

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051024\
 Data File : BG061161.D
 Acq On : 10 May 2024 12:47
 Operator : MA/JU
 Sample : PB160758TB
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB160758TB

Quant Time: May 10 13:19:44 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	7.928	152	18027	20.000	ng	-0.03
21) Naphthalene-d8	10.725	136	71346	20.000	ng	-0.03
39) Acenaphthene-d10	14.562	164	64879	20.000	ng	-0.03
64) Phenanthrene-d10	17.311	188	173932	20.000	ng	#-0.02
76) Chrysene-d12	21.571	240	185704	20.000	ng	-0.02
86) Perylene-d12	24.697	264	219412	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.519	112	142016	160.217	ng	-0.03
7) Phenol-d6	7.106	99	188864	144.032	ng	-0.03
23) Nitrobenzene-d5	9.086	82	129883	90.245	ng	-0.03
42) 2,4,6-Tribromophenol	16.054	330	170514	131.487	ng	-0.02
45) 2-Fluorobiphenyl	13.181	172	379014	86.384	ng	-0.03
79) Terphenyl-d14	19.932	244	1003072	89.834	ng	-0.02

Target Compounds	Qvalue

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

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