

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100625\
 Data File : VX048023.D
 Acq On : 06 Oct 2025 14:49
 Operator : JC/MD
 Sample : Q3263-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 251818

Quant Time: Oct 07 02:14:44 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	5.537	168	112174	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	225184	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	236785	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	118391	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	90992	50.961	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	101.920%
35) Dibromofluoromethane	5.379	113	74680	49.450	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.900%
50) Toluene-d8	8.634	98	253923	48.592	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.180%
62) 4-Bromofluorobenzene	11.067	95	109125	55.024	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	110.040%
Target Compounds						Qvalue
16) Acetone	2.379	43	2427	3.126	ug/l #	87
30) Chloroform	5.086	83	4268	1.491	ug/l	89
47) Bromodichloromethane	7.811	83	4455	1.729	ug/l	95
60) Dibromochloromethane	9.506	129	7005	3.818	ug/l	98
71) Bromoform	10.786	173	3741	2.699	ug/l #	98

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

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