

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV100625\
 Data File : VV039275.D
 Acq On : 06 Oct 2025 15:48
 Operator : SY/MD
 Sample : Q3269-06
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 EG1-TP2

Quant Time: Oct 07 03:36:37 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	447623	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	5.532	114	881706	50.000	ug/l	0.00
63) Chlorobenzene-d5	8.773	117	945653	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.172	152	434975	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	4.940	65	366605	56.589	ug/l	0.00
Spiked Amount	50.000	Range	81 - 118	Recovery	=	113.180%
35) Dibromofluoromethane	4.503	113	364819	51.468	ug/l	0.00
Spiked Amount	50.000	Range	80 - 119	Recovery	=	102.940%
50) Toluene-d8	7.236	98	951448	45.703	ug/l	0.00
Spiked Amount	50.000	Range	89 - 112	Recovery	=	91.400%
62) 4-Bromofluorobenzene	10.001	95	401720	46.874	ug/l	0.00
Spiked Amount	50.000	Range	85 - 114	Recovery	=	93.740%
Target Compounds						
						Qvalue
16) Acetone	2.124	43	65020	11.175	ug/l	98
18) Methyl Acetate	2.387	43	21692	2.260	ug/l	95
43) Isopropyl Acetate	5.217	43	37289	2.103	ug/l	97

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

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