

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110221\
 Data File : VU045496.D
 Acq On : 01 Nov 2021 18:53
 Operator : SY/MD
 Sample : M4431-03
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BP-VPB-RW9-GW-358-360

Quant Time: Nov 02 07:36:12 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U102821W.M
 Quant Title : SW846 8260
 QLast Update : Fri Oct 29 11:09:42 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.378	168	192444	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	380483	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	369283	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	180578	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.706	65	161860	51.662	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 103.320%	
35) Dibromofluoromethane	5.295	113	121492	50.218	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 100.440%	
50) Toluene-d8	7.902	98	487959	51.435	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 102.860%	
62) 4-Bromofluorobenzene	10.635	95	185834	51.405	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 102.800%	
Target Compounds						
					Qvalue	
3) Chloromethane	1.523	50	939	0.245	ug/l	90
16) Acetone	2.642	43	4311	3.031	ug/l	98
17) Carbon Disulfide	2.797	76	1737	0.271	ug/l #	75

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

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