

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM040517\
 Data File : BM009531.D
 Acq On : 04 Apr 2017 16:03
 Operator : SJ/MA
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Apr 05 02:14:53 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM032217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 31 12:10:26 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	136139	20.00	ng	-0.02
21) Naphthalene-d8	10.64	136	580563	20.00	ng	-0.02
38) Acenaphthene-d10	14.49	164	353003	20.00	ng	-0.02
63) Phenanthrene-d10	17.24	188	866067	20.00	ng	-0.01
75) Chrysene-d12	21.42	240	1147465	20.00	ng	-0.01
86) Perylene-d12	23.74	264	1046443	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	676786	82.86	ng	-0.02
7) Phenol-d6	7.00	99	987777	88.69	ng	-0.02
23) Nitrobenzene-d5	9.01	82	1264669	81.67	ng	-0.02
41) 2,4,6-Tribromophenol	15.98	330	386160	88.06	ng	-0.02
44) 2-Fluorobiphenyl	13.10	172	2339979	79.87	ng	-0.02
78) Terphenyl-d14	19.86	244	4492166	81.81	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.36	88	149768	36.47	ng	# 100
3) Pyridine	3.75	79	460079	39.64	ng	# 82
4) n-Nitrosodimethylamine	3.67	42	254810	41.24	ng	# 66
6) Aniline	7.18	93	645338	42.64	ng	# 89
8) 2-Chlorophenol	7.40	128	360167	43.49	ng	# 73
9) Benzaldehyde	6.99	77	329562	37.59	ng	# 79
10) Phenol	7.03	94	523342	43.36	ng	# 89
11) bis(2-Chloroethyl)ether	7.27	93	424130	42.88	ng	# 81
12) 1,3-Dichlorobenzene	7.73	146	404152	40.76	ng	# 82
13) 1,4-Dichlorobenzene	7.88	146	423324	41.32	ng	# 93
14) 1,2-Dichlorobenzene	8.19	146	414457	41.66	ng	# 92
15) Benzyl Alcohol	8.09	79	428511	44.01	ng	# 88
16) 2,2'-oxybis(1-Chloropropan	8.36	45	738889	42.98	ng	# 93
17) 2-Methylphenol	8.28	107	352713	44.67	ng	# 85
18) Hexachloroethane	8.92	117	178798	42.66	ng	# 90
19) n-Nitroso-di-n-propylamine	8.65	70	402816	44.86	ng	# 89
20) 3+4-Methylphenols	8.61	107	485489	45.22	ng	# 90
22) Acetophenone	8.67	105	658865	40.24	ng	# 89
24) Nitrobenzene	9.05	77	611651	40.32	ng	# 90
25) Isophorone	9.57	82	1003840	42.50	ng	# 90
26) 2-Nitrophenol	9.76	139	193473	42.29	ng	# 48
27) 2,4-Dimethylphenol	9.81	122	358069	42.78	ng	# 86
28) bis(2-Chloroethoxy)methane	10.05	93	597260	40.84	ng	# 99
29) 2,4-Dichlorophenol	10.29	162	367094	41.95	ng	# 94
30) 1,2,4-Trichlorobenzene	10.50	180	427303	39.36	ng	# 100
31) Naphthalene	10.69	128	1244066	40.37	ng	# 99
32) Benzoic acid	9.95	122	201085	34.76	ng	# 72
33) 4-Chloroaniline	10.81	127	509872	42.89	ng	# 82
34) Hexachlorobutadiene	10.96	225	286849	38.59	ng	# 99
35) Caprolactam	11.60	113	124893	46.75	ng	# 61
36) 4-Chloro-3-methylphenol	11.92	107	439181	44.36	ng	# 78
37) 2-Methylnaphthalene	12.30	142	879538	41.35	ng	# 93
39) 1,2,4,5-Tetrachlorobenzene	12.67	216	509174	38.21	ng	# 100
40) Hexachlorocyclopentadiene	12.64	237	267924	35.03	ng	# 94

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM040517\
 Data File : BM009531.D
 Acq On : 04 Apr 2017 16:03
 Operator : SJ/MA
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Apr 05 02:14:53 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM032217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 31 12:10:26 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.91	196	317008	41.60	ng	97
43) 2,4,5-Trichlorophenol	12.99	196	349659	41.12	ng #	87
45) 1,1'-Biphenyl	13.32	154	1187431	40.42	ng	95
46) 2-Chloronaphthalene	13.36	162	916824	40.18	ng	95
47) 2-Nitroaniline	13.57	65	382595	45.51	ng #	68
48) Acenaphthylene	14.21	152	1541801	41.48	ng	100
49) Dimethylphthalate	13.95	163	1134214	41.85	ng #	99
50) 2,6-Dinitrotoluene	14.07	165	260021	44.52	ng #	71
51) Acenaphthene	14.56	154	927614	41.37	ng	99
52) 3-Nitroaniline	14.40	138	266373	46.29	ng #	63
53) 2,4-Dinitrophenol	14.61	184	200970	53.00	ng	92
54) Dibenzofuran	14.89	168	1380201	41.62	ng	93
55) 4-Nitrophenol	14.70	139	207753	43.68	ng #	23
56) 2,4-Dinitrotoluene	14.86	165	360420	45.61	ng	96
57) Fluorene	15.54	166	1259784	42.26	ng	96
58) 2,3,4,6-Tetrachlorophenol	15.12	232	308878	42.95	ng #	100
59) Diethylphthalate	15.31	149	1189343	43.26	ng	100
60) 4-Chlorophenyl-phenylether	15.53	204	646958	41.56	ng	98
61) 4-Nitroaniline	15.57	138	292748	46.84	ng #	32
62) Azobenzene	15.82	77	1415060	44.09	ng	85
64) 4,6-Dinitro-2-methylphenol	15.62	198	217608	39.88	ng	91
65) n-Nitrosodiphenylamine	15.75	169	1044903	39.72	ng	99
66) 4-Bromophenyl-phenylether	16.43	248	390836	38.95	ng #	92
67) Hexachlorobenzene	16.54	284	425956	38.73	ng #	91
68) Atrazine	16.70	200	388396	39.26	ng	97
69) Pentachlorophenol	16.89	266	260114	38.71	ng	97
70) Phenanthrene	17.28	178	1991442	40.21	ng	99
71) Anthracene	17.37	178	2070514	41.29	ng	99
72) Carbazole	17.64	167	2045457	42.60	ng	99
73) Di-n-butylphthalate	18.19	149	2110876	43.04	ng #	98
74) Fluoranthene	19.30	202	2469841	42.10	ng	99
76) Benzidine	19.49	184	1152841	33.71	ng	100
77) Pyrene	19.66	202	2658745	40.85	ng	98
79) Butylbenzylphthalate	20.54	149	1018195	43.45	ng #	72
80) Benzo(a)anthracene	21.40	228	2746588	40.98	ng	99
81) 3,3'-Dichlorobenzidine	21.34	252	1056790	42.11	ng #	98
82) Chrysene	21.46	228	2604954	40.71	ng	98
83) Bis(2-ethylhexyl)phthalate	21.31	149	1516786	43.37	ng	98
84) Di-n-octyl phthalate	22.21	149	2594147	44.57	ng #	90
85) Indeno(1,2,3-cd)pyrene	26.11	276	2943542	42.18	ng #	100
87) Benzo(b)fluoranthene	23.04	252	2673346	40.30	ng #	94
88) Benzo(k)fluoranthene	23.09	252	2732920	42.84	ng	98
89) Benzo(a)pyrene	23.64	252	2583838	42.07	ng	98
90) Dibenzo(a,h)anthracene	26.12	278	2540636	42.22	ng #	95
91) Benzo(g,h,i)perylene	26.84	276	2456231	41.87	ng #	93

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

