

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070623\
 Data File : VX036431.D
 Acq On : 06 Jul 2023 13:30
 Operator : JC/MD
 Sample : 03509-15
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WC-WWS-03-G

Quant Time: Jul 07 01:46:48 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062323W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 23 15:03:28 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	174353	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	295872	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	268001	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	121538	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	113300	45.334	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	90.660%
35) Dibromofluoromethane	5.385	113	91466	46.884	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.760%
50) Toluene-d8	8.653	98	362028	49.684	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.360%
62) 4-Bromofluorobenzene	11.079	95	124733	48.599	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	97.200%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	232142	217.312	ug/l	97
20) Methylene Chloride	2.794	84	13051	5.674	ug/l #	94
25) 2-Butanone	4.562	43	117207	76.364	ug/l	94
37) Ethyl Acetate	4.720	43	40885	13.794	ug/l	97
43) Isopropyl Acetate	6.342	43	146952	31.079	ug/l	98
51) 4-Methyl-2-Pentanone	8.573	43	461931	151.638	ug/l	89

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

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