

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV102125\
 Data File : VV039296.D
 Acq On : 21 Oct 2025 11:38
 Operator : SY/MD
 Sample : Q3386-03
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 3443

Quant Time: Oct 21 12:37:51 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	902813	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	5.532	114	1732513	50.000	ug/l	0.00
63) Chlorobenzene-d5	8.773	117	1552214	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.172	152	805600	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	4.941	65	670925	52.300	ug/l	0.00
Spiked Amount	50.000	Range	81 - 118	Recovery	=	104.600%
35) Dibromofluoromethane	4.503	113	598709	43.858	ug/l	0.00
Spiked Amount	50.000	Range	80 - 119	Recovery	=	87.720%
50) Toluene-d8	7.236	98	2047012	44.079	ug/l	0.00
Spiked Amount	50.000	Range	89 - 112	Recovery	=	88.160%#
62) 4-Bromofluorobenzene	9.998	95	725044	43.838	ug/l	0.00
Spiked Amount	50.000	Range	85 - 114	Recovery	=	87.680%

Target Compounds						Qvalue
16) Acetone	2.118	43	886121	144.300	ug/l	100
30) Chloroform	4.282	83	366236	15.472	ug/l	100
43) Isopropyl Acetate	5.439	43	484664	18.443	ug/l #	60
52) Toluene	7.326	92	10034	0.302	ug/l	100
67) Ethyl Benzene	8.953	91	19145	0.354	ug/l	96
68) m/p-Xylenes	9.069	106	23228	1.082	ug/l	96
69) o-Xylene	9.474	106	15921	0.840	ug/l	95
84) 1,2,4-Trimethylbenzene	10.841	105	88328	2.187	ug/l	97

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

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