

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102822\
 Data File : BG055351.D
 Acq On : 28 Oct 2022 21:26
 Operator : CG/JU
 Sample : N5226-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20221093-198-SITE-WATER

Quant Time: Oct 29 04:56:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 29 04:52:38 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.264	152	29897	20.000	ng	0.00
21) Naphthalene-d8	11.090	136	125114	20.000	ng	0.00
39) Acenaphthene-d10	14.874	164	86758	20.000	ng	0.00
64) Phenanthrene-d10	17.606	188	199417	20.000	ng	0.00
76) Chrysene-d12	21.883	240	205895	20.000	ng	0.00
86) Perylene-d12	25.268	264	216184	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.785	112	112525	65.902	ng	0.00
7) Phenol-d6	7.406	99	115338	46.626	ng	0.00
23) Nitrobenzene-d5	9.433	82	204181	83.333	ng	0.00
42) 2,4,6-Tribromophenol	16.355	330	154027	163.314	ng	0.00
45) 2-Fluorobiphenyl	13.505	172	511670	96.530	ng	0.00
79) Terphenyl-d14	20.180	244	938473	94.701	ng	0.00

Target Compounds	Qvalue

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

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