

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111925\
 Data File : BP026161.D
 Acq On : 19 Nov 2025 16:52
 Operator : RC/JU
 Sample : Q3530-07
 Misc : LOD-WATER-2.5ng
 ALS Vial : 9 Sample Multiplier: 1

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

Quant Time: Nov 19 17:13:10 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 19 01:25:31 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.666	152	227573	20.000	ng	0.00
21) Naphthalene-d8	10.431	136	918716	20.000	ng	-0.02
39) Acenaphthene-d10	14.307	164	585119	20.000	ng	-0.02
64) Phenanthrene-d10	17.119	188	1211850	20.000	ng	-0.01
76) Chrysene-d12	21.572	240	1394641	20.000	ng	0.00
86) Perylene-d12	24.895	264	1555051	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.284	112	1675775	118.193	ng	0.00
7) Phenol-d6	6.843	99	2122595	117.596	ng	-0.01
23) Nitrobenzene-d5	8.802	82	1185906	73.322	ng	-0.01
42) 2,4,6-Tribromophenol	15.825	330	901206	118.729	ng	-0.01
45) 2-Fluorobiphenyl	12.913	172	2931566	73.429	ng	-0.02
79) Terphenyl-d14	19.866	244	4803875	70.236	ng	0.00
Target Compounds						
19) n-Nitroso-di-n-propyla...	8.460	70	23843	2.126	ng	Qvalue 98

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

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