

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102521\
 Data File : VU045383.D
 Acq On : 25 Oct 2021 14:39
 Operator : SY/MD
 Sample : M4290-11
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 TH-BBS

Quant Time: Oct 26 04:42:47 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 28 12:32:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.375	168	218213	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	417181	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.426	117	385884	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	163332	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	177615	48.897	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	97.800%
35) Dibromofluoromethane	5.292	113	128063	48.661	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	97.320%
50) Toluene-d8	7.902	98	520760	50.081	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.160%
62) 4-Bromofluorobenzene	10.642	95	186574	43.460	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	86.920%
Target Compounds						
					Qvalue	
16) Acetone	2.652	43	13197	6.121	ug/l	99
18) Methyl Acetate	2.960	43	8628	1.369	ug/l	94
20) Methylene Chloride	3.038	84	11435	3.521	ug/l	94
43) Isopropyl Acetate	5.922	43	31847	3.667	ug/l	99

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

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