

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072517\
 Data File : BF097021.D
 Acq On : 25 Jul 2017 11:37
 Operator : SJ/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Quant Time: Jul 25 13:21:24 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 13:17:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.50	152	141444	20.00	ng	0.00
21) Naphthalene-d8	7.79	136	589479	20.00	ng	0.00
38) Acenaphthene-d10	9.53	164	248178	20.00	ng	0.00
63) Phenanthrene-d10	11.00	188	400854	20.00	ng	0.00
75) Chrysene-d12	13.63	240	235182	20.00	ng	0.00
86) Perylene-d12	14.97	264	241086	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.10	112	876915	93.22	ng	0.00
7) Phenol-d6	6.17	99	996681	88.29	ng	0.00
23) Nitrobenzene-d5	7.08	82	975018	102.45	ng	0.00
41) 2,4,6-Tribromophenol	10.32	330	185481	87.56	ng	0.00
44) 2-Fluorobiphenyl	8.87	172	1390807	88.54	ng	0.00
78) Terphenyl-d14	12.59	244	1154866	85.16	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.05	88	198773	41.065	ng	# 73
3) Pyridine	2.66	79	630589	61.988	ng	# 64
4) n-Nitrosodimethylamine	2.63	42	336126	59.085	ng	# 82
6) Aniline	6.17	93	629901	45.299	ng	# 71
8) 2-Chlorophenol	6.29	128	470214	46.333	ng	# 92
9) Benzaldehyde	6.05	77	258852	39.672	ng	# 99
10) Phenol	6.19	94	602621	47.222	ng	# 96
11) bis(2-Chloroethyl)ether	6.25	93	480407	50.060	ng	# 89
12) 1,3-Dichlorobenzene	6.44	146	525468	48.711	ng	# 97
13) 1,4-Dichlorobenzene	6.52	146	519168	47.523	ng	# 96
14) 1,2-Dichlorobenzene	6.67	146	427520	43.025	ng	# 98
15) Benzyl Alcohol	6.66	79	310435	42.003	ng	# 96
16) 2,2'-oxybis(1-Chloropropan	6.80	45	978045	49.384	ng	# 95
17) 2-Methylphenol	6.79	107	370669	46.969	ng	# 90
18) Hexachloroethane	7.01	117	190154	49.088	ng	# 80
19) n-Nitroso-di-n-propylamine	6.94	70	329474	48.445	ng	# 84
20) 3+4-Methylphenols	6.95	107	481994	50.331	ng	# 64
22) Acetophenone	6.93	105	672788	51.065	ng	# 98
24) Nitrobenzene	7.10	77	511673	51.986	ng	# 92
25) Isophorone	7.34	82	952774	53.817	ng	# 97
26) 2-Nitrophenol	7.41	139	283797	51.855	ng	# 86
27) 2,4-Dimethylphenol	7.47	122	444540	49.373	ng	# 98
28) bis(2-Chloroethoxy)methane	7.56	93	624318	52.187	ng	# 99
29) 2,4-Dichlorophenol	7.66	162	383951	47.111	ng	# 98
30) 1,2,4-Trichlorobenzene	7.73	180	366548	47.303	ng	# 97
31) Naphthalene	7.81	128	1307259	46.813	ng	# 100
32) Benzoic acid	7.65	122	375529	65.257	ng	# 92
33) 4-Chloroaniline	7.87	127	536485	48.490	ng	# 98
34) Hexachlorobutadiene	7.93	225	201335	48.671	ng	# 99
35) Caprolactam	8.26	113	131628	50.405	ng	# 58
36) 4-Chloro-3-methylphenol	8.37	107	436043	49.758	ng	# 90
37) 2-Methylnaphthalene	8.50	142	834167	46.855	ng	# 98
39) 1,2,4,5-Tetrachlorobenzene	8.67	216	323025	48.149	ng	# 99
40) Hexachlorocyclopentadiene	8.66	237	114379	51.298	ng	# 99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.79	196	236705	47.832	ng	99
43) 2,4,5-Trichlorophenol	8.83	196	239738	48.035	ng	99
45) 1,1'-Biphenyl	8.97	154	978641	45.978	ng	98
46) 2-Chloronaphthalene	8.99	162	728817	46.501	ng	# 92
47) 2-Nitroaniline	9.09	65	277045	51.424	ng	97
48) Acenaphthylene	9.40	152	1097259	43.939	ng	99
49) Dimethylphthalate	9.28	163	896366	47.967	ng	99
50) 2,6-Dinitrotoluene	9.33	165	199385	49.260	ng	# 76
51) Acenaphthene	9.57	154	657145	44.545	ng	96
52) 3-Nitroaniline	9.50	138	237133	49.454	ng	90
53) 2,4-Dinitrophenol	9.62	184	100934	59.279	ng	# 72
54) Dibenzofuran	9.74	168	877328	42.439	ng	98
55) 4-Nitrophenol	9.69	139	153064	43.694	ng	# 63
56) 2,4-Dinitrotoluene	9.74	165	219201	42.145	ng	# 51
57) Fluorene	10.08	166	674394	44.146	ng	97
58) 2,3,4,6-Tetrachlorophenol	9.86	232	167093	48.274	ng	97
59) Diethylphthalate	9.97	149	841526	46.259	ng	99
60) 4-Chlorophenyl-phenylether	10.07	204	274390	42.720	ng	100
61) 4-Nitroaniline	10.12	138	238029	52.863	ng	90
62) Azobenzene	10.23	77	808119	50.355	ng	93
64) 4,6-Dinitro-2-methylphenol	10.15	198	141229	56.115	ng	90
65) n-Nitrosodiphenylamine	10.20	169	673407	46.996	ng	97
66) 4-Bromophenyl-phenylether	10.56	248	196735	48.071	ng	97
67) Hexachlorobenzene	10.63	284	216668	47.534	ng	# 88
68) Atrazine	10.73	200	189848	46.505	ng	95
69) Pentachlorophenol	10.83	266	101151	45.936	ng	97
70) Phenanthrene	11.03	178	1001231	45.937	ng	99
71) Anthracene	11.09	178	1021007	46.331	ng	99
72) Carbazole	11.25	167	1028383	48.189	ng	98
73) Di-n-butylphthalate	11.59	149	1289556	48.521	ng	98
74) Fluoranthene	12.21	202	971112	46.548	ng	98
76) Benzidine	12.33	184	423135	54.919	ng	98
77) Pyrene	12.43	202	1024539	48.001	ng	98
79) Butylbenzylphthalate	13.07	149	542989	53.543	ng	# 78
80) Benzo(a)anthracene	13.62	228	807346	55.095	ng	99
81) 3,3'-Dichlorobenzidine	13.59	252	300801	57.084	ng	# 96
82) Chrysene	13.66	228	771917	53.509	ng	99
83) Bis(2-ethylhexyl)phthalate	13.63	149	671671	54.386	ng	100
84) Di-n-octyl phthalate	14.24	149	1221356	59.827	ng	# 93
85) Indeno(1,2,3-cd)pyrene	16.21	276	640266	48.902	ng	98
87) Benzo(b)fluoranthene	14.62	252	666817	45.860	ng	97
88) Benzo(k)fluoranthene	14.64	252	669700	48.639	ng	96
89) Benzo(a)pyrene	14.93	252	647779	48.574	ng	99
90) Dibenzo(a,h)anthracene	16.22	278	505160	41.166	ng	97
91) Benzo(g,h,i)perylene	16.57	276	556604	42.358	ng	99

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

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