

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032822\
 Data File : VX027754.D
 Acq On : 28 Mar 2022 16:55
 Operator : JC/MD
 Sample : N2154-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-5

Quant Time: Mar 29 05:39:55 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032522W.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 25 16:26:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	68285	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	116633	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	104128	50.000	ug/l	# 0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	43813	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	46422	41.449	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	82.900%
35) Dibromofluoromethane	5.385	113	38522	50.295	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	100.600%
50) Toluene-d8	8.647	98	142048	47.977	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	95.960%
62) 4-Bromofluorobenzene	11.079	95	53816	44.795	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	89.580%
Target Compounds						
					Qvalue	
3) Chloromethane	1.295	50	1578	1.923	ug/l	98
5) Bromomethane	1.605	94	1713	5.412	ug/l	89
16) Acetone	2.380	43	103430	218.365	ug/l	96
25) 2-Butanone	4.556	43	13661	17.959	ug/l	# 88
51) 4-Methyl-2-Pentanone	8.574	43	7470	5.117	ug/l	94
52) Toluene	8.720	92	762	0.365	ug/l	86
59) 2-Hexanone	9.439	43	1268	1.111	ug/l	91

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

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