

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062722\
 Data File : VU049439.D
 Acq On : 28 Jun 2022 06:12
 Operator : SY/MD
 Sample : N3482-02
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-37

Quant Time: Jun 28 07:16:23 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 28 05:14:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.372	168	300551	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	514210	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	466351	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.813	152	221180	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.704	65	197939	42.764	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	85.520%
35) Dibromofluoromethane	5.289	113	161204	44.809	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	89.620%
50) Toluene-d8	7.896	98	620761	48.378	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	96.760%
62) 4-Bromofluorobenzene	10.633	95	229195	46.629	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	93.260%
Target Compounds						
					Qvalue	
16) Acetone	2.639	43	4842	2.074	ug/l	95
17) Carbon Disulfide	2.794	76	2588	0.254	ug/l	99
51) 4-Methyl-2-Pentanone	7.797	43	9293	1.329	ug/l	92

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

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