

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102521\
 Data File : VU045384.D
 Acq On : 25 Oct 2021 15:03
 Operator : SY/MD
 Sample : M4290-14
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 TH-BBD

Quant Time: Oct 26 04:43:08 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	5.372	168	216816	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	413720	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.427	117	385162	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	166024	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	175704	48.683	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	97.360%
35) Dibromofluoromethane	5.292	113	127448	48.833	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	97.660%
50) Toluene-d8	7.902	98	515807	50.020	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.040%
62) 4-Bromofluorobenzene	10.642	95	185843	43.652	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	87.300%
Target Compounds						
					Qvalue	
16) Acetone	2.645	43	16143	7.535	ug/l	96
18) Methyl Acetate	2.957	43	9075	1.449	ug/l	94
20) Methylene Chloride	3.041	84	11342	3.515	ug/l	94
43) Isopropyl Acetate	5.922	43	31187	3.621	ug/l	99

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

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