

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071023\
 Data File : VX036495.D
 Acq On : 10 Jul 2023 20:04
 Operator : JC/MD
 Sample : 03541-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WC-WWS-08-G

Quant Time: Jul 11 01:34:08 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	178716	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	307658	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	289867	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	131000	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	117380	45.820	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	91.640%
35) Dibromofluoromethane	5.385	113	97348	47.987	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.980%
50) Toluene-d8	8.647	98	382069	50.426	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.860%
62) 4-Bromofluorobenzene	11.079	95	136140	51.011	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	102.020%
Target Compounds						
					Qvalue	
16) Acetone	2.379	43	238191	217.531	ug/l	93
20) Methylene Chloride	2.788	84	9722	4.124	ug/l	93
25) 2-Butanone	4.562	43	110898	70.489	ug/l	95
37) Ethyl Acetate	4.720	43	18110	5.876	ug/l	99
43) Isopropyl Acetate	6.336	43	68462	13.925	ug/l	98
51) 4-Methyl-2-Pentanone	8.573	43	425820	134.429	ug/l	93

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

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