

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX083122\
 Data File : VX030971.D
 Acq On : 31 Aug 2022 12:25
 Operator : JC/MD
 Sample : N4381-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-10

Quant Time: Sep 01 00:55:59 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X083022W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 30 14:24:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	215379	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	363139	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	327104	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	155981	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	149792	48.490	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	96.980%
35) Dibromofluoromethane	5.391	113	137114	51.222	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.440%
50) Toluene-d8	8.653	98	514507	51.900	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	103.800%
62) 4-Bromofluorobenzene	11.085	95	177231	50.204	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	100.400%
Target Compounds						
16) Acetone	2.386	43	3249	2.330	ug/l	90
17) Carbon Disulfide	2.514	76	2355	0.341	ug/l	96
51) 4-Methyl-2-Pentanone	8.580	43	5306	1.239	ug/l	96

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

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