

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061025\
 Data File : VX046592.D
 Acq On : 10 Jun 2025 11:27
 Operator : JC/MD
 Sample : Q2262-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ARS20-0001

Quant Time: Jun 11 01:44:38 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060625W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 06 16:56:12 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.568	168	114799	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	225436	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	231982	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	110739	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	101949	49.916	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	99.840%		
35) Dibromofluoromethane	5.397	113	77958	47.024	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	94.040%		
50) Toluene-d8	8.653	98	300729	55.021	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	110.040%		
62) 4-Bromofluorobenzene	11.079	95	123194	54.477	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery =	108.960%		
Target Compounds					Qvalue	
16) Acetone	2.391	43	36562	49.344	ug/l	97
20) Methylene Chloride	2.806	84	4138	2.525	ug/l	94
25) 2-Butanone	4.592	43	6697m	5.890	ug/l	

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061025\
Data File : VX046592.D
Acq On : 10 Jun 2025 11:27
Operator : JC/MD
Sample : Q2262-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ARS20-0001

Quant Time: Jun 11 01:44:38 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060625W.M
Quant Title : SW846 8260
QLast Update : Fri Jun 06 16:56:12 2025
Response via : Initial Calibration

