

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100325\
 Data File : BG064493.D
 Acq On : 3 Oct 2025 18:59
 Operator : RC/JU
 Sample : Q3273-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW1

Quant Time: Oct 04 00:00:10 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	7.820	152	20058	20.000	ng	0.00
21) Naphthalene-d8	10.611	136	82373	20.000	ng	# 0.00
39) Acenaphthene-d10	14.459	164	60802	20.000	ng	0.01
64) Phenanthrene-d10	17.197	188	129947	20.000	ng	# 0.01
76) Chrysene-d12	21.416	240	136555	20.000	ng	# 0.00
86) Perylene-d12	24.389	264	149983	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.417	112	73920	68.203	ng	0.01
7) Phenol-d6	6.997	99	59110	40.584	ng	0.01
23) Nitrobenzene-d5	8.977	82	138717	90.549	ng	0.01
42) 2,4,6-Tribromophenol	15.946	330	147739	134.050	ng	0.01
45) 2-Fluorobiphenyl	13.079	172	378313	91.713	ng	0.01
79) Terphenyl-d14	19.812	244	622795	86.312	ng	0.00
Target Compounds						
78) Pyrene	19.606	202	41244	4.654	ng	# 91

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

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