

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051324\
 Data File : BG061176.D
 Acq On : 13 May 2024 15:24
 Operator : MA/JU
 Sample : P2430-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 1011UMR-SS001-1224-01

Quant Time: May 13 16:22:18 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG043024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 05:11:25 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.929	152	20541	20.000	ng	-0.03
21) Naphthalene-d8	10.720	136	72077	20.000	ng	#-0.04
39) Acenaphthene-d10	14.557	164	56351	20.000	ng	-0.03
64) Phenanthrene-d10	17.306	188	158972	20.000	ng	-0.03
76) Chrysene-d12	21.560	240	190442	20.000	ng	-0.03
86) Perylene-d12	24.686	264	216995	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.526	112	153125	151.607	ng	-0.02
7) Phenol-d6	7.112	99	155284	103.929	ng	-0.02
23) Nitrobenzene-d5	9.081	82	133528	91.836	ng	-0.04
42) 2,4,6-Tribromophenol	16.049	330	159241	141.377	ng	-0.03
45) 2-Fluorobiphenyl	13.176	172	332738	87.314	ng	-0.03
79) Terphenyl-d14	19.927	244	998191	87.173	ng	-0.03

Target Compounds	Qvalue

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

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