

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM090717\
 Data File : BM011460.D
 Acq On : 07 Sep 2017 18:27
 Operator : SJ/JU
 Sample : I5110-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP-2

Quant Time: Sep 08 08:17:30 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM083117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 31 18:39:59 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.98	152	403283	20.00	ng	-0.01
21) Naphthalene-d8	10.79	136	1656641	20.00	ng	-0.02
38) Acenaphthene-d10	14.61	164	849245	20.00	ng	-0.02
63) Phenanthrene-d10	17.34	188	1843210	20.00	ng	-0.02
75) Chrysene-d12	21.51	240	2066266	20.00	ng	-0.02
86) Perylene-d12	23.89	264	1862776	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	3513619	146.66	ng	0.00
7) Phenol-d6	7.13	99	4265064	128.38	ng	0.00
23) Nitrobenzene-d5	9.15	82	2636443	104.87	ng	-0.01
41) 2,4,6-Tribromophenol	16.09	330	1356050	138.01	ng	-0.02
44) 2-Fluorobiphenyl	13.23	172	5782105	98.14	ng	-0.02
78) Terphenyl-d14	19.96	244	8246856	99.03	ng	-0.02

Target Compounds Qvalue

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

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