

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032222\
 Data File : VX027639.D
 Acq On : 23 Mar 2022 02:09
 Operator : JC/MD
 Sample : N2062-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-4

Quant Time: Mar 23 07:41:53 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031722W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 17 15:02:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	358019	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	575622	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	495620	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	229244	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	196223	44.491	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	88.980%
35) Dibromofluoromethane	5.391	113	186515	50.363	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	100.720%
50) Toluene-d8	8.653	98	648396	46.267	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	92.540%
62) 4-Bromofluorobenzene	11.079	95	225721	43.821	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	87.640%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	7977	4.437	ug/l	90
17) Carbon Disulfide	2.514	76	7135	0.843	ug/l #	91
19) Methyl tert-butyl Ether	3.117	73	24829	2.214	ug/l	99
51) 4-Methyl-2-Pentanone	8.580	43	16589	3.202	ug/l	96

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

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