

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031922\
 Data File : VX027560.D
 Acq On : 19 Mar 2022 15:10
 Operator : JC/MD
 Sample : N2000-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE-132D1-20220315

Quant Time: Mar 21 01:22:34 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	337072	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	551092	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	498263	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	220698	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	199784	48.113	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	96.220%
35) Dibromofluoromethane	5.391	113	184510	52.039	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	104.080%
50) Toluene-d8	8.653	98	650013	48.447	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	96.900%
62) 4-Bromofluorobenzene	11.085	95	225764	45.780	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	91.560%
Target Compounds						
					Qvalue	
2) Dichlorodifluoromethane	1.166	85	1015	0.314	ug/l	97
9) 1,1,2-Trichlorotrifluo...	2.325	101	16421	5.034	ug/l	95
12) 1,1-Dichloroethene	2.319	96	2848	0.910	ug/l #	78
16) Acetone	2.386	43	4271	2.523	ug/l	93
19) Methyl tert-butyl Ether	3.117	73	11666	1.105	ug/l	99
22) Diisopropyl ether	3.764	45	5998	0.627	ug/l #	81
24) 1,1-Dichloroethane	3.617	63	3593	0.618	ug/l	95
27) cis-1,2-Dichloroethene	4.489	96	7934	2.049	ug/l	87
30) Chloroform	5.105	83	13002	2.026	ug/l	94
44) Trichloroethene	7.129	130	475705	112.731	ug/l	98
49) 1,4-Dioxane	7.671	88	1044	10.570	ug/l	99
64) Tetrachloroethene	9.275	164	12055	3.112	ug/l	97

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

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