

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071023\
 Data File : VX036488.D
 Acq On : 10 Jul 2023 17:19
 Operator : JC/MD
 Sample : 03541-10
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WC-WWS-09-G

Quant Time: Jul 11 01:31:38 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	5.550	168	167709	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	288700	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	267018	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	121501	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	112356	46.737	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	93.480%
35) Dibromofluoromethane	5.385	113	90027	47.293	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.580%
50) Toluene-d8	8.647	98	357781	50.321	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.640%
62) 4-Bromofluorobenzene	11.079	95	124964	49.898	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	99.800%
Target Compounds						
					Qvalue	
16) Acetone	2.379	43	248589	241.928	ug/l	94
18) Methyl Acetate	2.703	43	4573	1.082	ug/l	96
20) Methylene Chloride	2.788	84	10158	4.591	ug/l	95
25) 2-Butanone	4.556	43	114996	77.892	ug/l	95
37) Ethyl Acetate	4.721	43	20788	7.188	ug/l	98
43) Isopropyl Acetate	6.336	43	81848	17.740	ug/l	97
51) 4-Methyl-2-Pentanone	8.573	43	442818	148.975	ug/l	92

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

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