

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042922\
 Data File : VX028431.D
 Acq On : 29 Apr 2022 18:09
 Operator : JC/MD
 Sample : N2640-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MWBR3-B-117-137-042922

Quant Time: Apr 30 09:16:40 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	390066	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	674233	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	652767	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	311304	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	282201	55.486	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 110.980%	
35) Dibromofluoromethane	5.391	113	254115	53.743	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 107.480%	
50) Toluene-d8	8.653	98	947466	52.171	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 104.340%	
62) 4-Bromofluorobenzene	11.079	95	344458	49.606	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 99.220%	
Target Compounds						
						Qvalue
12) 1,1-Dichloroethene	2.325	96	190888	63.677	ug/l	95
16) Acetone	2.379	43	8437	4.058	ug/l	98
20) Methylene Chloride	2.788	84	1383	0.366	ug/l #	74
21) trans-1,2-Dichloroethene	3.093	96	13739	3.370	ug/l	96
24) 1,1-Dichloroethane	3.611	63	31194	4.611	ug/l	97
27) cis-1,2-Dichloroethene	4.489	96	129482	26.636	ug/l	85
30) Chloroform	5.092	83	2567	0.316	ug/l	93
42) 1,2-Dichloroethane	6.086	62	19449	2.902	ug/l	100
44) Trichloroethene	7.129	130	811557	155.010	ug/l	99
49) 1,4-Dioxane	7.671	88	2516	20.112	ug/l	99
52) Toluene	8.720	92	3745	0.299	ug/l	92
64) Tetrachloroethene	9.275	164	4326	0.848	ug/l	90

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

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