

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090424\
 Data File : BG062654.D
 Acq On : 4 Sep 2024 23:05
 Operator : MA/JU
 Sample : PB163109BL
 Misc : LOD-MDL-WATER-8 ng
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB163109BL

Quant Time: Sep 05 04:54:45 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	7.918	152	51429	20.000	ng	0.00
21) Naphthalene-d8	10.715	136	197880	20.000	ng	0.00
39) Acenaphthene-d10	14.545	164	157879	20.000	ng	0.00
64) Phenanthrene-d10	17.283	188	480092	20.000	ng	0.00
76) Chrysene-d12	21.531	240	552095	20.000	ng	0.00
86) Perylene-d12	24.598	264	621188	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.503	112	377179	126.771	ng	0.00
7) Phenol-d6	7.084	99	513205	119.399	ng	0.00
23) Nitrobenzene-d5	9.070	82	311424	84.759	ng	0.00
42) 2,4,6-Tribromophenol	16.032	330	316333	142.373	ng	0.00
45) 2-Fluorobiphenyl	13.165	172	805878	82.681	ng	0.00
79) Terphenyl-d14	19.904	244	2260128	81.172	ng	0.00

Target Compounds	Qvalue

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

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