

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
 Data File : VX028652.D
 Acq On : 12 May 2022 20:45
 Operator : JC/MD
 Sample : N2799-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-8BR

Quant Time: May 13 02:49:42 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 12 14:58:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	245729	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	453559	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	437995	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	197014	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	178543	53.006	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 106.020%	
35) Dibromofluoromethane	5.391	113	148624	50.014	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 100.020%	
50) Toluene-d8	8.653	98	548913	48.994	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 97.980%	
62) 4-Bromofluorobenzene	11.079	95	206088	49.249	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 98.500%	
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	3247	2.246	ug/l #	86
19) Methyl tert-butyl Ether	3.117	73	4182	0.463	ug/l #	89
20) Methylene Chloride	2.794	84	960	0.317	ug/l #	81
30) Chloroform	5.099	83	8475	1.545	ug/l	92
68) m/p-Xylenes	10.305	106	3880	0.598	ug/l	100
84) 1,2,4-Trimethylbenzene	11.756	105	3045	0.244	ug/l	95

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

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