

Data Path : Z:\HPCHEM1\BNA M\DATA\BM052617\
 Data File : BM010256.D
 Acq On : 26 May 2017 14:35
 Operator : SJ/MA
 Sample : I3239-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PURGE-WATER-B

Quant Time: May 27 00:29:33 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM051917.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 20 04:25:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.94	152	106959	20.00	ng	-0.04
21) Naphthalene-d8	10.75	136	374535	20.00	ng	-0.04
38) Acenaphthene-d10	14.57	164	263432	20.00	ng	-0.04
63) Phenanthrene-d10	17.32	188	688488	20.00	ng	-0.04
75) Chrysene-d12	21.48	240	981567	20.00	ng	-0.03
86) Perylene-d12	23.83	264	1054062	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	388458	65.59	ng	-0.04
7) Phenol-d6	7.13	99	285963	37.52	ng	-0.04
23) Nitrobenzene-d5	9.13	82	755450	108.82	ng	-0.04
41) 2,4,6-Tribromophenol	16.06	330	610187	152.51	ng	-0.04
44) 2-Fluorobiphenyl	13.19	172	2008487	98.02	ng	-0.04
78) Terphenyl-d14	19.92	244	4042146	93.64	ng	-0.03
Target Compounds						
10) Phenol	7.15	94	23808	2.96	ng	95
14) 1,2-Dichlorobenzene	8.29	146	18099	2.24	ng	96
51) Acenaphthene	14.63	154	69419	4.31	ng	97
57) Fluorene	15.62	166	61885	2.81	ng	99
70) Phenanthrene	17.36	178	191220	4.96	ng	98

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

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