

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122617\
 Data File : BF101719.D
 Acq On : 26 Dec 2017 16:54
 Operator : SJ/JU
 Sample : I6846-03DL 20X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-34972326-01DL

Quant Time: Dec 26 17:51:16 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 26 10:48:01 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	135625	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	507644	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	207604	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	334090	20.00	ng	0.00
75) Chrysene-d12	13.97	240	263950	20.00	ng	0.00
86) Perylene-d12	15.43	264	262682	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	69603	7.64	ng	0.05
7) Phenol-d6	6.43	99	87337	7.71	ng	0.00
23) Nitrobenzene-d5	7.36	82	63220	6.93	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	14818	7.41	ng	0.01
44) 2-Fluorobiphenyl	9.14	172	132719	9.92	ng	0.00
78) Terphenyl-d14	12.89	244	95551	8.73	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.44	94	758556	58.734	ng	99
17) 2-Methylphenol	7.05	107	181940	20.651	ng	100
20) 3+4-Methylphenols	7.20	107	558374	59.809	ng	89
27) 2,4-Dimethylphenol	7.74	122	103422	14.040	ng	96
31) Naphthalene	8.09	128	1186256	46.417	ng	100
37) 2-Methylnaphthalene	8.79	142	55734	3.347	ng	98
72) Carbazole	11.56	167	44779	2.520	ng	98

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

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