

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042022\
 Data File : VX028196.D
 Acq On : 20 Apr 2022 14:39
 Operator : JC/MD
 Sample : N2448-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 E14ST-B6-041522-GW

Quant Time: Apr 21 06:12:54 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041922W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 19 13:38:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	400972	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	685675	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	644086	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	299461	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	284677	54.450	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	108.900%
35) Dibromofluoromethane	5.391	113	262313	54.551	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	109.100%
50) Toluene-d8	8.653	98	940624	50.930	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	101.860%
62) 4-Bromofluorobenzene	11.079	95	343759	48.679	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	97.360%
Target Compounds						
					Qvalue	
16) Acetone	2.374	43	8695	4.072	ug/l	91
20) Methylene Chloride	2.788	84	1395	0.360	ug/l #	85
30) Chloroform	5.099	83	69271	8.295	ug/l	95
47) Bromodichloromethane	7.824	83	8312	1.227	ug/l	96
60) Dibromochloromethane	9.531	129	2300	0.407	ug/l	88

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

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