

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062922\
 Data File : VU049543.D
 Acq On : 30 Jun 2022 04:19
 Operator : SY/MD
 Sample : N3517-04
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-4-20220627

Quant Time: Jun 30 04:40:57 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	5.372	168	239594	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	425573	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	381894	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	174384	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	170613	46.238	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	92.480%
35) Dibromofluoromethane	5.288	113	135294	45.440	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	90.880%
50) Toluene-d8	7.896	98	509375	47.965	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	95.940%
62) 4-Bromofluorobenzene	10.632	95	178397	43.853	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	87.700%
Target Compounds						
					Qvalue	
16) Acetone	2.642	43	3514	1.888	ug/l	98
30) Chloroform	5.089	83	64564	11.316	ug/l	99
64) Tetrachloroethene	8.552	164	2509	1.024	ug/l	95

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

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