

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011624\
 Data File : BG060338.D
 Acq On : 16 Jan 2024 16:16
 Operator : MA/JU
 Sample : P1146-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 2ND-FLOOR-SAMPLE-A

Quant Time: Jan 16 16:49:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG010824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 09 04:07:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.931	152	231823	20.000	ng	0.00
21) Naphthalene-d8	10.728	136	943828	20.000	ng	0.00
39) Acenaphthene-d10	14.564	164	701291	20.000	ng	0.00
64) Phenanthrene-d10	17.308	188	1828189	20.000	ng	0.00
76) Chrysene-d12	21.574	240	1710481	20.000	ng	0.00
86) Perylene-d12	24.741	264	1935152	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	1232372	87.437	ng	0.00
7) Phenol-d6	7.091	99	1549292	74.233	ng	0.00
23) Nitrobenzene-d5	9.088	82	1051337	52.599	ng	0.00
42) 2,4,6-Tribromophenol	16.051	330	669657	64.902	ng	0.00
45) 2-Fluorobiphenyl	13.184	172	2758989	54.697	ng	0.00
79) Terphenyl-d14	19.929	244	4470541	58.579	ng	0.00

Target Compounds	Qvalue

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

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