

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100121\
 Data File : VU045079.D
 Acq On : 01 Oct 2021 17:35
 Operator : SY/MD
 Sample : M3951-02
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 WC-20210924

Quant Time: Oct 02 03:13:04 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	188423	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	348707	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	342225	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	159963	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.706	65	159316	50.794	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 101.580%			
35) Dibromofluoromethane	5.292	113	107742	48.979	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 97.960%			
50) Toluene-d8	7.902	98	449627	51.731	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 103.460%			
62) 4-Bromofluorobenzene	10.639	95	174947	48.754	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 97.500%			
Target Compounds					Qvalue	
16) Acetone	2.639	43	74958	40.261	ug/l	100
18) Methyl Acetate	2.954	43	6402	1.177	ug/l #	89
20) Methylene Chloride	3.047	84	17539	6.254	ug/l	97
25) 2-Butanone	4.716	43	16981	6.420	ug/l	97
43) Isopropyl Acetate	5.919	43	23285	3.208	ug/l	99

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

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