

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101325\
 Data File : VX048144.D
 Acq On : 13 Oct 2025 15:51
 Operator : JC/MD
 Sample : Q3326-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TP-1

Quant Time: Oct 14 01:56:28 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091625W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 17 06:39:58 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.544	168	111228	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	230199	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	247574	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	123164	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	95861	54.144	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 108.280%	
35) Dibromofluoromethane	5.379	113	77129	49.959	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 99.920%	
50) Toluene-d8	8.635	98	273065	51.117	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 102.240%	
62) 4-Bromofluorobenzene	11.061	95	114812	56.631	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 113.260%	
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	33860	43.985	ug/l	99
18) Methyl Acetate	2.709	43	4257	2.485	ug/l	99
20) Methylene Chloride	2.794	84	6395	3.925	ug/l	99
43) Isopropyl Acetate	6.336	43	10920	2.757	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101325\
Data File : VX048144.D
Acq On : 13 Oct 2025 15:51
Operator : JC/MD
Sample : Q3326-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TP-1

Quant Time: Oct 14 01:56:28 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091625W.M
Quant Title : SW846 8260
QLast Update : Wed Sep 17 06:39:58 2025
Response via : Initial Calibration

