

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072318\
 Data File : BM016044.D
 Acq On : 24 Jul 2018 02:21
 Operator : SJ/JU
 Sample : J3762-07
 Misc : LOD 8 PPM
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOD-MDL-WATER-01-QT3-2018

Quant Time: Jul 24 08:37:42 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 23 16:08:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.23	152	41478	20.00	ng	0.00
21) Naphthalene-d8	11.05	136	157224	20.00	ng	0.00
38) Acenaphthene-d10	14.86	164	82659	20.00	ng	0.00
63) Phenanthrene-d10	17.59	188	176016	20.00	ng	0.00
75) Chrysene-d12	21.70	240	189160	20.00	ng	0.00
86) Perylene-d12	24.07	264	201271	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.74	112	325210	130.24	ng	0.00
7) Phenol-d6	7.39	99	395134	121.17	ng	0.00
23) Nitrobenzene-d5	9.42	82	299721	88.67	ng	0.00
41) 2,4,6-Tribromophenol	16.34	330	135738	129.47	ng	0.00
44) 2-Fluorobiphenyl	13.48	172	595169	87.61	ng	0.00
78) Terphenyl-d14	20.15	244	1096858	121.38	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	13.02	237	6545	11.956	ng	98
53) 2,4-Dinitrophenol	15.02	184	2347	9.587	ng	99
69) Pentachlorophenol	17.24	266	5702	4.195	ng	92

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

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