

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092925\
 Data File : VV039234.D
 Acq On : 29 Sep 2025 16:08
 Operator : SY/MD
 Sample : Q3231-08
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 WC-2

Quant Time: Sep 30 01:27:59 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	4.603	168	432703	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	5.532	114	905184	50.000	ug/l	0.00
63) Chlorobenzene-d5	8.776	117	925002	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.172	152	393713	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	4.940	65	376726	60.156	ug/l	0.00
Spiked Amount	50.000	Range	81 - 118	Recovery	=	120.320%#
35) Dibromofluoromethane	4.503	113	360306	49.513	ug/l	0.00
Spiked Amount	50.000	Range	80 - 119	Recovery	=	99.020%
50) Toluene-d8	7.236	98	981084	45.904	ug/l	0.00
Spiked Amount	50.000	Range	89 - 112	Recovery	=	91.800%
62) 4-Bromofluorobenzene	10.001	95	358135	40.704	ug/l	0.00
Spiked Amount	50.000	Range	85 - 114	Recovery	=	81.400%#
Target Compounds						
					Qvalue	
16) Acetone	2.121	43	134393	29.029	ug/l	99
18) Methyl Acetate	2.388	43	20900	2.252	ug/l	95
20) Methylene Chloride	2.455	84	45472	4.849	ug/l	89
43) Isopropyl Acetate	5.214	43	29356	1.613	ug/l #	92

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092925\
Data File : VV039234.D
Acq On : 29 Sep 2025 16:08
Operator : SY/MD
Sample : Q3231-08
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
WC-2

Quant Time: Sep 30 01:27:59 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

