

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032217\
 Data File : BM009333.D
 Acq On : 23 Mar 2017 00:10
 Operator : SJ/MA
 Sample : I2296-05
 Misc : LOD 8 PPB
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOD-WATER-05-QT2-2017

Quant Time: Mar 23 16:33:08 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM032217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 22 16:43:02 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	186931	20.00	ng	0.00
21) Naphthalene-d8	10.66	136	701312	20.00	ng	0.00
38) Acenaphthene-d10	14.50	164	403972	20.00	ng	0.00
63) Phenanthrene-d10	17.25	188	916571	20.00	ng	0.00
75) Chrysene-d12	21.42	240	1189253	20.00	ng	0.00
86) Perylene-d12	23.75	264	1116481	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	1534604	136.84	ng	0.00
7) Phenol-d6	7.02	99	2089258	136.61	ng	0.00
23) Nitrobenzene-d5	9.03	82	1691055	90.41	ng	0.00
41) 2,4,6-Tribromophenol	15.99	330	720895	143.65	ng	0.00
44) 2-Fluorobiphenyl	13.12	172	3030069	90.38	ng	0.00
78) Terphenyl-d14	19.87	244	4607473	80.96	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	12.66	237	59922	6.85	ng	95
53) 2,4-Dinitrophenol	14.62	184	24221	10.47	ng	96
69) Pentachlorophenol	16.89	266	46666	6.56	ng	89

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

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