

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120925\
 Data File : BG064906.D
 Acq On : 9 Dec 2025 15:39
 Operator : RC/JU
 Sample : Q3800-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ETGI-288

Quant Time: Dec 09 16:22:07 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	8.150	152	25289	20.000	ng	-0.02
21) Naphthalene-d8	10.964	136	100169	20.000	ng	#-0.02
39) Acenaphthene-d10	14.765	164	91375	20.000	ng	-0.01
64) Phenanthrene-d10	17.492	188	223127	20.000	ng	#-0.01
76) Chrysene-d12	21.740	240	277372	20.000	ng	#-0.02
86) Perylene-d12	24.995	264	311452	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.682	112	27675	19.110	ng	-0.02
7) Phenol-d6	7.292	99	51793	26.889	ng	-0.02
23) Nitrobenzene-d5	9.313	82	69465	34.288	ng	-0.02
42) 2,4,6-Tribromophenol	16.240	330	116879	65.585	ng	-0.01
45) 2-Fluorobiphenyl	13.396	172	297032	45.358	ng	-0.01
79) Terphenyl-d14	20.077	244	804854	57.079	ng	-0.02

Target Compounds	Qvalue

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

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