

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080115\
 Data File : BF080762.D
 Acq On : 1 Aug 2015 5:38
 Operator : TP/UM
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

MMDadoda
 8/3/2015 3:40:46 PM

Quant Time: Aug 03 00:53:24 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF073015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 31 01:45:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	171718	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	618589	20.00	ng	-0.01
38) Acenaphthene-d10	9.89	164	313573	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	546564	20.00	ng	0.00
75) Chrysene-d12	14.00	240	354240	20.00	ng	0.00
86) Perylene-d12	15.41	264	341348	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.45	112	762679	76.19	ng	0.00
7) Phenol-d6	6.49	99	1075116	78.89	ng	0.00
23) Nitrobenzene-d5	7.42	82	1040065	81.42	ng	0.00
41) 2,4,6-Tribromophenol	10.68	330	246715	75.82	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	1650206	78.51	ng	0.00
78) Terphenyl-d14	12.96	244	1415907	75.60	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.50	88	183228	37.12	ng	99
3) Pyridine	3.22	79	536554	39.18	ng	98
4) n-Nitrosodimethylamine	3.17	42	322757	41.22	ng	95
6) Aniline	6.52	93	747083	38.08	ng	99
8) 2-Chlorophenol	6.64	128	410519	36.59	ng	98
9) Benzaldehyde	6.40	77	316618	37.03	ng	87
10) Phenol	6.50	94	666037	42.23	ng	95
11) bis(2-Chloroethyl)ether	6.59	93	388696	34.50	ng	99
12) 1,3-Dichlorobenzene	6.80	146	491387	39.34	ng	100
13) 1,4-Dichlorobenzene	6.88	146	524822	41.19	ng	99
14) 1,2-Dichlorobenzene	7.02	146	435217	37.37	ng	99
15) Benzyl Alcohol	7.00	79	414476	36.62	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.14	45	892925	37.10	ng	99
17) 2-Methylphenol	7.10	107	349990	38.00	ng	98
18) Hexachloroethane	7.37	117	185164	38.93	ng	96
19) n-Nitroso-di-n-propylamine	7.28	70	352327	38.39	ng	98
20) 3+4-Methylphenols	7.26	107	457786	40.29	ng	99
22) Acetophenone	7.26	105	579424	38.80	ng	98
24) Nitrobenzene	7.44	77	463861	36.24	ng	99
25) Isophorone	7.68	82	881716	38.81	ng	100
26) 2-Nitrophenol	7.76	139	233538m	44.95	ng	
27) 2,4-Dimethylphenol	7.79	122	379573	38.89	ng	98
28) bis(2-Chloroethoxy)methane	7.89	93	552728	41.07	ng	100
29) 2,4-Dichlorophenol	8.00	162	348118	40.13	ng	95
30) 1,2,4-Trichlorobenzene	8.08	180	357520	37.79	ng	100
31) Naphthalene	8.16	128	1129871	36.59	ng	100
32) Benzoic acid	7.90	122	236130	34.59	ng	99
33) 4-Chloroaniline	8.21	127	505871	38.82	ng	97
34) Hexachlorobutadiene	8.28	225	262850	43.04	ng	99
35) Caprolactam	8.59	113	108311	41.70	ng	# 67
36) 4-Chloro-3-methylphenol	8.69	107	390928	41.80	ng	# 84
37) 2-Methylnaphthalene	8.85	142	814900	42.04	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.01	216	338892	34.41	ng	98
40) Hexachlorocyclopentadiene	9.00	237	217893	35.97	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.13	196	251643m	36.73	ng	
43) 2,4,5-Trichlorophenol	9.16	196	262900m	38.69	ng	
45) 1,1'-Biphenyl	9.31	154	823241	34.24	ng	95
46) 2-Chloronaphthalene	9.33	162	693426	35.31	ng	# 90
47) 2-Nitroaniline	9.44	65	292203	38.24	ng	97
48) Acenaphthylene	9.76	152	1224605	39.22	ng	99
49) Dimethylphthalate	9.62	163	838345	37.94	ng	99
50) 2,6-Dinitrotoluene	9.68	165	180839	38.56	ng	97
51) Acenaphthene	9.93	154	764209	40.12	ng	99
52) 3-Nitroaniline	9.85	138	203509	37.61	ng	# 92
53) 2,4-Dinitrophenol	9.94	184	72600	29.07	ng	98
54) Dibenzofuran	10.10	168	982251	38.55	ng	98
55) 4-Nitrophenol	10.00	139	148132	37.15	ng	85
56) 2,4-Dinitrotoluene	10.08	165	257014	41.79	ng	93
57) Fluorene	10.44	166	726316	36.73	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.21	232	204233	39.86	ng	99
59) Diethylphthalate	10.32	149	815881	38.30	ng	99
60) 4-Chlorophenyl-phenylether	10.43	204	360680	42.06	ng	96
61) 4-Nitroaniline	10.45	138	183350	37.53	ng	97
62) Azobenzene	10.59	77	965441	41.84	ng	97
64) 4,6-Dinitro-2-methylphenol	10.48	198	106735	32.43	ng	97
65) n-Nitrosodiphenylamine	10.54	169	627944	35.05	ng	100
66) 4-Bromophenyl-phenylether	10.92	248	231318	41.00	ng	# 92
67) Hexachlorobenzene	10.98	284	235213	36.06	ng	92
68) Atrazine	11.07	200	165192	35.51	ng	99
69) Pentachlorophenol	11.17	266	147394	37.34	ng	99
70) Phenanthrene	11.39	178	957864	34.79	ng	100
71) Anthracene	11.45	178	1135844	41.98	ng	99
72) Carbazole	11.60	167	864443	36.81	ng	99
73) Di-n-butylphthalate	11.94	149	1217991	41.83	ng	98
74) Fluoranthene	12.58	202	1043442	39.19	ng	96
76) Benzidine	12.69	184	443984	34.42	ng	98
77) Pyrene	12.81	202	1012763	36.86	ng	100
79) Butylbenzylphthalate	13.43	149	412980	36.56	ng	87
80) Benzo(a)anthracene	13.99	228	757744	35.57	ng	99
81) 3,3'-Dichlorobenzidine	13.95	252	261589	35.97	ng	# 96
82) Chrysene	14.02	228	675794	32.68	ng	99
83) Bis(2-ethylhexyl)phthalate	13.99	149	520144	38.03	ng	# 94
84) Di-n-octyl phthalate	14.60	149	849841	37.06	ng	99
85) Indeno(1,2,3-cd)pyrene	16.80	276	901358	48.14	ng	# 89
87) Benzo(b)fluoranthene	15.00	252	767882	37.66	ng	# 97
88) Benzo(k)fluoranthene	15.04	252	568095m	33.85	ng	
89) Benzo(a)pyrene	15.36	252	655363	39.15	ng	98
90) Dibenzo(a,h)anthracene	16.81	278	742320	49.54	ng	# 96
91) Benzo(g,h,i)perylene	17.22	276	742576	50.36	ng	# 93

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

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