

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011915\
 Data File : BF076946.D
 Acq On : 20 Jan 2015 2:52
 Operator : IZ / TP
 Sample : G1076-11MSD
 Misc : W
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-3MSD

Quant Time: Jan 20 04:14:42 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 20 03:20:25 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.98	152	40070	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	178197	20.00	ng	0.00
38) Acenaphthene-d10	15.01	164	97925	20.00	ng	0.00
63) Phenanthrene-d10	17.77	188	163325	20.00	ng	0.00
75) Chrysene-d12	21.27	240	171112	20.00	ng	0.00
86) Perylene-d12	22.93	264	154200	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.04	112	29500	12.10	ng	0.00
7) Phenol-d6	7.37	99	34937	11.36	ng	0.00
23) Nitrobenzene-d5	9.27	82	24737	7.69	ng	0.00
41) 2,4,6-Tribromophenol	16.68	330	10029	12.62	ng	0.00
44) 2-Fluorobiphenyl	13.53	172	55319	8.60	ng	0.00
78) Terphenyl-d14	20.00	244	55230	7.43	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.29	88	3533	3.39	ng	91
4) n-Nitrosodimethylamine	1.77	42	3578	2.67	ng	# 93
6) Aniline	7.27	93	12293	2.89	ng	98
8) 2-Chlorophenol	7.51	128	9743	3.47	ng	94
10) Phenol	7.41	94	10386	3.18	ng	93
11) bis(2-Chloroethyl)ether	7.50	93	10173	3.98	ng	90
12) 1,3-Dichlorobenzene	7.82	146	11226	3.51	ng	# 82
13) 1,4-Dichlorobenzene	8.01	146	12117	3.57	ng	96
14) 1,2-Dichlorobenzene	8.32	146	11961	3.83	ng	90
15) Benzyl Alcohol	8.41	79	7842	3.15	ng	87
16) 2,2'-oxybis(1-Chloropropan	8.76	45	14206	4.14	ng	92
17) 2-Methylphenol	8.76	107	7584	3.30	ng	# 90
18) Hexachloroethane	9.08	117	5118	4.34	ng	# 84
19) n-Nitroso-di-n-propylamine	9.03	70	7420	3.45	ng	92
20) 3+4-Methylphenols	9.14	107	9848	3.24	ng	91
22) Acetophenone	8.96	105	15420	3.53	ng	94
24) Nitrobenzene	9.32	77	11575	3.53	ng	99
25) Isophorone	9.91	82	19787	3.38	ng	96
26) 2-Nitrophenol	10.06	139	4048	4.09	ng	# 92
27) 2,4-Dimethylphenol	10.33	122	8143	3.21	ng	88
28) bis(2-Chloroethoxy)methane	10.53	93	12019	3.61	ng	93
29) 2,4-Dichlorophenol	10.68	162	7111	3.01	ng	# 72
30) 1,2,4-Trichlorobenzene	10.79	180	9656	3.68	ng	99
31) Naphthalene	10.93	128	32725	3.57	ng	99
33) 4-Chloroaniline	11.18	127	11412	3.08	ng	# 96
34) Hexachlorobutadiene	11.29	225	5407	3.47	ng	95
35) Caprolactam	11.96	113	1432	2.06	ng	# 74
36) 4-Chloro-3-methylphenol	12.47	107	9106	3.37	ng	98
37) 2-Methylnaphthalene	12.59	142	23033	3.68	ng	95
39) 1,2,4,5-Tetrachlorobenzene	12.99	216	9945	4.11	ng	91
40) Hexachlorocyclopentadiene	12.97	237	4261	7.66	ng	91
42) 2,4,6-Trichlorophenol	13.34	196	5555	3.15	ng	99
43) 2,4,5-Trichlorophenol	13.43	196	5333	2.96	ng	# 82
45) 1,1'-Biphenyl	13.74	154	28362	3.91	ng	95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 2-Chloronaphthalene	13.72	162	21742	3.64	ng	94
47) 2-Nitroaniline	14.06	65	5838	3.22	ng	95
48) Acenaphthylene	14.67	152	34199	3.39	ng	96
49) Dimethylphthalate	14.60	163	25949	3.74	ng	99
50) 2,6-Dinitrotoluene	14.70	165	5180	3.73	ng	91
51) Acenaphthene	15.09	154	18558	3.25	ng	96
52) 3-Nitroaniline	15.05	138	5298	4.63	ng	# 73
53) 2,4-Dinitrophenol	15.35	184	1256	13.40	ng	# 71
54) Dibenzofuran	15.51	168	32967	3.96	ng	97
55) 4-Nitrophenol	15.66	139	5248	4.82	ng	# 79
56) 2,4-Dinitrotoluene	15.64	165	6121	4.92	ng	# 92
57) Fluorene	16.22	166	27077	3.84	ng	97
58) 2,3,4,6-Tetrachlorophenol	15.85	232	3978	3.04	ng	# 87
59) Diethylphthalate	16.21	149	26073	3.71	ng	98
60) 4-Chlorophenyl-phenylether	16.31	204	10971	3.80	ng	97
61) 4-Nitroaniline	16.36	138	5018	4.21	ng	99
62) Azobenzene	16.60	77	25472	3.48	ng	95
64) 4,6-Dinitro-2-methylphenol	16.44	198	2165	5.23	ng	77
65) n-Nitrosodiphenylamine	16.55	169	20349	3.49	ng	94
66) 4-Bromophenyl-phenylether	17.16	248	6507	3.93	ng	# 83
67) Hexachlorobenzene	17.17	284	7324	4.20	ng	# 87
68) Atrazine	17.51	200	6557	3.71	ng	92
69) Pentachlorophenol	17.53	266	3030	10.37	ng	97
70) Phenanthrene	17.81	178	37929	3.72	ng	98
71) Anthracene	17.89	178	36141	3.64	ng	97
72) Carbazole	18.17	167	34527	3.58	ng	95
73) Di-n-butylphthalate	18.76	149	39013	3.50	ng	98
74) Fluoranthene	19.45	202	38342	3.67	ng	98
77) Pvrene	19.74	202	40298	3.44	ng	99
79) Butylbenzylphthalate	20.67	149	17057	3.21	ng	96
80) Benzo(a)anthracene	21.26	228	38963	3.63	ng	99
81) 3,3'-Dichlorobenzidine	21.27	252	11442	3.35	ng	# 87
82) Chrysene	21.31	228	35600	3.64	ng	98
83) Bis(2-ethylhexyl)phthalate	21.40	149	23977	3.26	ng	# 94
84) Di-n-octyl phthalate	22.16	149	40152	3.15	ng	100
85) Indeno(1,2,3-cd)pyrene	24.05	276	36191	3.49	ng	# 82
87) Benzo(b)fluoranthene	22.51	252	36499	3.51	ng	98
88) Benzo(k)fluoranthene	22.54	252	32239m	3.55	ng	
89) Benzo(a)pyrene	22.87	252	34364	3.75	ng	99
90) Dibenzo(a,h)anthracene	24.07	278	30923	3.68	ng	# 92
91) Benzo(g,h,i)perylene	24.35	276	29055	3.48	ng	98

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

