

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050924\
 Data File : BG061148.D
 Acq On : 9 May 2024 12:13
 Operator : MA/JU
 Sample : PB160779BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB160779BL

Quant Time: May 09 12:51:38 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	7.935	152	19930	20.000	ng	-0.03
21) Naphthalene-d8	10.731	136	77812	20.000	ng	#-0.03
39) Acenaphthene-d10	14.562	164	63422	20.000	ng	-0.03
64) Phenanthrene-d10	17.312	188	176354	20.000	ng	-0.02
76) Chrysene-d12	21.571	240	188665	20.000	ng	-0.02
86) Perylene-d12	24.697	264	216731	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.526	112	154144	157.295	ng	-0.02
7) Phenol-d6	7.112	99	206875	142.703	ng	-0.02
23) Nitrobenzene-d5	9.092	82	142955	91.073	ng	-0.03
42) 2,4,6-Tribromophenol	16.060	330	171370	135.182	ng	-0.02
45) 2-Fluorobiphenyl	13.187	172	399095	93.051	ng	-0.02
79) Terphenyl-d14	19.932	244	1041912	91.848	ng	-0.02

Target Compounds	Qvalue

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

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