

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
 Data File : BG055687.D
 Acq On : 18 Nov 2022 12:20
 Operator : CG/JU
 Sample : N5598-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 FRAC-TANK-0760533

Quant Time: Nov 18 23:54:50 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	8.252	152	33462	20.000	ng	0.00
21) Naphthalene-d8	11.078	136	128007	20.000	ng	0.00
39) Acenaphthene-d10	14.873	164	85783	20.000	ng	0.00
64) Phenanthrene-d10	17.611	188	210091	20.000	ng	0.00
76) Chrysene-d12	21.912	240	215703	20.000	ng	0.00
86) Perylene-d12	25.326	264	228307	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.784	112	257498	139.093	ng	0.00
7) Phenol-d6	7.400	99	321199	130.339	ng	0.01
23) Nitrobenzene-d5	9.421	82	198835	82.100	ng	0.00
42) 2,4,6-Tribromophenol	16.354	330	173026	143.241	ng	0.00
45) 2-Fluorobiphenyl	13.499	172	447538	78.159	ng	0.00
79) Terphenyl-d14	20.202	244	796038	73.813	ng	0.00

Target Compounds	Qvalue

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

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