

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100121\
 Data File : VU045075.D
 Acq On : 01 Oct 2021 16:02
 Operator : SY/MD
 Sample : M3999-04
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 219

Quant Time: Oct 02 03:12:12 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	196819	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	367492	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	351758	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	158969	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	165746	50.589	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	101.180%
35) Dibromofluoromethane	5.292	113	112290	48.437	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.880%
50) Toluene-d8	7.899	98	461610	50.395	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.800%
62) 4-Bromofluorobenzene	10.639	95	175123	46.308	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	92.620%
Target Compounds						
					Qvalue	
16) Acetone	2.639	43	16462	8.465	ug/l	94
18) Methyl Acetate	2.957	43	7157	1.259	ug/l	96
20) Methylene Chloride	3.047	84	20489	6.995	ug/l	98
25) 2-Butanone	4.719	43	18311	6.627	ug/l	96
43) Isopropyl Acetate	5.922	43	24711	3.230	ug/l	98

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

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