

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092925\
 Data File : VV039236.D
 Acq On : 29 Sep 2025 16:53
 Operator : SY/MD
 Sample : Q3231-16
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 WC-4

Quant Time: Sep 30 01:28:50 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\82V092225W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 24 08:08:08 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	4.600	168	421100	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	5.535	114	839716	50.000	ug/l	0.00
63) Chlorobenzene-d5	8.776	117	896203	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.175	152	377074	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	4.944	65	369582	60.641	ug/l	0.00
Spiked Amount	50.000	Range	81 - 118	Recovery	=	121.280%#
35) Dibromofluoromethane	4.500	113	351028	51.999	ug/l	0.00
Spiked Amount	50.000	Range	80 - 119	Recovery	=	104.000%
50) Toluene-d8	7.236	98	944338	47.630	ug/l	0.00
Spiked Amount	50.000	Range	89 - 112	Recovery	=	95.260%
62) 4-Bromofluorobenzene	10.001	95	347955	42.631	ug/l	0.00
Spiked Amount	50.000	Range	85 - 114	Recovery	=	85.260%
Target Compounds						
16) Acetone	2.124	43	135663	30.285	ug/l	Qvalue 97
18) Methyl Acetate	2.384	43	19386	2.147	ug/l #	86
20) Methylene Chloride	2.462	84	43965	4.805	ug/l	87
43) Isopropyl Acetate	5.211	43	31524	1.867	ug/l	97

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092925\
Data File : VV039236.D
Acq On : 29 Sep 2025 16:53
Operator : SY/MD
Sample : Q3231-16
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
WC-4

Quant Time: Sep 30 01:28:50 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\82V092225W.M
Quant Title : SW846 8260
QLast Update : Wed Sep 24 08:08:08 2025
Response via : Initial Calibration

