

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062722\
 Data File : VU049430.D
 Acq On : 28 Jun 2022 02:40
 Operator : SY/MD
 Sample : N3386-04
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW181S-20220614

Quant Time: Jun 28 07:13:47 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	309060	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.250	114	532001	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	469098	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	212387	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	205556	43.187	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	86.380%		
35) Dibromofluoromethane	5.289	113	164912	44.307	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	88.620%		
50) Toluene-d8	7.899	98	626799	47.215	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	94.420%		
62) 4-Bromofluorobenzene	10.632	95	227402	44.717	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	89.440%		
Target Compounds						
						Qvalue
12) 1,1-Dichloroethene	2.581	96	1526	0.455	ug/l	90
16) Acetone	2.636	43	4121	1.717	ug/l #	84
24) 1,1-Dichloroethane	3.874	63	3989	0.536	ug/l #	93
27) cis-1,2-Dichloroethene	4.671	96	3367	0.805	ug/l	97
30) Chloroform	5.089	83	3540	0.481	ug/l	93
44) Trichloroethene	6.539	130	4538	1.172	ug/l	88
64) Tetrachloroethene	8.552	164	4089	1.359	ug/l #	90

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

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