

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062722\
 Data File : VU049438.D
 Acq On : 28 Jun 2022 05:49
 Operator : SY/MD
 Sample : N3482-01
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-35

Quant Time: Jun 28 07:16:06 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	304464	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.250	114	523298	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	464463	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.813	152	210624	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	201358	42.943	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	85.880%
35) Dibromofluoromethane	5.289	113	163522	44.664	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	89.320%
50) Toluene-d8	7.896	98	622943	47.705	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	95.400%
62) 4-Bromofluorobenzene	10.632	95	221874	44.355	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	88.720%
Target Compounds						
						Qvalue
11) Tert butyl alcohol	3.199	59	141330	128.775	ug/l	# 53
19) Methyl tert-butyl Ether	3.366	73	42034	3.418	ug/l	98
22) Diisopropyl ether	3.993	45	7495	0.587	ug/l	# 50
30) Chloroform	5.086	83	10822	1.493	ug/l	97
47) Bromodichloromethane	7.105	83	3206	0.563	ug/l	89
60) Dibromochloromethane	8.816	129	1708	0.417	ug/l	89

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

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