

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110525\
 Data File : BG064691.D
 Acq On : 5 Nov 2025 18:45
 Operator : RC/JU
 Sample : PB170402BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB170402BL

Quant Time: Nov 06 00:08:41 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102425.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 03 18:16:26 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.205	152	19318	20.000	ng	-0.02
21) Naphthalene-d8	11.031	136	77931	20.000	ng	#-0.02
39) Acenaphthene-d10	14.826	164	74381	20.000	ng	-0.02
64) Phenanthrene-d10	17.564	188	240875	20.000	ng	#-0.02
76) Chrysene-d12	21.836	240	296369	20.000	ng	#-0.02
86) Perylene-d12	25.167	264	292507	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.737	112	168297	146.229	ng	-0.01
7) Phenol-d6	7.341	99	221270	140.107	ng	-0.01
23) Nitrobenzene-d5	9.368	82	134601	103.739	ng	-0.02
42) 2,4,6-Tribromophenol	16.307	330	215582	200.814	ng	-0.01
45) 2-Fluorobiphenyl	13.457	172	436331	88.916	ng	-0.02
79) Terphenyl-d14	20.155	244	1661997	105.830	ng	-0.02

Target Compounds	Qvalue

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

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