

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG113022\
 Data File : BG055849.D
 Acq On : 30 Nov 2022 19:53
 Operator : CG/JU
 Sample : N5809-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 HOWE-DRUM-A-1129

Quant Time: Dec 01 02:14:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 01 02:10:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.223	152	44315	20.000	ng	0.00
21) Naphthalene-d8	11.055	136	174394	20.000	ng	0.00
39) Acenaphthene-d10	14.850	164	111088	20.000	ng	0.00
64) Phenanthrene-d10	17.594	188	220762	20.000	ng	# 0.00
76) Chrysene-d12	21.913	240	192607	20.000	ng	# 0.03
86) Perylene-d12	25.403	264	192236	20.000	ng	# 0.14
System Monitoring Compounds						
5) 2-Fluorophenol	5.767	112	274241	111.857	ng	0.02
7) Phenol-d6	7.383	99	345208	105.775	ng	0.00
23) Nitrobenzene-d5	9.398	82	121733	36.894	ng	0.00
42) 2,4,6-Tribromophenol	16.337	330	148370	94.850	ng	0.00
45) 2-Fluorobiphenyl	13.470	172	487253	65.711	ng	0.00
79) Terphenyl-d14	20.185	244	597937	62.093	ng	0.02
Target Compounds						
						Qvalue
75) Fluoranthene	19.639	202	46356	3.198	ng	# 47
78) Pyrene	20.003	202	45835	3.516	ng	# 47

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

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