

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110625\
 Data File : BG064708.D
 Acq On : 6 Nov 2025 18:30
 Operator : RC/JU
 Sample : Q3532-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-1

Quant Time: Nov 07 00:32:27 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	8.198	152	31168	20.000	ng	-0.02
21) Naphthalene-d8	11.024	136	119813	20.000	ng	#-0.02
39) Acenaphthene-d10	14.820	164	100517	20.000	ng	-0.02
64) Phenanthrene-d10	17.558	188	244672	20.000	ng	#-0.02
76) Chrysene-d12	21.823	240	305345	20.000	ng	#-0.04
86) Perylene-d12	25.149	264	330107	20.000	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.736	112	216961	116.840	ng	-0.01
7) Phenol-d6	7.340	99	263828	103.541	ng	-0.01
23) Nitrobenzene-d5	9.361	82	203880	102.206	ng	-0.02
42) 2,4,6-Tribromophenol	16.300	330	212931	146.772	ng	-0.02
45) 2-Fluorobiphenyl	13.451	172	570949	86.096	ng	-0.02
79) Terphenyl-d14	20.149	244	1422992	87.947	ng	-0.02

Target Compounds	Qvalue

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

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