

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102925\
 Data File : BG064619.D
 Acq On : 29 Oct 2025 12:03
 Operator : RC/JU
 Sample : PB170285BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB170285BL

Quant Time: Oct 29 12:43:09 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102425.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 24 16:40:26 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.216	152	31966	20.000	ng	0.00
21) Naphthalene-d8	11.042	136	126218	20.000	ng	-0.01
39) Acenaphthene-d10	14.838	164	100016	20.000	ng	0.00
64) Phenanthrene-d10	17.576	188	253171	20.000	ng	0.00
76) Chrysene-d12	21.847	240	318569	20.000	ng	-0.02
86) Perylene-d12	25.191	264	365282	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.749	112	281567	147.847	ng	0.00
7) Phenol-d6	7.353	99	368638	141.063	ng	0.00
23) Nitrobenzene-d5	9.380	82	225643	107.375	ng	-0.01
42) 2,4,6-Tribromophenol	16.319	330	234059	162.143	ng	0.00
45) 2-Fluorobiphenyl	13.469	172	627891	95.157	ng	0.00
79) Terphenyl-d14	20.167	244	1533731	90.856	ng	0.00

Target Compounds	Qvalue

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

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