

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090524\
Data File : BG062669.D
Acq On : 5 Sep 2024 16:31
Operator : MA/JU
Sample : PB163138BL
Misc : MDL-WATER-8 ng
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PB163138BL

Quant Time: Sep 05 17:48:51 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG083024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Aug 31 02:36:01 2024
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.917	152	45157	20.000	ng	0.00
21) Naphthalene-d8	10.714	136	170932	20.000	ng	0.00
39) Acenaphthene-d10	14.545	164	134891	20.000	ng	0.00
64) Phenanthrene-d10	17.288	188	399629	20.000	ng	0.00
76) Chrysene-d12	21.530	240	482411	20.000	ng	0.00
86) Perylene-d12	24.597	264	550716	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.502	112	360806	138.111	ng	0.00
7) Phenol-d6	7.083	99	466162	123.518	ng	0.00
23) Nitrobenzene-d5	9.074	82	284428	89.616	ng	0.00
42) 2,4,6-Tribromophenol	16.031	330	269933	142.194	ng	0.00
45) 2-Fluorobiphenyl	13.164	172	726325	87.218	ng	0.00
79) Terphenyl-d14	19.909	244	2085294	85.711	ng	0.00

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

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