

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100325\
 Data File : BG064491.D
 Acq On : 3 Oct 2025 17:38
 Operator : RC/JU
 Sample : Q3263-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 251818

Quant Time: Oct 03 18:29:12 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	7.826	152	14769	20.000	ng	0.00
21) Naphthalene-d8	10.611	136	62942	20.000	ng	# 0.00
39) Acenaphthene-d10	14.447	164	49465	20.000	ng	0.00
64) Phenanthrene-d10	17.185	188	123783	20.000	ng	# 0.00
76) Chrysene-d12	21.410	240	133977	20.000	ng	0.00
86) Perylene-d12	24.383	264	141843	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.411	112	42580	53.356	ng	0.00
7) Phenol-d6	6.991	99	31420	29.298	ng	0.00
23) Nitrobenzene-d5	8.977	82	88968	76.003	ng	0.01
42) 2,4,6-Tribromophenol	15.940	330	129213	144.111	ng	0.00
45) 2-Fluorobiphenyl	13.072	172	253941	75.672	ng	0.00
79) Terphenyl-d14	19.806	244	711442	100.495	ng	0.00

Target Compounds	Qvalue

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

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