

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV102125\
 Data File : VV039297.D
 Acq On : 21 Oct 2025 12:00
 Operator : SY/MD
 Sample : Q3386-04
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 3444

Quant Time: Oct 21 12:37:36 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\82V100725W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 08 05:01:26 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	4.600	168	1014747	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	5.532	114	1869889	50.000	ug/l	0.00
63) Chlorobenzene-d5	8.773	117	1729629	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.172	152	880932	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	4.941	65	732497	50.801	ug/l	0.00
Spiked Amount	50.000	Range	81 - 118	Recovery	=	101.600%
35) Dibromofluoromethane	4.500	113	652945	44.317	ug/l	0.00
Spiked Amount	50.000	Range	80 - 119	Recovery	=	88.640%
50) Toluene-d8	7.236	98	2244819	44.787	ug/l	0.00
Spiked Amount	50.000	Range	89 - 112	Recovery	=	89.580%
62) 4-Bromofluorobenzene	9.998	95	827110	46.335	ug/l	0.00
Spiked Amount	50.000	Range	85 - 114	Recovery	=	92.680%
Target Compounds						
					Qvalue	
16) Acetone	2.121	43	228800	24.882	ug/l	97
17) Carbon Disulfide	2.256	76	11440	0.296	ug/l #	90
30) Chloroform	4.285	83	29210	1.098	ug/l	90
52) Toluene	7.310	92	467116	13.028	ug/l	99
67) Ethyl Benzene	8.947	91	35984	0.597	ug/l	97
68) m/p-Xylenes	9.066	106	49401	2.065	ug/l	98
69) o-Xylene	9.474	106	20105	0.952	ug/l	99

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

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