

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102521\
 Data File : VU045392.D
 Acq On : 25 Oct 2021 18:39
 Operator : SY/MD
 Sample : M4300-08
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 TH-CCS

Quant Time: Oct 26 04:45:47 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	218093	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	414906	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	384597	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	171406	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.707	65	177676	48.941	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	97.880%
35) Dibromofluoromethane	5.292	113	128973	49.276	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.560%
50) Toluene-d8	7.903	98	522293	50.504	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	101.000%
62) 4-Bromofluorobenzene	10.636	95	185915	43.544	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	87.080%
Target Compounds						
					Qvalue	
16) Acetone	2.639	43	11915	5.529	ug/l	99
20) Methylene Chloride	3.051	84	11441	3.525	ug/l	94
43) Isopropyl Acetate	5.916	43	28634	3.315	ug/l	98

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

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