

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102615\
 Data File : BF082522.D
 Acq On : 26 Oct 2015 20:22
 Operator : UM/IZ
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Quant Time: Oct 27 01:37:15 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 27 01:37:08 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	98539	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	383527	20.00	ng	0.00
38) Acenaphthene-d10	9.77	164	184474	20.00	ng	0.00
63) Phenanthrene-d10	11.26	188	354697	20.00	ng	0.00
75) Chrysene-d12	13.91	240	305938	20.00	ng	0.00
86) Perylene-d12	15.36	264	283372	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.31	112	486212	100.04	ng	0.00
7) Phenol-d6	6.39	99	632813	100.47	ng	0.01
23) Nitrobenzene-d5	7.30	82	583425	102.40	ng	0.00
41) 2,4,6-Tribromophenol	10.57	330	147143	86.64	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	1018411	92.68	ng	0.00
78) Terphenyl-d14	12.85	244	979809	85.53	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.23	88	111561	45.43	ng	96
3) Pyridine	2.90	79	300312	52.01	ng	98
4) n-Nitrosodimethylamine	2.89	42	132556	54.37	ng	100
6) Aniline	6.40	93	405641	49.91	ng	# 86
8) 2-Chlorophenol	6.51	128	293300	49.22	ng	98
9) Benzaldehyde	6.27	77	150919	46.90	ng	100
10) Phenol	6.40	94	373177	52.17	ng	98
11) bis(2-Chloroethyl)ether	6.47	93	277885	45.75	ng	99
12) 1,3-Dichlorobenzene	6.66	146	323016	46.76	ng	99
13) 1,4-Dichlorobenzene	6.74	146	334775	46.51	ng	99
14) 1,2-Dichlorobenzene	6.90	146	315752	46.20	ng	97
15) Benzyl Alcohol	6.89	79	224822	51.55	ng	# 89
16) 2,2'-oxybis(1-Chloropropan	7.02	45	379335	44.20	ng	97
17) 2-Methylphenol	7.01	107	225897	47.79	ng	98
18) Hexachloroethane	7.23	117	117036	47.23	ng	99
19) n-Nitroso-di-n-propylamine	7.15	70	198570	45.07	ng	98
20) 3+4-Methylphenols	7.17	107	279707	49.33	ng	97
22) Acetophenone	7.15	105	385958	50.16	ng	# 98
24) Nitrobenzene	7.33	77	298294	46.87	ng	95
25) Isophorone	7.57	82	547117	47.29	ng	99
26) 2-Nitrophenol	7.65	139	157741	51.72	ng	96
27) 2,4-Dimethylphenol	7.69	122	271288	54.19	ng	95
28) bis(2-Chloroethoxy)methane	7.78	93	340397	51.37	ng	99
29) 2,4-Dichlorophenol	7.89	162	245803	49.57	ng	98
30) 1,2,4-Trichlorobenzene	7.97	180	278427	48.61	ng	99
31) Naphthalene	8.05	128	801849	48.10	ng	99
32) Benzoic acid	7.85	122	149419m	46.63	ng	
33) 4-Chloroaniline	8.10	127	342667	49.86	ng	97
34) Hexachlorobutadiene	8.15	225	154205	43.73	ng	98
35) Caprolactam	8.50	113	76266	53.72	ng	# 78
36) 4-Chloro-3-methylphenol	8.59	107	238891	48.55	ng	96
37) 2-Methylnaphthalene	8.73	142	537399	46.70	ng	100
39) 1,2,4,5-Tetrachlorobenzene	8.90	216	279169	50.43	ng	99
40) Hexachlorocyclopentadiene	8.88	237	65801	31.98	ng	100

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42) 2,4,6-Trichlorophenol	9.02	196	169210	46.11	nq	99
43) 2,4,5-Trichlorophenol	9.06	196	186180	47.21	nq	96
45) 1,1'-Biphenyl	9.20	154	684246	47.75	nq	98
46) 2-Chloronaphthalene	9.22	162	530139	47.68	nq	98
47) 2-Nitroaniline	9.34	65	159096	51.97	nq	# 53
48) Acenaphthylene	9.63	152	786543	45.69	nq	100
49) Dimethylphthalate	9.51	163	598987	45.30	nq	100
50) 2,6-Dinitrotoluene	9.58	165	141822	51.22	nq	# 74
51) Acenaphthene	9.81	154	450813	43.97	nq	99
52) 3-Nitroaniline	9.75	138	158529	50.73	nq	89
53) 2,4-Dinitrophenol	9.86	184	59017m	46.98	nq	
54) Dibenzofuran	9.98	168	666099	46.95	nq	99
55) 4-Nitrophenol	9.93	139	59102	35.75	nq	99
56) 2,4-Dinitrotoluene	9.98	165	170783	50.15	nq	92
57) Fluorene	10.32	166	523388	44.93	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.10	232	141860	44.22	nq	97
59) Diethylphthalate	10.21	149	621011	50.53	nq	98
60) 4-Chlorophenyl-phenylether	10.32	204	273399	50.62	nq	95
61) 4-Nitroaniline	10.37	138	155529	51.20	nq	95
62) Azobenzene	10.48	77	588517	49.50	nq	93
64) 4,6-Dinitro-2-methylphenol	10.40	198	107718	54.51	nq	# 59
65) n-Nitrosodiphenylamine	10.45	169	530546	50.30	nq	99
66) 4-Bromophenyl-phenylether	10.80	248	154134	43.95	nq	96
67) Hexachlorobenzene	10.87	284	175219	45.98	nq	# 91
68) Atrazine	10.97	200	153374	44.96	nq	99
69) Pentachlorophenol	11.07	266	65308	35.26	nq	96
70) Phenanthrene	11.28	178	820501	47.22	nq	100
71) Anthracene	11.34	178	906079	50.47	nq	99
72) Carbazole	11.50	167	767701	46.85	nq	99
73) Di-n-butylphthalate	11.83	149	913992	46.02	nq	# 97
74) Fluoranthene	12.47	202	851751	46.16	nq	99
76) Benzidine	12.61	184	416721	47.39	nq	98
77) Pyrene	12.70	202	863222	47.40	nq	99
79) Butylbenzylphthalate	13.33	149	419331	50.89	nq	# 77
80) Benzo(a)anthracene	13.90	228	788254	49.42	nq	100
81) 3,3'-Dichlorobenzidine	13.86	252	281621	46.60	nq	# 96
82) Chrysene	13.94	228	830254	49.42	nq	100
83) Bis(2-ethylhexyl)phthalate	13.89	149	522081	44.78	nq	# 96
84) Di-n-octyl phthalate	14.53	149	885460	44.18	nq	99
85) Indeno(1,2,3-cd)pyrene	16.73	276	661114	48.72	nq	100
87) Benzo(b)fluoranthene	14.96	252	692769m	41.20	nq	
88) Benzo(k)fluoranthene	15.00	252	755019m	51.49	nq	
89) Benzo(a)pyrene	15.30	252	669170	45.15	nq	# 95
90) Dibenzo(a,h)anthracene	16.74	278	549393	45.11	nq	99
91) Benzo(g,h,i)perylene	17.14	276	534637	45.05	nq	99

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

