

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062922\
 Data File : VU049507.D
 Acq On : 29 Jun 2022 13:03
 Operator : SY/MD
 Sample : N3482-01
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-35

Quant Time: Jun 30 00:48:13 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 28 10:35:43 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.375	168	280713	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	491908	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	443947	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	203011	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	191654	44.332	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	88.660%
35) Dibromofluoromethane	5.288	113	155197	45.096	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	90.200%
50) Toluene-d8	7.896	98	585518	47.700	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	95.400%
62) 4-Bromofluorobenzene	10.632	95	211812	45.046	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	90.100%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	3.199	59	141546	139.885	ug/l #	54
19) Methyl tert-butyl Ether	3.363	73	40291	3.553	ug/l	99
22) Diisopropyl ether	3.996	45	7344	0.624	ug/l #	53
30) Chloroform	5.089	83	9627	1.440	ug/l	99
47) Bromodichloromethane	7.108	83	3430	0.640	ug/l	91
60) Dibromochloromethane	8.809	129	1428	0.371	ug/l	85

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

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