

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110221\
 Data File : VU045513.D
 Acq On : 02 Nov 2021 14:49
 Operator : SY/MD
 Sample : M4431-02RE
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BP-VPB-RW9-GW-338-340RE

Quant Time: Nov 02 15:16:53 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	182640	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	362996	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	372013	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	183075	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.706	65	136861	46.028	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	92.060%
35) Dibromofluoromethane	5.295	113	110853	48.028	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.060%
50) Toluene-d8	7.902	98	489916	54.129	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	108.260%
62) 4-Bromofluorobenzene	10.635	95	187864	54.470	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	108.940%
Target Compounds						
16) Acetone	2.642	43	67664	50.125	ug/l	99
25) 2-Butanone	4.716	43	15828	7.767	ug/l	98

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

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