

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX110920\
 Data File : VX019371.D
 Acq On : 09 Nov 2020 16:28
 Operator : JC/MD
 Sample : L4637-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TRE-1059

Quant Time: Nov 10 02:51:15 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102720W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 27 13:55:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	297488	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	527813	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	509744	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	252126	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	210903	54.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.94%	
35) Dibromofluoromethane	5.46	113	163839	49.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.84%	
50) Toluene-d8	8.70	98	658486	52.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.38%	
62) 4-Bromofluorobenzene	11.12	95	253696	53.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.20%	
Target Compounds						
16) Acetone	2.42	43	14581	11.017	ug/l	100
18) Methyl Acetate	2.75	43	3553	1.104	ug/l #	89
20) Methylene Chloride	2.84	84	9990	3.007	ug/l	96
43) Isopropyl Acetate	6.41	43	74006	9.767	ug/l	99

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

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