

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050621\
 Data File : VU043589.D
 Acq On : 06 May 2021 17:30
 Operator : SY/MD
 Sample : M2267-03
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DI-WATER

Quant Time: May 07 02:06:25 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U050621W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 06 07:26:23 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.382	168	113208	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	201547	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	197439	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	87095	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	118197	51.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.16%	
35) Dibromofluoromethane	5.295	113	80960	50.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.74%	
50) Toluene-d8	7.906	98	284063	49.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.68%	
62) 4-Bromofluorobenzene	10.639	95	98164	45.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.54%	
Target Compounds						
						Qvalue
16) Acetone	2.633	43	10890	8.56	ug/l	99
17) Carbon Disulfide	2.800	76	2029	0.49	ug/l	98
30) Chloroform	5.096	83	14174	4.50	ug/l	97
65) Chlorobenzene	9.456	112	1631	0.37	ug/l	96

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

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