

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060615\
 Data File : BF079790.D
 Acq On : 7 Jun 2015 1:25
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Jun 08 19:49:36 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 08 15:29:07 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	68231	20.00	ng	-0.02
21) Naphthalene-d8	8.76	136	269424	20.00	ng	-0.02
38) Acenaphthene-d10	10.55	164	118614	20.00	ng	-0.01
63) Phenanthrene-d10	12.06	188	223864	20.00	ng	-0.01
75) Chrysene-d12	14.93	240	285846	20.00	ng	-0.01
86) Perylene-d12	16.81	264	219024	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	6.09	112	286126	70.28	ng	-0.01
7) Phenol-d6	7.07	99	444436	84.53	ng	-0.01
23) Nitrobenzene-d5	8.03	82	363706	82.88	ng	-0.01
41) 2,4,6-Tribromophenol	11.34	330	81066	71.46	ng	-0.02
44) 2-Fluorobiphenyl	9.84	172	617214	70.50	ng	-0.02
78) Terphenyl-d14	13.69	244	919122	73.64	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	77045	42.77	ng	98
3) Pyridine	4.25	79	169104	33.56	ng	94
4) n-Nitrosodimethylamine	4.17	42	97475	44.58	ng	97
6) Aniline	7.13	93	315771	41.38	ng	100
8) 2-Chlorophenol	7.26	128	189100	42.91	ng	98
9) Benzaldehyde	7.03	77	129433	38.01	ng	# 76
10) Phenol	7.08	94	244458	42.71	ng	98
11) bis(2-Chloroethyl)ether	7.19	93	192052	43.27	ng	98
12) 1,3-Dichlorobenzene	7.42	146	207532	37.62	ng	96
13) 1,4-Dichlorobenzene	7.50	146	230273	39.34	ng	98
14) 1,2-Dichlorobenzene	7.66	146	198296	36.26	ng	98
15) Benzyl Alcohol	7.60	79	164233	37.79	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.74	45	274350	40.87	ng	99
17) 2-Methylphenol	7.70	107	160660	39.09	ng	98
18) Hexachloroethane	8.00	117	71660	42.13	ng	93
19) n-Nitroso-di-n-propylamine	7.86	70	149955	41.10	ng	99
20) 3+4-Methylphenols	7.85	107	215337	41.11	ng	98
22) Acetophenone	7.87	105	280554	43.65	ng	99
24) Nitrobenzene	8.04	77	185461	42.37	ng	97
25) Isophorone	8.28	82	394324	44.02	ng	99
26) 2-Nitrophenol	8.38	139	70714	38.36	ng	93
27) 2,4-Dimethylphenol	8.39	122	174473	42.06	ng	97
28) bis(2-Chloroethoxy)methane	8.49	93	218484	39.47	ng	# 96
29) 2,4-Dichlorophenol	8.62	162	151904	40.81	ng	86
30) 1,2,4-Trichlorobenzene	8.71	180	195910	43.35	ng	97
31) Naphthalene	8.79	128	573695	40.59	ng	100
32) Benzoic acid	8.44	122	57440	37.61	ng	94
33) 4-Chloroaniline	8.82	127	226342m	31.52	ng	
34) Hexachlorobutadiene	8.91	225	105518	38.38	ng	98
35) Caprolactam	9.16	113	45457	40.10	ng	# 66
36) 4-Chloro-3-methylphenol	9.29	107	137118	32.97	ng	92
37) 2-Methylnaphthalene	9.48	142	324440	36.11	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.66	216	154014	41.69	ng	98
40) Hexachlorocyclopentadiene	9.64	237	73472	46.26	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.76	196	99608	40.54	ng	100
43) 2,4,5-Trichlorophenol	9.79	196	102432	39.63	ng	96
45) 1,1'-Biphenyl	9.95	154	371109	35.96	ng	99
46) 2-Chloronaphthalene	9.98	162	296935	35.25	ng	99
47) 2-Nitroaniline	10.06	65	65405	37.01	ng	93
48) Acenaphthylene	10.41	152	508480	36.50	ng	98
49) Dimethylphthalate	10.23	163	337986	36.99	ng	100
50) 2,6-Dinitrotoluene	10.30	165	62345	42.36	ng	95
51) Acenaphthene	10.58	154	295268	37.43	ng	99
52) 3-Nitroaniline	10.47	138	71832	37.23	ng	98
53) 2,4-Dinitrophenol	10.57	184	12831	36.84	ng	# 1
54) Dibenzofuran	10.75	168	428964	38.39	ng	99
55) 4-Nitrophenol	10.60	139	57523	39.91	ng	87
56) 2,4-Dinitrotoluene	10.71	165	77260	43.53	ng	97
57) Fluorene	11.10	166	359888	37.22	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.87	232	75715	36.19	ng	98
59) Diethylphthalate	10.94	149	343884	38.49	ng	99
60) 4-Chlorophenyl-phenylether	11.08	204	160389	36.60	ng	# 75
61) 4-Nitroaniline	11.08	138	73552	39.63	ng	98
62) Azobenzene	11.23	77	318305	35.25	ng	98
64) 4,6-Dinitro-2-methylphenol	11.12	198	23177	42.17	ng	98
65) n-Nitrosodiphenylamine	11.19	169	291161	38.94	ng	99
66) 4-Bromophenyl-phenylether	11.58	248	100216	41.34	ng	96
67) Hexachlorobenzene	11.66	284	104439	38.31	ng	# 63
68) Atrazine	11.71	200	87736	39.04	ng	95
69) Pentachlorophenol	11.85	266	44798	39.44	ng	97
70) Phenanthrene	12.08	178	539643	41.17	ng	99
71) Anthracene	12.12	178	518245	40.10	ng	98
72) Carbazole	12.27	167	498207	40.56	ng	98
73) Di-n-butylphthalate	12.59	149	538799	38.69	ng	# 80
74) Fluoranthene	13.31	202	575893	40.13	ng	97
76) Benzidine	13.42	184	328684	36.43	ng	99
77) Pyrene	13.57	202	682260	38.93	ng	98
79) Butylbenzylphthalate	14.22	149	272359	38.99	ng	# 76
80) Benzo(a)anthracene	14.91	228	683507	39.06	ng	99
81) 3,3'-Dichlorobenzidine	14.85	252	230349	38.75	ng	# 97
82) Chrysene	14.95	228	589924	36.47	ng	98
83) Bis(2-ethylhexyl)phthalate	14.86	149	388109	35.16	ng	# 97
84) Di-n-octyl phthalate	15.62	149	665780	34.25	ng	99
85) Indeno(1,2,3-cd)pyrene	18.79	276	572474	30.57	ng	100
87) Benzo(h)fluoranthene	16.24	252	536512m	39.43	ng	
88) Benzo(k)fluoranthene	16.29	252	534057m	40.92	ng	
89) Benzo(a)pyrene	16.73	252	490912	39.08	ng	# 98
90) Dibenzo(a,h)anthracene	18.81	278	463305	38.22	ng	98
91) Benzo(g,h,i)perylene	19.38	276	469773	37.48	ng	98

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

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