

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081717\
 Data File : BF097813.D
 Acq On : 18 Aug 2017 4:35
 Operator : SJ/JU
 Sample : I4752-01
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RMAB-MW-12-GW

Quant Time: Aug 18 05:12:21 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 15 13:25:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	147094	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	512456	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	187560	20.00	ng	0.00
63) Phenanthrene-d10	11.29	188	314338	20.00	ng	-0.01
75) Chrysene-d12	13.93	240	325271	20.00	ng	-0.01
86) Perylene-d12	15.36	264	227849	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	264652	27.37	ng	0.00
7) Phenol-d6	6.41	99	580783	44.20	ng	0.00
23) Nitrobenzene-d5	7.35	82	840110	89.06	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	192752	118.44	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	1113486	87.22	ng	0.00
78) Terphenyl-d14	12.89	244	893589	57.29	ng	0.00
Target Compounds						
17) 2-Methylphenol	7.03	107	75270	8.375	ng	Qvalue # 65
27) 2,4-Dimethylphenol	7.72	122	162551	19.870	ng	90
31) Naphthalene	8.09	128	2444721	91.892	ng	98
32) Benzoic acid	7.83	122	56612	14.635	ng	# 83
37) 2-Methylnaphthalene	8.77	142	485814	29.785	ng	97

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

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