

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092225\
 Data File : BG064432.D
 Acq On : 23 Sep 2025 2:30
 Operator : RC/JU
 Sample : Q3151-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 1402

Quant Time: Sep 23 03:53:38 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 28 17:41:58 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.838	152	15997	20.000	ng	-0.02
21) Naphthalene-d8	10.623	136	62473	20.000	ng	#-0.03
39) Acenaphthene-d10	14.459	164	48564	20.000	ng	-0.02
64) Phenanthrene-d10	17.203	188	124373	20.000	ng	#-0.02
76) Chrysene-d12	21.428	240	149935	20.000	ng	-0.03
86) Perylene-d12	24.430	264	164471	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.429	112	60446	67.791	ng	-0.02
7) Phenol-d6	7.009	99	45847	37.427	ng	-0.02
23) Nitrobenzene-d5	8.989	82	134388	119.970	ng	-0.03
42) 2,4,6-Tribromophenol	15.952	330	152381	236.892	ng	-0.02
45) 2-Fluorobiphenyl	13.084	172	373102	117.265	ng	-0.03
79) Terphenyl-d14	19.824	244	870579	121.141	ng	-0.02

Target Compounds	Qvalue

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

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