

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092225\
 Data File : BG064421.D
 Acq On : 22 Sep 2025 18:59
 Operator : RC/JU
 Sample : PB169754TB
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB169754TB

Quant Time: Sep 23 01:51:37 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	7.838	152	17631	20.000	ng	-0.02
21) Naphthalene-d8	10.628	136	68244	20.000	ng	#-0.02
39) Acenaphthene-d10	14.465	164	51178	20.000	ng	-0.02
64) Phenanthrene-d10	17.203	188	135776	20.000	ng	#-0.02
76) Chrysene-d12	21.433	240	156064	20.000	ng	#-0.02
86) Perylene-d12	24.430	264	171396	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	168183	171.140	ng	0.00
7) Phenol-d6	7.015	99	192749	142.767	ng	-0.02
23) Nitrobenzene-d5	8.989	82	132892	108.602	ng	-0.03
42) 2,4,6-Tribromophenol	15.957	330	140266	206.920	ng	-0.02
45) 2-Fluorobiphenyl	13.090	172	357045	106.486	ng	-0.02
79) Terphenyl-d14	19.829	244	854443	114.227	ng	-0.02

Target Compounds Qvalue

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

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