

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
 Data File : BG055690.D
 Acq On : 18 Nov 2022 14:23
 Operator : CG/JU
 Sample : N5598-12
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 FRAC-TANK-SLURRY-0760533

Quant Time: Nov 18 23:56:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 17 02:02:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.255	152	39731	20.000	ng	0.00
21) Naphthalene-d8	11.081	136	164077	20.000	ng	0.00
39) Acenaphthene-d10	14.877	164	113978	20.000	ng	0.00
64) Phenanthrene-d10	17.615	188	248030	20.000	ng	0.00
76) Chrysene-d12	21.916	240	233174	20.000	ng	0.00
86) Perylene-d12	25.335	264	245511	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.787	112	245621	111.743	ng	0.00
7) Phenol-d6	7.397	99	316230	108.075	ng	0.00
23) Nitrobenzene-d5	9.424	82	197837	63.730	ng	0.00
42) 2,4,6-Tribromophenol	16.357	330	169072	105.344	ng	0.00
45) 2-Fluorobiphenyl	13.502	172	513582	67.505	ng	0.00
79) Terphenyl-d14	20.206	244	842015	72.227	ng	0.00

Target Compounds	Qvalue

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

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