

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021022\
 Data File : VX026943.D
 Acq On : 11 Feb 2022 01:12
 Operator : JC/MD
 Sample : N1481-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RW7B-GW-20220209

Quant Time: Feb 11 04:07:21 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020822W.M
 Quant Title : SW846 8260
 QLast Update : Tue Feb 08 07:07:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	123072	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	206220	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	185021	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	76456	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	79381	53.732	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 107.460%	
35) Dibromofluoromethane	5.391	113	64456	52.462	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 104.920%	
50) Toluene-d8	8.653	98	243296	53.947	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 107.900%	
62) 4-Bromofluorobenzene	11.079	95	86474	54.462	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 108.920%	
Target Compounds						
					Qvalue	
2) Dichlorodifluoromethane	1.173	85	355	0.265	ug/l	84
7) Trichlorofluoromethane	1.892	101	443	0.231	ug/l #	77
9) 1,1,2-Trichlorotrifluo...	2.337	101	37979	30.212	ug/l	99
12) 1,1-Dichloroethene	2.325	96	25400	20.166	ug/l	97
16) Acetone	2.386	43	1113	1.715	ug/l	96
24) 1,1-Dichloroethane	3.611	63	4958	2.040	ug/l	99
27) cis-1,2-Dichloroethene	4.495	96	3764	2.473	ug/l	88
30) Chloroform	5.099	83	2046	0.810	ug/l	85
32) 1,1,1-Trichloroethane	5.391	97	3584	1.610	ug/l	97
38) Carbon Tetrachloride	5.684	117	3808	1.874	ug/l	99
44) Trichloroethene	7.129	130	628666	370.648	ug/l	96
55) 1,1,2-Trichloroethane	9.153	97	1331	0.983	ug/l	89
64) Tetrachloroethene	9.275	164	645	0.347	ug/l	96

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

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