

Data Path : Z:\HPCHEM1\BNA G\DATA\BG011017\
 Data File : BG025451.D
 Acq On : 10 Jan 2017 10:35
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Jan 10 13:05:25 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 10 11:26:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.20	152	64464	20.00	ng	0.00
21) Naphthalene-d8	11.03	136	259641	20.00	ng	0.00
38) Acenaphthene-d10	14.82	164	215693	20.00	ng	0.00
63) Phenanthrene-d10	17.57	188	487311	20.00	ng	0.00
75) Chrysene-d12	21.86	240	674006	20.00	ng	0.00
86) Perylene-d12	25.27	264	737354	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.74	112	287032	80.93	ng	0.00
7) Phenol-d6	7.36	99	418751	83.84	ng	0.00
23) Nitrobenzene-d5	9.38	82	513702	87.40	ng	0.00
41) 2,4,6-Tribromophenol	16.32	330	264830	76.26	ng	0.00
44) 2-Fluorobiphenyl	13.45	172	1090808	81.88	ng	0.00
78) Terphenyl-d14	20.15	244	1983777	78.04	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.58	88	62910	41.57	ng	95
3) Pyridine	3.99	79	189075	43.10	ng	94
4) n-Nitrosodimethylamine	3.91	42	105914	40.68	ng	# 99
6) Aniline	7.52	93	284510	42.04	ng	97
8) 2-Chlorophenol	7.76	128	159078	39.76	ng	97
9) Benzaldehyde	7.34	77	132563	36.27	ng	98
10) Phenol	7.39	94	232280	41.95	ng	93
11) bis(2-Chloroethyl)ether	7.61	93	176754	41.43	ng	89
12) 1,3-Dichlorobenzene	8.09	146	200166	41.34	ng	98
13) 1,4-Dichlorobenzene	8.24	146	203767	40.61	ng	96
14) 1,2-Dichlorobenzene	8.56	146	188530	39.37	ng	94
15) Benzyl Alcohol	8.44	79	182658	38.31	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.72	45	332822	42.12	ng	96
17) 2-Methylphenol	8.65	107	155385	42.73	ng	98
18) Hexachloroethane	9.28	117	77822	40.39	ng	87
19) n-Nitroso-di-n-propylamine	9.00	70	173297	41.08	ng	98
20) 3+4-Methylphenols	8.98	107	215330	42.79	ng	95
22) Acetophenone	9.03	105	306047	42.42	ng	# 93
24) Nitrobenzene	9.42	77	278704	43.24	ng	98
25) Isophorone	9.94	82	464515	43.65	ng	97
26) 2-Nitrophenol	10.14	139	102178	41.44	ng	91
27) 2,4-Dimethylphenol	10.18	122	175092	43.20	ng	94
28) bis(2-Chloroethoxy)methane	10.41	93	229787	41.58	ng	98
29) 2,4-Dichlorophenol	10.68	162	190881	44.01	ng	97
30) 1,2,4-Trichlorobenzene	10.88	180	215682	41.16	ng	97
31) Naphthalene	11.08	128	550592	42.46	ng	98
32) Benzoic acid	10.35	122	126602	38.84	ng	87
33) 4-Chloroaniline	11.20	127	236277	43.33	ng	95
34) Hexachlorobutadiene	11.33	225	185979	43.48	ng	99
35) Caprolactam	11.97	113	64492	39.57	ng	# 89
36) 4-Chloro-3-methylphenol	12.31	107	224403	41.91	ng	91
37) 2-Methylnaphthalene	12.67	142	419322	43.64	ng	95
39) 1,2,4,5-Tetrachlorobenzene	13.03	216	288771	40.55	ng	# 97
40) Hexachlorocyclopentadiene	12.99	237	94451	40.96	ng	91

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42) 2,4,6-Trichlorophenol	13.27	196	181045	40.38	nq	95
43) 2,4,5-Trichlorophenol	13.36	196	177345	37.78	nq	98
45) 1,1'-Biphenyl	13.66	154	595174	41.02	nq	99
46) 2-Chloronaphthalene	13.71	162	465241	41.05	nq	96
47) 2-Nitroaniline	13.93	65	209576	43.81	nq	95
48) Acenaphthylene	14.55	152	709433	40.38	nq	98
49) Dimethylphthalate	14.27	163	634324	40.34	nq	98
50) 2,6-Dinitrotoluene	14.41	165	138147	40.22	nq	97
51) Acenaphthene	14.89	154	467340	40.67	nq	96
52) 3-Nitroaniline	14.75	138	129592	39.24	nq	92
53) 2,4-Dinitrophenol	14.97	184	79763	38.24	nq	93
54) Dibenzofuran	15.22	168	756369	41.64	nq	96
55) 4-Nitrophenol	15.07	139	94319	38.24	nq	92
56) 2,4-Dinitrotoluene	15.21	165	200161	40.02	nq	# 89
57) Fluorene	15.87	166	612810	40.30	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.45	232	193063	38.40	nq	92
59) Diethylphthalate	15.62	149	665898	40.38	nq	98
60) 4-Chlorophenyl-phenylether	15.85	204	347764	40.35	nq	94
61) 4-Nitroaniline	15.92	138	136482	41.07	nq	# 84
62) Azobenzene	16.15	77	662335	41.65	nq	99
64) 4,6-Dinitro-2-methylphenol	15.97	198	142839	42.33	nq	82
65) n-Nitrosodiphenylamine	16.07	169	538139	40.86	nq	98
66) 4-Bromophenyl-phenylether	16.75	248	252857	42.06	nq	96
67) Hexachlorobenzene	16.87	284	279064	40.45	nq	91
68) Atrazine	17.01	200	204963	35.51	nq	95
69) Pentachlorophenol	17.23	266	133380	37.29	nq	95
70) Phenanthrene	17.62	178	1036004	41.26	nq	96
71) Anthracene	17.71	178	1056553	41.41	nq	99
72) Carbazole	17.99	167	937124	41.06	nq	99
73) Di-n-butylphthalate	18.49	149	1139015	40.56	nq	99
74) Fluoranthene	19.61	202	1366971	41.21	nq	94
76) Benzidine	19.79	184	569334	33.86	nq	99
77) Pyrene	19.98	202	1425019	40.45	nq	96
79) Butylbenzylphthalate	20.82	149	588000	41.57	nq	96
80) Benzo(a)anthracene	21.84	228	1563303	41.58	nq	98
81) 3,3'-Dichlorobenzidine	21.75	252	605804	41.49	nq	98
82) Chrysene	21.92	228	1486256	41.23	nq	98
83) Bis(2-ethylhexyl)phthalate	21.69	149	840317	42.69	nq	99
84) Di-n-octyl phthalate	22.94	149	1417810	43.39	nq	95
85) Indeno(1,2,3-cd)pyrene	29.20	276	2010524	43.87	nq	# 55
87) Benzo(b)fluoranthene	24.17	252	1667049	41.44	nq	98
88) Benzo(k)fluoranthene	24.24	252	1594341	41.04	nq	96
89) Benzo(a)pyrene	25.10	252	1620618	41.71	nq	98
90) Dibenzo(a,h)anthracene	29.23	278	1730124	42.01	nq	# 97
91) Benzo(g,h,i)perylene	30.42	276	1723169	42.11	nq	98

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

