

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020218\
 Data File : BF102701.D
 Acq On : 2 Feb 2018 11:37
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 2/2/2018 5:41:48 PM

Quant Time: Feb 02 13:43:22 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 31 18:38:00 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	189923	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	781043	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	352435	20.00	ng	-0.01
63) Phenanthrene-d10	11.22	188	594697	20.00	ng	-0.01
75) Chrysene-d12	13.86	240	374971	20.00	ng	0.00
86) Perylene-d12	15.29	264	330726	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.32	112	1064015	84.95	ng	0.00
7) Phenol-d6	6.36	99	1193251	79.02	ng	0.00
23) Nitrobenzene-d5	7.27	82	1107134	81.46	ng	0.00
41) 2,4,6-Tribromophenol	10.53	330	236790	67.81	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	1919419	80.08	ng	-0.01
78) Terphenyl-d14	12.80	244	1376990	82.79	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.32	88	260787	43.820	ng	96
3) Pyridine	3.04	79	637488	40.989	ng	96
4) n-Nitrosodimethylamine	3.02	42	243783	39.982	ng	99
6) Aniline	6.37	93	724896	36.292	ng	# 59
8) 2-Chlorophenol	6.49	128	543797	41.415	ng	98
9) Benzaldehyde	6.26	77	310368	35.715	ng	97
10) Phenol	6.38	94	735895	44.637	ng	# 49
11) bis(2-Chloroethyl)ether	6.44	93	623232	49.704	ng	99
12) 1,3-Dichlorobenzene	6.63	146	622772	39.881	ng	99
13) 1,4-Dichlorobenzene	6.71	146	672008	42.960	ng	99
14) 1,2-Dichlorobenzene	6.86	146	545860	38.150	ng	98
15) Benzyl Alcohol	6.86	79	398056	37.360	ng	98
16) 2,2'-oxybis(1-Chloropropan	6.97	45	827905	39.368	ng	85
17) 2-Methylphenol	6.98	107	432034	37.886	ng	# 87
18) Hexachloroethane	7.19	117	215701	39.571	ng	94
19) n-Nitroso-di-n-propylamine	7.12	70	382023	41.338	ng	96
20) 3+4-Methylphenols	7.13	107	608653	45.382	ng	# 86
22) Acetophenone	7.12	105	779384	41.860	ng	# 96
24) Nitrobenzene	7.29	77	607096	43.648	ng	98
25) Isophorone	7.52	82	991480	40.920	ng	99
26) 2-Nitrophenol	7.60	139	300405	41.037	ng	93
27) 2,4-Dimethylphenol	7.65	122	415900	39.127	ng	98
28) bis(2-Chloroethoxy)methane	7.73	93	619414	41.384	ng	99
29) 2,4-Dichlorophenol	7.86	162	438933	40.539	ng	98
30) 1,2,4-Trichlorobenzene	7.92	180	443402	38.202	ng	99
31) Naphthalene	8.00	128	1550536	39.761	ng	99
32) Benzoic acid	7.82	122	258984	29.081	ng	99
33) 4-Chloroaniline	8.07	127	665723	40.597	ng	99
34) Hexachlorobutadiene	8.10	225	236501	38.151	ng	99
35) Caprolactam	8.45	113	149874	41.912	ng	# 83
36) 4-Chloro-3-methylphenol	8.56	107	565778	49.573	ng	96
37) 2-Methylnaphthalene	8.69	142	1150599	44.304	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.86	216	451373	42.569	ng	99
40) Hexachlorocyclopentadiene	8.83	237	91572	19.298	ng	99

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42) 2,4,6-Trichlorophenol	8.98	196	277779	39.134	nq	99
43) 2,4,5-Trichlorophenol	9.03	196	309802	38.764	nq	99
45) 1,1'-Biphenyl	9.15	154	1288761	40.896	nq	99
46) 2-Chloronaphthalene	9.18	162	966576	39.460	nq	99
47) 2-Nitroaniline	9.30	65	313986	41.115	nq	99
48) Acenaphthylene	9.59	152	1491632	39.087	nq	99
49) Dimethylphthalate	9.46	163	1077859	37.000	nq	100
50) 2,6-Dinitrotoluene	9.53	165	242910	38.676	nq #	87
51) Acenaphthene	9.76	154	873487	39.192	nq	99
52) 3-Nitroaniline	9.72	138	292181	39.325	nq	98
53) 2,4-Dinitrophenol	9.84	184	75150	29.132	nq #	25
54) Dibenzofuran	9.94	168	1183819	39.247	nq	94
55) 4-Nitrophenol	9.94	139	692092	141.114	nq #	18
56) 2,4-Dinitrotoluene	9.95	165	303026	39.234	nq #	92
57) Fluorene	10.28	166	955025	39.957	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.07	232	210393	36.870	nq	98
59) Diethylphthalate	10.16	149	1085233	38.336	nq	98
60) 4-Chlorophenyl-phenylether	10.27	204	413658	39.918	nq	97
61) 4-Nitroaniline	10.34	138	284772	38.410	nq	90
62) Azobenzene	10.43	77	1088960	42.016	nq	98
64) 4,6-Dinitro-2-methylphenol	10.36	198	141162	35.579	nq	94
65) n-Nitrosodiphenylamine	10.40	169	876188	40.141	nq	99
66) 4-Bromophenyl-phenylether	10.76	248	249243	38.933	nq	92
67) Hexachlorobenzene	10.83	284	257831	36.011	nq	95
68) Atrazine	10.93	200	237797	36.887	nq	98
69) Pentachlorophenol	11.04	266	124815m	33.185	nq	
70) Phenanthrene	11.24	178	1313655	37.907	nq	99
71) Anthracene	11.30	178	1365462	38.604	nq	99
72) Carbazole	11.46	167	1349357	39.103	nq	100
73) Di-n-butylphthalate	11.77	149	1636617	37.943	nq	100
74) Fluoranthene	12.43	202	1275497	36.093	nq	99
76) Benzidine	12.57	184	346892	27.538	nq	98
77) Pyrene	12.66	202	1288163	44.433	nq	99
79) Butylbenzylphthalate	13.27	149	632251	43.822	nq	99
80) Benzo(a)anthracene	13.85	228	926149	40.118	nq	100
81) 3,3'-Dichlorobenzidine	13.82	252	308115	36.383	nq	98
82) Chrysene	13.89	228	946428	40.371	nq	99
83) Bis(2-ethylhexyl)phthalate	13.83	149	758913	39.141	nq	99
84) Di-n-octyl phthalate	14.44	149	1195732	37.922	nq	99
85) Indeno(1,2,3-cd)pyrene	16.70	276	758690	39.186	nq	97
87) Benzo(b)fluoranthene	14.88	252	770285m	35.934	nq	
88) Benzo(k)fluoranthene	14.91	252	831285	41.046	nq	99
89) Benzo(a)pyrene	15.23	252	740710	38.313	nq	98
90) Dibenzo(a,h)anthracene	16.70	278	630588	40.266	nq #	95
91) Benzo(g,h,i)perylene	17.12	276	668283	43.218	nq	95

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