

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060618\
 Data File : BM015513.D
 Acq On : 07 Jun 2018 06:31
 Operator : SJ/JU
 Sample : J3306-03
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A&M-WATER

Quant Time: Jun 07 07:57:45 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM052518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 06 14:01:06 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	143105	20.00	ng	0.00
21) Naphthalene-d8	11.19	136	577562	20.00	ng	0.00
38) Acenaphthene-d10	14.97	164	302831	20.00	ng	0.00
63) Phenanthrene-d10	17.69	188	645016	20.00	ng	0.00
75) Chrysene-d12	21.79	240	655282	20.00	ng	-0.01
86) Perylene-d12	24.21	264	646646	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.86	112	572330	68.03	ng	0.00
7) Phenol-d6	7.53	99	706411	55.55	ng	0.02
23) Nitrobenzene-d5	9.54	82	562258	52.95	ng	0.00
41) 2,4,6-Tribromophenol	16.44	330	308698	70.93	ng	0.00
44) 2-Fluorobiphenyl	13.60	172	1160777	51.47	ng	0.00
78) Terphenyl-d14	20.24	244	1669782	58.82	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.58	88	9606	3.062	ng	Qvalue # 100
10) Phenol	7.55	94	92778	7.194	ng	81
20) 3+4-Methylphenols	9.13	107	151231	12.052	ng	89
32) Benzoic acid	10.53	122	243687	36.490	ng	94
49) Dimethylphthalate	14.42	163	121694	4.680	ng	100

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

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