

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070623\
 Data File : VX036429.D
 Acq On : 06 Jul 2023 12:42
 Operator : JC/MD
 Sample : 03509-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WC-WWS-01-G

Quant Time: Jul 07 01:46:00 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	174892	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	295058	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	268044	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	120311	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	112958	45.058	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	90.120%
35) Dibromofluoromethane	5.391	113	91487	47.024	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.040%
50) Toluene-d8	8.653	98	358160	49.289	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	98.580%
62) 4-Bromofluorobenzene	11.079	95	126840	49.556	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	99.120%
Target Compounds						
						Qvalue
16) Acetone	2.392	43	238658	222.723	ug/l	91
20) Methylene Chloride	2.788	84	9183	3.980	ug/l	92
25) 2-Butanone	4.562	43	122479	79.553	ug/l	95
37) Ethyl Acetate	4.714	43	35067	11.864	ug/l	98
43) Isopropyl Acetate	6.342	43	123042	26.094	ug/l	98
51) 4-Methyl-2-Pentanone	8.573	43	461774	152.005	ug/l	89

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

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