

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112125\
 Data File : BG064839.D
 Acq On : 21 Nov 2025 10:49
 Operator : RC/JU
 Sample : PB170631BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB170631BL

Quant Time: Nov 21 11:24:58 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	8.159	152	44450	20.000	ng	0.00
21) Naphthalene-d8	10.979	136	174541	20.000	ng	0.00
39) Acenaphthene-d10	14.769	164	146167	20.000	ng	0.00
64) Phenanthrene-d10	17.501	188	361264	20.000	ng	0.00
76) Chrysene-d12	21.749	240	444014	20.000	ng	0.00
86) Perylene-d12	25.004	264	434648	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.703	112	329551	129.464	ng	0.00
7) Phenol-d6	7.307	99	425828	125.775	ng	0.00
23) Nitrobenzene-d5	9.323	82	262931	74.481	ng	0.00
42) 2,4,6-Tribromophenol	16.250	330	335625	117.734	ng	0.00
45) 2-Fluorobiphenyl	13.406	172	785262	74.963	ng	0.00
79) Terphenyl-d14	20.086	244	1938900	85.898	ng	0.00

Target Compounds	Qvalue

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

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