

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102425\
 Data File : VX048348.D
 Acq On : 24 Oct 2025 12:47
 Operator : JC/MD
 Sample : Q3412-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TP-6

Quant Time: Oct 24 23:03:39 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102025W.M
 Quant Title : SW846 8260
 QLast Update : Mon Oct 20 14:13:59 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.544	168	125875	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	252480	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	253174	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	119697	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	98047	55.555	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	111.120%
35) Dibromofluoromethane	5.373	113	68796	39.913	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	79.820%
50) Toluene-d8	8.635	98	283134	46.864	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	93.720%
62) 4-Bromofluorobenzene	11.061	95	118583	51.593	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	103.180%
Target Compounds						
					Qvalue	
16) Acetone	2.374	43	24886	41.103	ug/l	94
18) Methyl Acetate	2.703	43	3556	2.812	ug/l	95
20) Methylene Chloride	2.788	84	3271	2.217	ug/l	93
43) Isopropyl Acetate	6.324	43	9978	2.991	ug/l	96

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102425\
Data File : VX048348.D
Acq On : 24 Oct 2025 12:47
Operator : JC/MD
Sample : Q3412-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TP-6

Quant Time: Oct 24 23:03:39 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102025W.M
Quant Title : SW846 8260
QLast Update : Mon Oct 20 14:13:59 2025
Response via : Initial Calibration

