

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050118\
 Data File : BG034107.D
 Acq On : 2 May 2018 9:17
 Operator : SJ/JU
 Sample : J2133-09
 Misc : 8PPM MDL
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MDL-WATER-03-QT2-2018

Quant Time: May 02 13:57:17 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 01 19:08:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	63939	20.00	ng	0.00
21) Naphthalene-d8	10.89	136	271578	20.00	ng	0.00
38) Acenaphthene-d10	14.72	164	210353	20.00	ng	0.00
63) Phenanthrene-d10	17.47	188	561398	20.00	ng	0.00
75) Chrysene-d12	21.76	240	612587	20.00	ng	0.00
86) Perylene-d12	25.05	264	601721	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	530557	134.06	ng	0.00
7) Phenol-d6	7.22	99	801617	130.11	ng	0.00
23) Nitrobenzene-d5	9.23	82	586291	82.83	ng	0.00
41) 2,4,6-Tribromophenol	16.20	330	383653	131.71	ng	0.00
44) 2-Fluorobiphenyl	13.34	172	1232366	85.37	ng	0.00
78) Terphenyl-d14	20.09	244	2330033	83.32	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	12.89	237	37109	7.694	ng	89
53) 2,4-Dinitrophenol	14.81	184	9865	7.198	ng	# 56
69) Pentachlorophenol	17.11	266	35200	7.138	ng	90

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

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