

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090524\
Data File : BG062670.D
Acq On : 5 Sep 2024 17:12
Operator : MA/JU
Sample : PB163139BL
Misc : MDL-WATER-8 ng
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PB163139BL

Quant Time: Sep 05 17:49:14 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	7.919	152	44900	20.000	ng	0.00
21) Naphthalene-d8	10.715	136	166560	20.000	ng	0.00
39) Acenaphthene-d10	14.540	164	130204	20.000	ng	0.00
64) Phenanthrene-d10	17.284	188	386128	20.000	ng	0.00
76) Chrysene-d12	21.532	240	474388	20.000	ng	0.00
86) Perylene-d12	24.599	264	535371	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	379458	146.082	ng	0.00
7) Phenol-d6	7.084	99	488137	130.081	ng	0.00
23) Nitrobenzene-d5	9.070	82	288592	93.315	ng	0.00
42) 2,4,6-Tribromophenol	16.033	330	281182	153.452	ng	0.00
45) 2-Fluorobiphenyl	13.165	172	742005	92.308	ng	0.00
79) Terphenyl-d14	19.905	244	2172078	90.788	ng	0.00

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

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