

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100118\
 Data File : BF109490.D
 Acq On : 1 Oct 2018 23:01
 Operator : JU/SJ
 Sample : J5155-04
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 QNWP8-MW9-GW

Quant Time: Oct 02 04:59:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	96348	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	235989	20.00	ng	0.01
38) Acenaphthene-d10	9.74	164	150196	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	246247	20.00	ng	0.00
75) Chrysene-d12	13.85	240	233391	20.00	ng	0.00
86) Perylene-d12	15.24	264	167743	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	42874	7.33	ng	0.02
7) Phenol-d6	6.37	99	76130	10.20	ng	0.01
23) Nitrobenzene-d5	7.27	82	118044	29.53	ng	0.00
41) 2,4,6-Tribromophenol	10.53	330	28081	18.97	ng	0.00
44) 2-Fluorobiphenyl	9.06	172	196689	19.75	ng	-0.01
78) Terphenyl-d14	12.79	244	105011	11.04	ng	-0.01

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.36	88	7498	2.783	ng	93
3) Pyridine	3.10	79	88586	12.776	ng	92
6) Aniline	6.38	93	174071	18.228	ng	# 23
10) Phenol	6.38	94	197414	23.354	ng	# 63
15) Benzyl Alcohol	6.87	79	19924	3.487	ng	89
17) 2-Methylphenol	6.99	107	160327	30.461	ng	# 92
20) 3+4-Methylphenols	7.14	107	425391	66.941	ng	# 50
22) Acetophenone	7.13	105	1026604	188.160	ng	# 93
27) 2,4-Dimethylphenol	7.68	122	371525	118.445	ng	98
31) Naphthalene	8.06	128	13219542	1198.758	ng	# 93
37) 2-Methylnaphthalene	8.72	142	1507909	212.111	ng	96
45) 1,1'-Biophenyl	9.16	154	259516	21.207	ng	97
48) Acenaphthylene	9.60	152	51857	3.516	ng	# 87
49) Dimethylphthalate	9.47	163	274854	25.160	ng	99
51) Acenaphthene	9.77	154	336586	37.748	ng	99
54) Dibenzofuran	9.95	168	431360	32.529	ng	98
57) Fluorene	10.28	166	192097	20.303	ng	# 97
70) Phenanthrene	11.24	178	349241	28.405	ng	99
71) Anthracene	11.29	178	44659	3.581	ng	99
72) Carbazole	11.45	167	211051	17.010	ng	99
74) Fluoranthene	12.42	202	45877	3.492	ng	97

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

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