

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN102418\
 Data File : BN003287.D
 Acq On : 25 Oct 2018 00:24
 Operator : MJ/SJ
 Sample : J5164-07
 Misc : LOD 8PPM
 ALS Vial : 23 Sample Multiplier: 1

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

Quant Time: Oct 25 07:11:18 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN102418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 24 14:21:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	107179	20.00	ng	0.00
21) Naphthalene-d8	10.37	136	498959	20.00	ng	0.00
39) Acenaphthene-d10	14.24	164	320364	20.00	ng	0.00
64) Phenanthrene-d10	17.00	188	753985	20.00	ng	0.00
76) Chrysene-d12	21.20	240	852516	20.00	ng	0.00
87) Perylene-d12	23.41	264	885711	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	758240	123.22	ng	0.00
7) Phenol-d6	6.80	99	1065131	121.73	ng	0.00
23) Nitrobenzene-d5	8.75	82	623848	71.70	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	516172	119.66	ng	0.00
45) 2-Fluorobiphenyl	12.85	172	1638439	69.37	ng	0.00
79) Terphenyl-d14	19.64	244	2648543	66.44	ng	0.00
Target Compounds						
41) Hexachlorocyclopentadiene	12.39	237	13850	11.606	ng	99
54) 2,4-Dinitrophenol	14.43	184	8203	8.693	ng	# 90
70) Pentachlorophenol	16.68	266	15363	9.420	ng	97

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

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