

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082218\
 Data File : BM016527.D
 Acq On : 22 Aug 2018 20:21
 Operator : SJ/JU
 Sample : J4545-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RB-BASELINE-WATER

Quant Time: Aug 23 03:44:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 21 16:31:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	104672	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	312335	20.00	ng	0.00
38) Acenaphthene-d10	14.66	164	236622	20.00	ng	-0.01
63) Phenanthrene-d10	17.40	188	791073	20.00	ng	0.00
75) Chrysene-d12	21.54	240	1066506	20.00	ng	0.00
86) Perylene-d12	23.83	264	1234199	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	709832	132.89	ng	0.00
7) Phenol-d6	7.19	99	732674	103.66	ng	0.00
23) Nitrobenzene-d5	9.20	82	826605	120.84	ng	0.00
41) 2,4,6-Tribromophenol	16.15	330	1290947	195.17	ng	0.00
44) 2-Fluorobiphenyl	13.27	172	2878516	116.93	ng	0.00
78) Terphenyl-d14	19.99	244	7633442	151.49	ng	0.00

Target Compounds Qvalue

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

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