

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101719\
 Data File : BG043268.D
 Acq On : 17 Oct 2019 19:00
 Operator : JU
 Sample : K5407-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DEVELOPMENT-WATER

Quant Time: Oct 18 02:05:50 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 09 01:11:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	60407	20.00	ng	-0.02
21) Naphthalene-d8	10.85	136	235680	20.00	ng	-0.02
39) Acenaphthene-d10	14.67	164	168063	20.00	ng	-0.02
64) Phenanthrene-d10	17.41	188	365467	20.00	ng	0.00
76) Chrysene-d12	21.68	240	394750	20.00	ng	0.00
87) Perylene-d12	24.91	264	469335	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	228903	67.63	ng	-0.02
7) Phenol-d6	7.21	99	221708	45.48	ng	-0.02
23) Nitrobenzene-d5	9.20	82	305368	62.98	ng	-0.02
42) 2,4,6-Tribromophenol	16.16	330	312006	107.96	ng	0.00
45) 2-Fluorobiphenyl	13.29	172	799993	69.01	ng	-0.02
79) Terphenyl-d14	20.02	244	1417284	76.51	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.50	88	16053	11.375	ng	93
10) Phenol	7.24	94	14688	3.284	ng	97
50) Dimethylphthalate	14.11	163	78715	6.251	ng	98

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

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