

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110124\
 Data File : BG063146.D
 Acq On : 1 Nov 2024 13:37
 Operator : RC/JU
 Sample : PB164382BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB164382BL

Quant Time: Nov 01 14:31:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG103124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 01 05:13:40 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.843	152	97652	20.000	ng	0.00
21) Naphthalene-d8	10.634	136	385106	20.000	ng	# 0.00
39) Acenaphthene-d10	14.477	164	302732	20.000	ng	0.00
64) Phenanthrene-d10	17.227	188	757559	20.000	ng	0.00
76) Chrysene-d12	21.480	240	928440	20.000	ng	0.00
86) Perylene-d12	24.512	264	1037693	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	877724	139.375	ng	0.00
7) Phenol-d6	7.021	99	1166575	141.360	ng	0.00
23) Nitrobenzene-d5	8.995	82	840439	94.917	ng	0.00
42) 2,4,6-Tribromophenol	15.969	330	669309	130.157	ng	0.00
45) 2-Fluorobiphenyl	13.102	172	1886365	92.290	ng	0.00
79) Terphenyl-d14	19.859	244	4216446	93.456	ng	0.00

Target Compounds	Qvalue

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

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