13	77	1251358	43.12	nq	93
64) 4,6-Dinitro-2-methylphenol	15.93	198	267615	41.25	nq	94
65) n-Nitrosodiphenylamine	16.05	169	1038821	41.31	nq	96
66) 4-Bromophenyl-phenylether	16.73	248	398058	35.05	nq	99
67) Hexachlorobenzene	16.86	284	432090	35.00	nq	# 93
68) Atrazine	17.00	200	410069	36.76	nq	96
69) Pentachlorophenol	17.20	266	347378	43.44	nq	99
70) Phenanthrene	17.60	178	1729390	40.43	nq	98
71) Anthracene	17.69	178	1778931	41.07	nq	99
72) Carbazole	17.97	167	1553131	41.50	nq	99
73) Di-n-butylphthalate	18.51	149	1836939	44.13	nq	99
74) Fluoranthene	19.61	202	1968229	37.50	nq	97
76) Benzidine	19.79	184	1573745	65.07	nq	96
77) Pyrene	19.98	202	2019681	42.60	nq	99
79) Butylbenzylphthalate	20.85	149	876582	46.68	nq	98
80) Benzo(a)anthracene	21.86	228	1903619	40.33	nq	97
81) 3,3'-Dichlorobenzidine	21.77	252	796248	38.43	nq	# 96
82) Chrysene	21.93	228	1806501	40.80	nq	97
83) Bis(2-ethylhexyl)phthalate	21.73	149	1219768	46.78	nq	# 99
84) Di-n-octyl phthalate	23.01	149	2017577	46.40	nq	99
85) Indeno(1,2,3-cd)pyrene	29.17	276	2042932	36.19	nq	# 100
87) Benzo(b)fluoranthene	24.19	252	1955240	40.99	nq	# 98
88) Benzo(k)fluoranthene	24.26	252	1874002	41.40	nq	# 98
89) Benzo(a)pyrene	25.11	252	1864958	40.79	nq	# 97
90) Dibenzo(a,h)anthracene	29.23	278	1790776	39.46	nq	# 95
91) Benzo(g,h,i)perylene	30.39	276	1784364	39.26	nq	# 97

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

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