

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052615\
 Data File : BF079482.D
 Acq On : 26 May 2015 17:49
 Operator : TP/IZ
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

apatel
 5/27/2015 2:00:35 PM

Quant Time: May 26 18:58:23 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 26 18:36:44 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.02	152	88349	20.00	ng	0.00
21) Naphthalene-d8	8.59	136	368987	20.00	ng	0.00
38) Acenaphthene-d10	10.76	164	181858	20.00	ng	0.00
63) Phenanthrene-d10	12.60	188	341535	20.00	ng	0.01
75) Chrysene-d12	15.89	240	330678	20.00	ng	0.01
86) Perylene-d12	17.63	264	355234	20.00	ng	0.05

System Monitoring Compounds

5) 2-Fluorophenol	5.30	112	487946m	94.55	ng	0.00
7) Phenol-d6	6.62	99	626432	91.57	ng	0.00
23) Nitrobenzene-d5	7.72	82	631919	97.58	ng	0.01
41) 2,4,6-Tribromophenol	11.75	330	194378	97.95	ng	0.00
44) 2-Fluorobiphenyl	9.94	172	1108864	84.24	ng	0.00
78) Terphenyl-d14	14.59	244	1262283	89.97	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.85	88	115899m	50.68	ng	
3) Pyridine	2.49	79	266004m	41.18	ng	
4) n-Nitrosodimethylamine	2.43	42	132547m	47.99	ng	
6) Aniline	6.60	93	431577m	45.09	ng	
8) 2-Chlorophenol	6.75	128	287380	48.27	ng	97
9) Benzaldehyde	6.47	77	169690	37.89	ng	99
10) Phenol	6.63	94	367386	46.32	ng	94
11) bis(2-Chloroethyl)ether	6.72	93	268964	46.14	ng	96
12) 1,3-Dichlorobenzene	6.94	146	308365	46.48	ng	98
13) 1,4-Dichlorobenzene	7.04	146	319053	46.87	ng	99
14) 1,2-Dichlorobenzene	7.22	146	293171	46.36	ng	98
15) Benzyl Alcohol	7.22	79	227354	45.02	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.39	45	380861	43.22	ng	99
17) 2-Methylphenol	7.37	107	222483	46.84	ng	98
18) Hexachloroethane	7.63	117	110430	47.20	ng	98
19) n-Nitroso-di-n-propylamine	7.55	70	223762	48.93	ng	# 89
20) 3+4-Methylphenols	7.58	107	310650	49.16	ng	99
22) Acetophenone	7.54	105	387814	46.13	ng	# 89
24) Nitrobenzene	7.75	77	299972	46.71	ng	# 82
25) Isophorone	8.04	82	567266	47.36	ng	98
26) 2-Nitrophenol	8.14	139	152195	56.33	ng	95
27) 2,4-Dimethylphenol	8.22	122	260536	48.99	ng	98
28) bis(2-Chloroethoxy)methane	8.33	93	353518	48.09	ng	100
29) 2,4-Dichlorophenol	8.44	162	236957	49.83	ng	98
30) 1,2,4-Trichlorobenzene	8.54	180	244324	47.06	ng	# 93
31) Naphthalene	8.63	128	838994	47.19	ng	99
32) Benzoic acid	8.36	122	154109	49.63	ng	96
33) 4-Chloroaniline	8.71	127	357507	47.40	ng	99
34) Hexachlorobutadiene	8.78	225	137681	45.90	ng	99
35) Caprolactam	9.15	113	78393	50.93	ng	97
36) 4-Chloro-3-methylphenol	9.34	107	243868	49.35	ng	# 76
37) 2-Methylnaphthalene	9.48	142	593292	48.29	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.69	216	237317	45.75	ng	# 100
40) Hexachlorocyclopentadiene	9.67	237	62228	26.65	ng	95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.84	196	166334	46.71	nq	100
43) 2,4,5-Trichlorophenol	9.90	196	170691	47.70	nq	96
45) 1,1'-Biphenyl	10.07	154	681715	46.58	nq	99
46) 2-Chloronaphthalene	10.08	162	513592	44.71	nq	98
47) 2-Nitroaniline	10.23	65	167169	50.63	nq	# 51
48) Acenaphthylene	10.59	152	878151	46.77	nq	99
49) Dimethylphthalate	10.46	163	616534	45.63	nq	99
50) 2,6-Dinitrotoluene	10.54	165	139952	52.47	nq	# 87
51) Acenaphthene	10.80	154	489686	44.20	nq	100
52) 3-Nitroaniline	10.74	138	173906	52.00	nq	# 74
53) 2,4-Dinitrophenol	10.88	184	44820	59.40	nq	93
54) Dibenzofuran	11.02	168	720384	44.82	nq	99
55) 4-Nitrophenol	10.98	139	91409m	36.86	nq	
56) 2,4-Dinitrotoluene	11.03	165	178895	50.67	nq	92
57) Fluorene	11.44	166	599055	45.06	nq	100
58) 2,3,4,6-Tetrachlorophenol	11.19	232	124186	42.84	nq	# 100
59) Diethylphthalate	11.34	149	626016	46.91	nq	97
60) 4-Chlorophenyl-phenylether	11.45	204	260482	44.45	nq	94
61) 4-Nitroaniline	11.50	138	168926	51.31	nq	88
62) Azobenzene	11.64	77	582945	44.71	nq	99
64) 4,6-Dinitro-2-methylphenol	11.53	198	82459	58.53	nq	82
65) n-Nitrosodiphenylamine	11.61	169	552623	48.23	nq	99
66) 4-Bromophenyl-phenylether	12.06	248	165962	45.90	nq	# 92
67) Hexachlorobenzene	12.11	284	189632	46.01	nq	94
68) Atrazine	12.28	200	149194	44.77	nq	97
69) Pentachlorophenol	12.38	266	80823	41.26	nq	97
70) Phenanthrene	12.63	178	862762	44.89	nq	99
71) Anthracene	12.70	178	904136	47.54	nq	99
72) Carbazole	12.90	167	852652	47.12	nq	99
73) Di-n-butylphthalate	13.36	149	999444	46.35	nq	100
74) Fluoranthene	14.10	202	940947	46.72	nq	96
76) Benzidine	14.30	184	357549	41.55	nq	98
77) Pyrene	14.38	202	956042	49.20	nq	99
79) Butylbenzylphthalate	15.21	149	442668	52.42	nq	92
80) Benzo(a)anthracene	15.87	228	873051	47.61	nq	100
81) 3,3'-Dichlorobenzidine	15.86	252	333954	50.98	nq	# 99
82) Chrysene	15.92	228	857041	48.31	nq	99
83) Bis(2-ethylhexyl)phthalate	15.94	149	629557	49.63	nq	98
84) Di-n-octyl phthalate	16.74	149	1088766	48.39	nq	100
85) Indeno(1,2,3-cd)pyrene	18.97	276	1105825	49.13	nq	99
87) Benzo(b)fluoranthene	17.19	252	874366	44.73	nq	99
88) Benzo(k)fluoranthene	17.22	252	917155m	47.98	nq	
89) Benzo(a)pyrene	17.58	252	871290	47.98	nq	100
90) Dibenzo(a,h)anthracene	18.99	278	893419	46.58	nq	98
91) Benzo(g,h,i)perylene	19.36	276	908032	46.96	nq	99

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

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