

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061021\
 Data File : VU044157.D
 Acq On : 10 Jun 2021 15:47
 Operator : SY/MD
 Sample : M2644-02
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 1894-1900

Quant Time: Jun 11 01:20:43 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U060921W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 10 05:11:37 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.382	168	142497	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	296820	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	308561	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	143267	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	127561	56.292	ug/l	0.00
Spiked Amount	50.000		Recovery	=	112.580%	
35) Dibromofluoromethane	5.298	113	102135	49.466	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.940%	
50) Toluene-d8	7.906	98	403268	50.856	ug/l	0.00
Spiked Amount	50.000		Recovery	=	101.720%	
62) 4-Bromofluorobenzene	10.639	95	147891	46.539	ug/l	0.00
Spiked Amount	50.000		Recovery	=	93.080%	
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	13167	9.191	ug/l	99
20) Methylene Chloride	3.054	84	12663	5.125	ug/l	97
43) Isopropyl Acetate	5.922	43	15323	2.731	ug/l	99
51) 4-Methyl-2-Pentanone	7.800	43	4251	1.047	ug/l	95

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

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