

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100121\
 Data File : VU045078.D
 Acq On : 01 Oct 2021 17:12
 Operator : SY/MD
 Sample : M3823-02
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DKM-ENV-IC

Quant Time: Oct 02 03:12:52 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	184769	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	341711	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	329575	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	151099	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.706	65	156118	50.758	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	101.520%
35) Dibromofluoromethane	5.295	113	106801	49.545	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.100%
50) Toluene-d8	7.899	98	440530	51.722	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	103.440%
62) 4-Bromofluorobenzene	10.636	95	163575	46.518	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	93.040%
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	71314	39.062	ug/l	99
18) Methyl Acetate	2.951	43	21492	4.028	ug/l	100
20) Methylene Chloride	3.051	84	11404	4.147	ug/l	92
25) 2-Butanone	4.716	43	15685	6.047	ug/l	100
43) Isopropyl Acetate	5.915	43	23797	3.345	ug/l	98

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

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