

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072518\
 Data File : BN002156.D
 Acq On : 26 Jul 2018 03:09
 Operator : JU/SJ
 Sample : J3762-07
 Misc : LOD 8 PPM
 ALS Vial : 17 Sample Multiplier: 1

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

Quant Time: Jul 26 07:28:16 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN072518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 25 20:08:02 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	204320	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	793887	20.00	ng	0.00
38) Acenaphthene-d10	14.32	164	437480	20.00	ng	0.00
63) Phenanthrene-d10	17.07	188	896189	20.00	ng	0.00
75) Chrysene-d12	21.27	240	858553	20.00	ng	0.00
86) Perylene-d12	23.49	264	841385	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	1599568	126.04	ng	0.00
7) Phenol-d6	6.88	99	2111694	118.57	ng	0.00
23) Nitrobenzene-d5	8.84	82	1376061	88.89	ng	0.00
41) 2,4,6-Tribromophenol	15.82	330	711674	125.28	ng	0.00
44) 2-Fluorobiphenyl	12.94	172	2944324	87.66	ng	0.00
78) Terphenyl-d14	19.71	244	4247649	103.19	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	12.48	237	67270	8.097	ng	91
53) 2,4-Dinitrophenol	14.45	184	20347	4.760	ng	86
69) Pentachlorophenol	16.73	266	34382	4.603	ng	94

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

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