

Data Path : Z:\HPCHEM1\BNA N\DATA\BN041818\
 Data File : BN000756.D
 Acq On : 19 Apr 2018 00:51
 Operator : JU/SJ
 Sample : J2133-07
 Misc : 8PPM LOD
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 LOD-MDL-WATER-01-QT2-2018

Quant Time: Apr 19 07:16:23 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN041818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 18 16:58:01 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	105106	20.00	ng	0.00
21) Naphthalene-d8	10.39	136	439056	20.00	ng	0.00
38) Acenaphthene-d10	14.26	164	226682	20.00	ng	0.00
63) Phenanthrene-d10	17.00	188	464434	20.00	ng	0.00
75) Chrysene-d12	21.21	240	464550	20.00	ng	-0.01
86) Perylene-d12	23.40	264	469187	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	971815	137.35	ng	0.00
7) Phenol-d6	6.81	99	1269671	131.68	ng	0.00
23) Nitrobenzene-d5	8.78	82	727043	86.32	ng	0.00
41) 2,4,6-Tribromophenol	15.75	330	332741	145.81	ng	0.00
44) 2-Fluorobiphenyl	12.88	172	1440124	87.70	ng	0.00
78) Terphenyl-d14	19.66	244	1764325	81.89	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	12.42	237	28727	6.968	ng	91
53) 2,4-Dinitrophenol	14.37	184	8404	9.606	ng	97
69) Pentachlorophenol	16.66	266	19693	5.579	ng	99

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

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