

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011515\
 Data File : BF076903.D
 Acq On : 16 Jan 2015 4:08
 Operator : IZ / TP
 Sample : G1076-11MS
 Misc : S
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-3MS

Quant Time: Jan 16 11:39:17 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 15 11:52:04 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	38211	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	167038	20.00	ng	0.00
38) Acenaphthene-d10	15.07	164	90506	20.00	ng	-0.01
63) Phenanthrene-d10	17.82	188	148095	20.00	ng	0.00
75) Chrysene-d12	21.31	240	150633	20.00	ng	0.00
86) Perylene-d12	22.95	264	145196	20.00	ng	-0.05

System Monitoring Compounds

5) 2-Fluorophenol	5.10	112	27088	11.66	ng	0.01
7) Phenol-d6	7.42	99	32465	11.07	ng	0.00
23) Nitrobenzene-d5	9.32	82	23977	7.95	ng	0.00
41) 2,4,6-Tribromophenol	16.72	330	8190	11.15	ng	0.00
44) 2-Fluorobiphenyl	13.58	172	51272	8.62	ng	-0.01
78) Terphenyl-d14	20.04	244	51542	7.88	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.33	88	3270	3.29	ng	# 87
3) Pyridine	1.89	79	6050	2.35	ng	90
4) n-Nitrosodimethylamine	1.81	42	3661	2.87	ng	# 93
6) Aniline	7.32	93	14143	3.49	ng	95
8) 2-Chlorophenol	7.56	128	10181	3.80	ng	98
10) Phenol	7.44	94	10631	3.41	ng	91
11) bis(2-Chloroethyl)ether	7.56	93	9687	3.98	ng	85
12) 1,3-Dichlorobenzene	7.86	146	11523	3.78	ng	92
13) 1,4-Dichlorobenzene	8.06	146	11903	3.68	ng	91
14) 1,2-Dichlorobenzene	8.37	146	11562	3.88	ng	95
15) Benzyl Alcohol	8.45	79	8539	3.60	ng	# 91
16) 2,2'-oxybis(1-Chloropropan	8.80	45	13478	4.12	ng	80
17) 2-Methylphenol	8.79	107	7632	3.49	ng	88
18) Hexachloroethane	9.12	117	3888	3.46	ng	91
19) n-Nitroso-di-n-propylamine	9.08	70	6749	3.29	ng	# 98
20) 3+4-Methylphenols	9.18	107	10257	3.54	ng	89
22) Acetophenone	9.00	105	15192	3.71	ng	97
24) Nitrobenzene	9.35	77	10934	3.56	ng	# 91
25) Isophorone	9.94	82	19663	3.58	ng	# 92
26) 2-Nitrophenol	10.10	139	4141	4.30	ng	# 78
27) 2,4-Dimethylphenol	10.37	122	8670	3.65	ng	98
28) bis(2-Chloroethoxy)methane	10.57	93	12262	3.93	ng	# 98
29) 2,4-Dichlorophenol	10.71	162	7288	3.29	ng	95
30) 1,2,4-Trichlorobenzene	10.84	180	9725	3.96	ng	97
31) Naphthalene	10.97	128	33685	3.92	ng	97
33) 4-Chloroaniline	11.21	127	12618	3.63	ng	93
34) Hexachlorobutadiene	11.34	225	5325	3.65	ng	96
36) 4-Chloro-3-methylphenol	12.52	107	9265	3.66	ng	95
37) 2-Methylnaphthalene	12.63	142	21111	3.59	ng	98
39) 1,2,4,5-Tetrachlorobenzene	13.03	216	9070	4.05	ng	93
40) Hexachlorocyclopentadiene	13.02	237	2159	6.27	ng	92
42) 2,4,6-Trichlorophenol	13.39	196	4466	2.74	ng	92
43) 2,4,5-Trichlorophenol	13.48	196	5385m	3.24	ng	
45) 1,1'-Biphenyl	13.79	154	26589	3.96	ng	95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 2-Chloronaphthalene	13.76	162	21594	3.91	nq	93
47) 2-Nitroaniline	14.11	65	5221	3.12	nq	91
48) Acenaphthylene	14.71	152	33525	3.60	nq	# 91
49) Dimethylphthalate	14.64	163	25215	3.93	nq	100
50) 2,6-Dinitrotoluene	14.75	165	4311	3.36	nq	88
51) Acenaphthene	15.15	154	19154	3.62	nq	91
52) 3-Nitroaniline	15.10	138	5241	4.83	nq	# 73
54) Dibenzofuran	15.57	168	27958	3.63	nq	96
55) 4-Nitrophenol	15.72	139	4383m	4.36	nq	
56) 2,4-Dinitrotoluene	15.68	165	4959	4.58	nq	# 91
57) Fluorene	16.27	166	24101	3.70	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.90	232	3841	3.17	nq	# 94
59) Diethylphthalate	16.25	149	25035	3.86	nq	98
60) 4-Chlorophenyl-phenylether	16.36	204	10146	3.80	nq	95
61) 4-Nitroaniline	16.39	138	4352	4.04	nq	# 79
62) Azobenzene	16.63	77	22219	3.29	nq	95
64) 4,6-Dinitro-2-methylphenol	16.48	198	228	3.53	nq	# 1
65) n-Nitrosodiphenylamine	16.60	169	19718	3.73	nq	92
66) 4-Bromophenyl-phenylether	17.19	248	6437	4.29	nq	93
67) Hexachlorobenzene	17.21	284	5866	3.71	nq	# 83
68) Atrazine	17.55	200	6355	3.97	nq	96
69) Pentachlorophenol	17.58	266	1397	8.67	nq	# 84
70) Phenanthrene	17.84	178	34537	3.74	nq	96
71) Anthracene	17.92	178	33557	3.73	nq	96
72) Carbazole	18.21	167	32578	3.73	nq	96
73) Di-n-butylphthalate	18.79	149	37033	3.66	nq	100
74) Fluoranthene	19.49	202	34620	3.66	nq	95
77) Pvrene	19.77	202	36691	3.55	nq	98
79) Butylbenzylphthalate	20.70	149	14331	3.07	nq	94
80) Benzo(a)anthracene	21.29	228	33796	3.58	nq	97
81) 3,3'-Dichlorobenzidine	21.31	252	11791	3.93	nq	# 92
82) Chrysene	21.34	228	32658	3.79	nq	99
83) Bis(2-ethylhexyl)phthalate	21.43	149	22282	3.44	nq	# 97
84) Di-n-octyl phthalate	22.20	149	35062	3.12	nq	# 95
85) Indeno(1,2,3-cd)pvrene	24.05	276	34447m	3.77	nq	
87) Benzo(b)fluoranthene	22.54	252	31248	3.20	nq	# 91
88) Benzo(k)fluoranthene	22.57	252	32342m	3.79	nq	
89) Benzo(a)pyrene	22.88	252	32314	3.75	nq	# 98
90) Dibenzo(a,h)anthracene	24.08	278	29183m	3.69	nq	
91) Benzo(g,h,i)perylene	24.35	276	30259m	3.85	nq	

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

