

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092521\
 Data File : VU044873.D
 Acq On : 26 Sep 2021 06:15
 Operator : SY/MD
 Sample : M3888-09
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 40456

Quant Time: Sep 27 03:17:59 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092421W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 25 09:30:46 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.379	168	208609	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	364455	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	336549	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	154475	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.707	65	166378	45.698	ug/l	0.00
Spiked Amount	50.000		Recovery	=	91.400%	
35) Dibromofluoromethane	5.292	113	115252	46.097	ug/l	0.00
Spiked Amount	50.000		Recovery	=	92.200%	
50) Toluene-d8	7.903	98	462231	48.951	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.900%	
62) 4-Bromofluorobenzene	10.636	95	165482	47.183	ug/l	0.00
Spiked Amount	50.000		Recovery	=	94.360%	
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	2437	1.187	ug/l	98
18) Methyl Acetate	2.961	43	2995	0.437	ug/l	92
19) Methyl tert-butyl Ether	3.369	73	6041	0.644	ug/l	98
22) Diisopropyl ether	4.002	45	4227	0.439	ug/l #	72
29) Tetrahydrofuran	5.054	42	127360	64.968	ug/l	98
30) Chloroform	5.096	83	2884	0.513	ug/l	92
95) Naphthalene	14.092	128	12913	2.012	ug/l	99

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

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