

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062922\
 Data File : VX029891.D
 Acq On : 30 Jun 2022 04:03
 Operator : JC/MD
 Sample : N3481-17
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-31

Quant Time: Jun 30 06:27:58 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062822W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 29 05:54:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	231648	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	400994	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	334347	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	132499	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	129063	44.015	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	88.020%
35) Dibromofluoromethane	5.391	113	115096	45.928	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	91.860%
50) Toluene-d8	8.653	98	455469	48.448	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	96.900%
62) 4-Bromofluorobenzene	11.079	95	143414	43.156	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	86.320%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	18537	12.922	ug/l	96
25) 2-Butanone	4.574	43	10591	4.254	ug/l	99
27) cis-1,2-Dichloroethene	4.501	96	6255	2.025	ug/l	94
44) Trichloroethene	7.129	130	3100	1.068	ug/l	84
51) 4-Methyl-2-Pentanone	8.574	43	28756	6.353	ug/l	99
64) Tetrachloroethene	9.275	164	6290	3.119	ug/l	98

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

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