

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092925\
 Data File : BG064457.D
 Acq On : 29 Sep 2025 13:02
 Operator : RC/JU
 Sample : PB169873BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB169873BL

Quant Time: Sep 29 18:11:10 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG092525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 29 18:04:12 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.838	152	15751	20.000	ng	0.00
21) Naphthalene-d8	10.623	136	63617	20.000	ng	# 0.00
39) Acenaphthene-d10	14.460	164	53384	20.000	ng	0.00
64) Phenanthrene-d10	17.203	188	149845	20.000	ng	# 0.00
76) Chrysene-d12	21.428	240	170237	20.000	ng	0.00
86) Perylene-d12	24.424	264	187310	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.429	112	118221	143.429	ng	0.00
7) Phenol-d6	7.010	99	158449	140.318	ng	0.00
23) Nitrobenzene-d5	8.990	82	106882	91.052	ng	0.00
42) 2,4,6-Tribromophenol	15.952	330	134606	144.434	ng	0.00
45) 2-Fluorobiphenyl	13.085	172	284697	78.750	ng	0.00
79) Terphenyl-d14	19.824	244	803405	93.032	ng	0.00

Target Compounds	Qvalue

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

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