

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050818\
 Data File : BF105139.D
 Acq On : 8 May 2018 11:40
 Operator : JU/SJ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Quant Time: May 08 13:40:41 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 08 13:40:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	199009	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	813794	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	405755	20.00	ng	0.00
63) Phenanthrene-d10	11.31	188	796132	20.00	ng	0.00
75) Chrysene-d12	14.00	240	675288	20.00	ng	-0.01
86) Perylene-d12	15.55	264	627781	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.39	112	254955	19.59	ng	-0.01
7) Phenol-d6	6.40	99	302204	18.90	ng	-0.02
23) Nitrobenzene-d5	7.35	82	193769	15.55	ng	-0.01
41) 2,4,6-Tribromophenol	10.61	330	78116	18.56	ng	-0.01
44) 2-Fluorobiphenyl	9.15	172	624655	24.32	ng	-0.01
78) Terphenyl-d14	12.90	244	633098	21.37	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.51	88	65805	10.851	ng	92
3) Pyridine	3.24	79	171140	10.037	ng	# 84
4) n-Nitrosodimethylamine	3.18	42	85052	14.270	ng	83
6) Aniline	6.45	93	219614	9.885	ng	97
8) 2-Chlorophenol	6.57	128	135012	9.828	ng	93
9) Benzaldehyde	6.34	77	122918	15.403	ng	99
10) Phenol	6.42	94	168831	9.950	ng	94
11) bis(2-Chloroethyl)ether	6.52	93	142723	10.541	ng	91
12) 1,3-Dichlorobenzene	6.73	146	164880	10.595	ng	99
13) 1,4-Dichlorobenzene	6.81	146	167841	10.819	ng	98
14) 1,2-Dichlorobenzene	6.96	146	156609	10.894	ng	95
15) Benzyl Alcohol	6.93	79	107933	8.886	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.07	45	262168	14.418	ng	96
17) 2-Methylphenol	7.03	107	115351	9.660	ng	97
18) Hexachloroethane	7.30	117	50315	9.097	ng	97
19) n-Nitroso-di-n-propylamine	7.20	70	98179	9.395	ng	90
20) 3+4-Methylphenols	7.19	107	142297	9.608	ng	95
22) Acetophenone	7.20	105	212503	10.607	ng	# 94
24) Nitrobenzene	7.37	77	126071	9.162	ng	95
25) Isophorone	7.61	82	279763	10.615	ng	99
26) 2-Nitrophenol	7.69	139	17392	3.105	ng	# 86
27) 2,4-Dimethylphenol	7.72	122	121158	10.417	ng	97
28) bis(2-Chloroethoxy)methane	7.82	93	176120	10.930	ng	98
29) 2,4-Dichlorophenol	7.92	162	104528	9.419	ng	93
30) 1,2,4-Trichlorobenzene	8.02	180	138879	11.403	ng	96
31) Naphthalene	8.10	128	430034	10.612	ng	98
32) Benzoic acid	7.79	122	19235	2.073	ng	89
33) 4-Chloroaniline	8.14	127	175383	10.102	ng	95
34) Hexachlorobutadiene	8.21	225	79251	11.880	ng	99
35) Caprolactam	8.47	113	26791	6.920	ng	# 66
36) 4-Chloro-3-methylphenol	8.61	107	113261	9.518	ng	99
37) 2-Methylnaphthalene	8.79	142	301147	11.489	ng	96
39) 1,2,4,5-Tetrachlorobenzene	8.95	216	140178	12.069	ng	98
40) Hexachlorocyclopentadiene	8.94	237	53360	8.206	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.06	196	69312	8.596	nq	98
43) 2,4,5-Trichlorophenol	9.09	196	76848	8.517	nq	90
45) 1,1'-Biphenyl	9.25	154	368929	10.823	nq	98
46) 2-Chloronaphthalene	9.27	162	271140	10.071	nq	98
47) 2-Nitroaniline	9.37	65	34052	5.114	nq	95
48) Acenaphthylene	9.69	152	421836	9.986	nq	99
49) Dimethylphthalate	9.55	163	292023	9.127	nq	99
50) 2,6-Dinitrotoluene	9.61	165	46169	7.798	nq	96
51) Acenaphthene	9.86	154	264732	10.271	nq	99
52) 3-Nitroaniline	9.77	138	49377	7.124	nq	# 94
53) 2,4-Dinitrophenol	9.87	184	4728	2.627	nq	# 29
54) Dibenzofuran	10.03	168	399781	11.130	nq	97
55) 4-Nitrophenol	9.92	139	32872	6.037	nq	100
56) 2,4-Dinitrotoluene	10.00	165	51508	6.632	nq	# 86
57) Fluorene	10.37	166	309593	11.842	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.15	232	67632	10.047	nq	88
59) Diethylphthalate	10.25	149	290103	9.104	nq	96
60) 4-Chlorophenyl-phenylether	10.37	204	148189	12.728	nq	93
61) 4-Nitroaniline	10.38	138	50135	7.398	nq	93
62) Azobenzene	10.53	77	309113	10.153	nq	96
64) 4,6-Dinitro-2-methylphenol	10.41	198	8010	2.316	nq	91
65) n-Nitrosodiphenylamine	10.49	169	270757	9.809	nq	100
66) 4-Bromophenyl-phenylether	10.86	248	91760	10.836	nq	91
67) Hexachlorobenzene	10.92	284	102861	10.795	nq	93
68) Atrazine	11.01	200	77241	9.270	nq	99
69) Pentachlorophenol	11.11	266	36198	6.141	nq	98
70) Phenanthrene	11.33	178	472306	10.653	nq	98
71) Anthracene	11.39	178	478640	10.620	nq	99
72) Carbazole	11.54	167	426312	9.680	nq	99
73) Di-n-butylphthalate	11.88	149	399406	7.424	nq	98
74) Fluoranthene	12.52	202	488785	10.552	nq	98
76) Benzidine	12.65	184	212414	9.088	nq	100
77) Pyrene	12.75	202	504800	10.085	nq	99
79) Butylbenzylphthalate	13.40	149	72325	3.067	nq	99
80) Benzo(a)anthracene	13.99	228	447672	11.097	nq	99
81) 3,3'-Dichlorobenzidine	13.96	252	128396	8.536	nq	# 96
82) Chrysene	14.03	228	439146	10.598	nq	98
83) Bis(2-ethylhexyl)phthalate	14.00	149	144439	4.762	nq	100
84) Di-n-octyl phthalate	14.69	149	167983	3.314	nq	# 83
85) Indeno(1,2,3-cd)pyrene	16.99	276	307672	8.740	nq	94
87) Benzo(b)fluoranthene	15.12	252	366730	9.249	nq	98
88) Benzo(k)fluoranthene	15.15	252	419607	11.210	nq	# 96
89) Benzo(a)pyrene	15.49	252	350711	9.886	nq	97
90) Dibenzo(a,h)anthracene	17.02	278	254728	8.742	nq	96
91) Benzo(g,h,i)perylene	17.42	276	261992	9.281	nq	94

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

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