

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011624\
 Data File : BG060337.D
 Acq On : 16 Jan 2024 15:36
 Operator : MA/JU
 Sample : P1146-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 1ST-FLOOR-SAMPLE-B

Quant Time: Jan 16 16:10:56 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	7.926	152	229313	20.000	ng	0.00
21) Naphthalene-d8	10.728	136	946945	20.000	ng	0.00
39) Acenaphthene-d10	14.565	164	710160	20.000	ng	0.00
64) Phenanthrene-d10	17.309	188	1788471	20.000	ng	0.00
76) Chrysene-d12	21.574	240	1632799	20.000	ng	0.00
86) Perylene-d12	24.741	264	1856049	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.505	112	1431637	102.687	ng	0.00
7) Phenol-d6	7.091	99	1747787	84.661	ng	0.00
23) Nitrobenzene-d5	9.089	82	1162478	57.968	ng	0.00
42) 2,4,6-Tribromophenol	16.051	330	794254	76.016	ng	0.00
45) 2-Fluorobiphenyl	13.184	172	3048972	59.691	ng	0.00
79) Terphenyl-d14	19.929	244	4610078	65.236	ng	0.00

Target Compounds	Qvalue

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

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