

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101625\
 Data File : BG064517.D
 Acq On : 16 Oct 2025 15:34
 Operator : RC/JU
 Sample : PB170133BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB170133BL

Quant Time: Oct 16 16:25:57 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 02 16:40:42 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.823	152	14441	20.000	ng	# 0.00
21) Naphthalene-d8	10.608	136	57385	20.000	ng	# 0.00
39) Acenaphthene-d10	14.444	164	43972	20.000	ng	0.00
64) Phenanthrene-d10	17.188	188	106337	20.000	ng	# 0.00
76) Chrysene-d12	21.413	240	129197	20.000	ng	0.00
86) Perylene-d12	24.397	264	136933	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.414	112	123308	158.025	ng	0.00
7) Phenol-d6	7.000	99	152191	145.137	ng	0.01
23) Nitrobenzene-d5	8.968	82	105929	99.256	ng	0.00
42) 2,4,6-Tribromophenol	15.937	330	123174	154.537	ng	0.00
45) 2-Fluorobiphenyl	13.070	172	284117	95.240	ng	0.00
79) Terphenyl-d14	19.809	244	661947	96.963	ng	0.00

Target Compounds Qvalue

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

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