

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111925\
 Data File : BG064818.D
 Acq On : 19 Nov 2025 16:42
 Operator : RC/JU
 Sample : Q3530-07
 Misc : LOD-WATER-2.5PPM
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOD-MDL-WATER-01-QT4-2025

Quant Time: Nov 19 17:19:48 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	8.164	152	42956	20.000	ng	0.00
21) Naphthalene-d8	10.979	136	174592	20.000	ng	0.00
39) Acenaphthene-d10	14.774	164	145306	20.000	ng	0.00
64) Phenanthrene-d10	17.500	188	366976	20.000	ng	0.00
76) Chrysene-d12	21.748	240	452120	20.000	ng	0.00
86) Perylene-d12	25.009	264	439502	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.697	112	290064	117.915	ng	0.00
7) Phenol-d6	7.307	99	385769	117.906	ng	0.00
23) Nitrobenzene-d5	9.322	82	252943	71.631	ng	0.00
42) 2,4,6-Tribromophenol	16.249	330	337873	119.225	ng	0.00
45) 2-Fluorobiphenyl	13.405	172	737049	70.777	ng	0.00
79) Terphenyl-d14	20.086	244	1768980	76.965	ng	0.00
Target Compounds						
19) n-Nitroso-di-n-propyla...	8.963	70	4722	2.134	ng	Qvalue # 85

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

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