

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050217\
 Data File : BF094807.D
 Acq On : 2 May 2017 22:32
 Operator : SJ/MA
 Sample : I2928-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-1-COMP

Quant Time: May 03 02:51:55 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 02 17:18:57 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	87603	20.00	ng	0.00
21) Naphthalene-d8	9.75	136	361040	20.00	ng	0.00
38) Acenaphthene-d10	12.58	164	155000	20.00	ng	0.00
63) Phenanthrene-d10	14.98	188	280968	20.00	ng	0.00
75) Chrysene-d12	18.64	240	207004	20.00	ng	-0.01
86) Perylene-d12	20.31	264	163606	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.68	112	558634	106.32	ng	0.00
7) Phenol-d6	7.27	99	693315	109.97	ng	0.00
23) Nitrobenzene-d5	8.62	82	380109	66.39	ng	-0.01
41) 2,4,6-Tribromophenol	13.87	330	144783	114.06	ng	-0.01
44) 2-Fluorobiphenyl	11.52	172	794935	73.82	ng	0.00
78) Terphenyl-d14	17.31	244	684169	72.80	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.05	88	118	0.046	ng	# 73
6) Aniline	7.29	93	118	0.015	ng	# 1
8) 2-Chlorophenol	7.41	128	137	0.023	ng	# 1
9) Benzaldehyde	7.27	77	661	0.172	ng	# 1
10) Phenol	7.28	94	14864	2.237	ng	# 84
11) bis(2-Chloroethyl)ether	7.39	93	608	0.111	ng	# 1
15) Benzyl Alcohol	7.97	79	5938	1.464	ng	# 42
16) 2,2'-oxybis(1-Chloropropan	8.31	45	121	0.016	ng	# 70
24) Nitrobenzene	8.62	77	863	0.144	ng	# 39
25) Isophorone	9.17	82	108	0.011	ng	# 67
45) 1,1'-Biphenyl	11.52	154	1338	0.103	ng	# 1
49) Dimethylphthalate	12.21	163	271745	21.356	ng	# 98
50) 2,6-Dinitrotoluene	12.21	165	2825	1.088	ng	# 11
51) Acenaphthene	11.52	154	1338	0.103	ng	# 1
59) Diethylphthalate	13.37	149	1052	0.086	ng	# 60
62) Azobenzene	13.87	77	620	0.063	ng	# 9
64) 4,6-Dinitro-2-methylphenol	13.87	198	126	0.067	ng	# 1
65) n-Nitrosodiphenylamine	13.87	169	4919	0.482	ng	# 43
70) Phenanthrene	15.01	178	662	0.041	ng	# 59
71) Anthracene	15.01	178	662	0.040	ng	# 59
73) Di-n-butylphthalate	15.95	149	4598	0.246	ng	# 88
74) Fluoranthene	16.78	202	1358	0.082	ng	# 62
77) Pyrene	17.07	202	1656	0.107	ng	# 68
79) Butylbenzylphthalate	17.96	149	119	0.017	ng	# 32
80) Benzo(a)anthracene	18.63	228	1049	0.087	ng	# 50
82) Chrysene	18.67	228	983	0.084	ng	# 49
83) Bis(2-ethylhexyl)phthalate	18.71	149	4493	0.438	ng	# 97
84) Di-n-octyl phthalate	19.53	149	864	0.051	ng	# 1
87) Benzo(b)fluoranthene	19.92	252	825	0.082	ng	# 59
88) Benzo(k)fluoranthene	19.92	252	825	0.085	ng	# 56
89) Benzo(a)pyrene	20.25	252	516	0.055	ng	# 29
91) Benzo(g,h,i)perylene	22.03	276	116	0.015	ng	# 47

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

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