

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101718\
 Data File : BM017175.D
 Acq On : 18 Oct 2018 01:23
 Operator : MJ/SJ
 Sample : J5164-07
 Misc : LOD 8 PPM
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOD-MDL-WATER-01-QT4-2018

Quant Time: Oct 18 13:25:09 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM101718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 17 14:23:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	221715	20.00	ng	0.00
21) Naphthalene-d8	10.45	136	923883	20.00	ng	0.00
39) Acenaphthene-d10	14.31	164	544136	20.00	ng	0.00
64) Phenanthrene-d10	17.06	188	1281083	20.00	ng	0.00
76) Chrysene-d12	21.26	240	1525069	20.00	ng	0.00
87) Perylene-d12	23.47	264	1572117	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	1575447	120.18	ng	0.00
7) Phenol-d6	6.85	99	2061305	115.46	ng	0.00
23) Nitrobenzene-d5	8.82	82	1217163	77.05	ng	-0.01
42) 2,4,6-Tribromophenol	15.81	330	886824	120.51	ng	0.00
45) 2-Fluorobiphenyl	12.93	172	2945036	72.01	ng	0.00
79) Terphenyl-d14	19.70	244	4667775	68.99	ng	0.00
Target Compounds						
41) Hexachlorocyclopentadiene	12.46	237	74009	7.989	ng	100
54) 2,4-Dinitrophenol	14.44	184	23038	12.039	ng	94
69) Atrazine	16.54	200	86315	6.222	ng	96

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

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