

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092225\
 Data File : BG064419.D
 Acq On : 22 Sep 2025 17:38
 Operator : RC/JU
 Sample : PB169783BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB169783BL

Quant Time: Sep 22 18:22:32 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 28 17:41:58 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.842	152	19046	20.000	ng	-0.02
21) Naphthalene-d8	10.627	136	76167	20.000	ng	#-0.02
39) Acenaphthene-d10	14.463	164	56503	20.000	ng	-0.02
64) Phenanthrene-d10	17.207	188	147930	20.000	ng	#-0.02
76) Chrysene-d12	21.437	240	169803	20.000	ng	#-0.02
86) Perylene-d12	24.434	264	186504	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.433	112	153456	144.552	ng	-0.01
7) Phenol-d6	7.013	99	195036	133.728	ng	-0.02
23) Nitrobenzene-d5	8.993	82	122948	90.024	ng	-0.02
42) 2,4,6-Tribromophenol	15.956	330	126556	169.100	ng	-0.02
45) 2-Fluorobiphenyl	13.088	172	329310	88.959	ng	-0.02
79) Terphenyl-d14	19.828	244	781675	96.043	ng	-0.02

Target Compounds Qvalue

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

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