

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031723\
 Data File : VX034617.D
 Acq On : 18 Mar 2023 07:15
 Operator : JC/MD
 Sample : 01964-11
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MAPLE-6

Quant Time: Mar 20 09:00:44 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
 Quant Title : SW846 8260
 QLast Update : Mon Mar 13 10:29:32 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	228099	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	392600	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	401646	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	185040	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	160065	48.236	ug/l	0.01
Spiked Amount	50.000	Range	78 - 117	Recovery	=	96.480%
35) Dibromofluoromethane	5.385	113	122920	47.420	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.840%
50) Toluene-d8	8.647	98	520242	53.763	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	107.520%
62) 4-Bromofluorobenzene	11.079	95	238094	60.915	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	121.840%
Target Compounds						
					Qvalue	
16) Acetone	2.373	43	54516	34.876	ug/l	95
18) Methyl Acetate	2.703	43	5931	1.161	ug/l	99
20) Methylene Chloride	2.788	84	19643	6.475	ug/l	95
25) 2-Butanone	4.568	43	11936	5.337	ug/l	91
37) Ethyl Acetate	4.714	43	42108	9.681	ug/l	99
43) Isopropyl Acetate	6.342	43	176628	23.971	ug/l	99

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

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