

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120925\
 Data File : BG064909.D
 Acq On : 9 Dec 2025 17:41
 Operator : RC/JU
 Sample : PB170870TB
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB170870TB

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/10/2025
 Supervised By :Sohil Jodhani 12/12/2025

Quant Time: Dec 10 00:14:21 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 12 17:10:56 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.152	152	20536	20.000	ng	-0.01
21) Naphthalene-d8	10.973	136	77670	20.000	ng	#-0.01
39) Acenaphthene-d10	14.762	164	73436	20.000	ng	-0.01
64) Phenanthrene-d10	17.494	188	177293m	20.000	ng	0.00
76) Chrysene-d12	21.742	240	221194	20.000	ng	#-0.01
86) Perylene-d12	24.992	264	243364	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.691	112	150916	128.327	ng	0.00
7) Phenol-d6	7.295	99	195249	124.826	ng	-0.01
23) Nitrobenzene-d5	9.316	82	124297	79.124	ng	-0.01
42) 2,4,6-Tribromophenol	16.237	330	222570	155.401	ng	-0.01
45) 2-Fluorobiphenyl	13.393	172	459926	87.390	ng	-0.01
79) Terphenyl-d14	20.080	244	1160309	103.187	ng	-0.01

Target Compounds Qvalue

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

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