

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050515\
 Data File : BF078943.D
 Acq On : 5 May 2015 16:30
 Operator : TP/IZ
 Sample : G2032-08MSD
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B4(5-7)MSD

Quant Time: May 06 01:54:01 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 06 00:45:00 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.36	152	81310	20.00	ng	0.00
21) Naphthalene-d8	8.94	136	341699	20.00	ng	0.00
38) Acenaphthene-d10	11.11	164	162974	20.00	ng	0.00
63) Phenanthrene-d10	12.96	188	312683	20.00	ng	0.00
75) Chrysene-d12	16.30	240	321359	20.00	ng	0.02
86) Perylene-d12	18.18	264	326521	20.00	ng	0.09

System Monitoring Compounds

5) 2-Fluorophenol	5.66	112	604337	127.94	ng	0.01
7) Phenol-d6	6.90	99	802236	128.49	ng	0.00
23) Nitrobenzene-d5	8.04	82	438824	73.81	ng	0.00
41) 2,4,6-Tribromophenol	12.09	330	220566	125.10	ng	0.00
44) 2-Fluorobiphenyl	10.28	172	933796	79.83	ng	0.00
78) Terphenyl-d14	14.96	244	1087871	81.05	ng	0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.34	88	66071	32.17	ng	# 28
3) Pyridine	3.11	79	145973	24.29	ng	99
4) n-Nitrosodimethylamine	3.02	42	106274m	41.27	ng	
6) Aniline	6.95	93	153128	17.37	ng	99
8) 2-Chlorophenol	7.08	128	225279	42.05	ng	100
9) Benzaldehyde	6.81	77	15820	3.76	ng	100
10) Phenol	6.92	94	311163	42.96	ng	95
11) bis(2-Chloroethyl)ether	7.04	93	214800	40.45	ng	98
12) 1,3-Dichlorobenzene	7.28	146	231794	38.49	ng	97
13) 1,4-Dichlorobenzene	7.38	146	236361	38.14	ng	98
14) 1,2-Dichlorobenzene	7.56	146	229677	39.81	ng	98
15) Benzyl Alcohol	7.53	79	177593	38.32	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.71	45	329840	40.60	ng	100
17) 2-Methylphenol	7.68	107	182539	42.34	ng	94
18) Hexachloroethane	7.98	117	80969	37.93	ng	# 77
19) n-Nitroso-di-n-propylamine	7.87	70	167399	39.69	ng	# 82
20) 3+4-Methylphenols	7.86	107	239061	41.61	ng	94
22) Acetophenone	7.86	105	320847	41.56	ng	# 98
24) Nitrobenzene	8.07	77	228433	38.76	ng	97
25) Isophorone	8.36	82	422689	38.35	ng	99
26) 2-Nitrophenol	8.47	139	104053	42.66	ng	96
27) 2,4-Dimethylphenol	8.52	122	195917	40.14	ng	94
28) bis(2-Chloroethoxy)methane	8.65	93	269818	39.71	ng	99
29) 2,4-Dichlorophenol	8.75	162	180659	41.62	ng	96
30) 1,2,4-Trichlorobenzene	8.87	180	194057	40.70	ng	97
31) Naphthalene	8.96	128	658617	40.45	ng	99
32) Benzoic acid	8.58	122	30145	14.87	ng	92
33) 4-Chloroaniline	9.03	127	125501	18.22	ng	98
34) Hexachlorobutadiene	9.12	225	101509	36.97	ng	97
35) Caprolactam	9.45	113	54704	38.98	ng	95
36) 4-Chloro-3-methylphenol	9.63	107	202054	44.32	ng	91
37) 2-Methylnaphthalene	9.82	142	453168	40.17	ng	98
39) 1,2,4,5-Tetrachlorobenzene	10.02	216	179345	39.07	ng	# 100
40) Hexachlorocyclopentadiene	10.01	237	157018	68.03	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

42) 2,4,6-Trichlorophenol	10.17	196	125816	39.87	nq	98
43) 2,4,5-Trichlorophenol	10.22	196	131821	41.32	nq	# 87
45) 1,1'-Biphenyl	10.40	154	532838	41.13	nq	100
46) 2-Chloronaphthalene	10.42	162	408869	40.46	nq	99
47) 2-Nitroaniline	10.55	65	121834	41.61	nq	98
48) Acenaphthylene	10.94	152	663848	40.01	nq	99
49) Dimethylphthalate	10.79	163	545483	45.61	nq	99
50) 2,6-Dinitrotoluene	10.86	165	95044	40.26	nq	98
51) Acenaphthene	11.15	154	395665	39.99	nq	98
52) 3-Nitroaniline	11.06	138	86956	29.49	nq	# 98
53) 2,4-Dinitrophenol	11.19	184	51593	61.44	nq	# 82
54) Dibenzofuran	11.37	168	598169	41.85	nq	97
55) 4-Nitrophenol	11.27	139	205274	87.96	nq	96
56) 2,4-Dinitrotoluene	11.36	165	134389	43.06	nq	# 67
57) Fluorene	11.79	166	483250	41.19	nq	99
58) 2,3,4,6-Tetrachlorophenol	11.52	232	106760	40.95	nq	# 100
59) Diethylphthalate	11.67	149	471406	39.78	nq	100
60) 4-Chlorophenyl-phenylether	11.81	204	205773	39.54	nq	95
61) 4-Nitroaniline	11.82	138	109114	36.99	nq	93
62) Azobenzene	12.00	77	496744	42.54	nq	96
64) 4,6-Dinitro-2-methylphenol	11.85	198	46077	38.48	nq	93
65) n-Nitrosodiphenylamine	11.94	169	400313	38.44	nq	98
66) 4-Bromophenyl-phenylether	12.41	248	132608	40.69	nq	93
67) Hexachlorobenzene	12.47	284	145106	38.95	nq	93
68) Atrazine	12.62	200	124659	41.17	nq	99
69) Pentachlorophenol	12.72	266	150729	70.09	nq	99
70) Phenanthrene	12.98	178	675113	38.59	nq	100
71) Anthracene	13.05	178	681315	39.70	nq	100
72) Carbazole	13.26	167	690722	42.23	nq	98
73) Di-n-butylphthalate	13.70	149	779031	39.71	nq	99
74) Fluoranthene	14.47	202	763333	41.88	nq	96
76) Benzidine	14.64	184	172306	19.89	nq	99
77) Pyrene	14.75	202	789518	42.63	nq	99
79) Butylbenzylphthalate	15.59	149	345551	42.69	nq	# 77
80) Benzo(a)anthracene	16.29	228	728242	41.38	nq	99
81) 3,3'-Dichlorobenzidine	16.26	252	211079	33.67	nq	100
82) Chrysene	16.33	228	674242	39.84	nq	99
83) Bis(2-ethylhexyl)phthalate	16.34	149	509588	41.56	nq	99
84) Di-n-octyl phthalate	17.20	149	869862	40.28	nq	# 100
85) Indeno(1,2,3-cd)pyrene	19.71	276	888303	41.19	nq	# 100
87) Benzo(b)fluoranthene	17.69	252	778832	43.58	nq	99
88) Benzo(k)fluoranthene	17.73	252	671784	38.83	nq	99
89) Benzo(a)pyrene	18.11	252	708492	43.05	nq	# 96
90) Dibenzo(a,h)anthracene	19.75	278	704333	40.27	nq	98
91) Benzo(g,h,i)perylene	20.17	276	736277	42.00	nq	99

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

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