

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
 Data File : VU045030.D
 Acq On : 30 Sep 2021 06:36
 Operator : SY/MD
 Sample : M3971-05
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 RO-COMP-2

Quant Time: Sep 30 08:29:30 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	5.379	168	199819	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	361426	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	346283	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	157842	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.704	65	161286	48.49	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	96.98%		
35) Dibromofluoromethane	5.292	113	112375	49.29	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	98.58%		
50) Toluene-d8	7.903	98	463108	51.41	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	102.82%		
62) 4-Bromofluorobenzene	10.636	95	171004	45.98	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	91.96%		
Target Compounds					Qvalue	
16) Acetone	2.630	43	38088	19.29	ug/l	98
17) Carbon Disulfide	2.790	76	19663	2.81	ug/l	99
18) Methyl Acetate	2.954	43	6465	1.12	ug/l	96
20) Methylene Chloride	3.048	84	10501	3.53	ug/l	94
43) Isopropyl Acetate	5.916	43	20413	2.71	ug/l	100

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

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