

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090822\
 Data File : VX031229.D
 Acq On : 08 Sep 2022 22:09
 Operator : JC/MD
 Sample : N4360-02RE
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TT17411-20220823RE

Quant Time: Sep 09 01:12:11 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090722W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 07 23:22:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	61598	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	108385	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	94445	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	40502	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	85784	58.822	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	117.640%
35) Dibromofluoromethane	5.391	113	58111	49.201	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.400%
50) Toluene-d8	8.653	98	195862	45.643	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	91.280%
62) 4-Bromofluorobenzene	11.079	95	77504	49.284	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	98.560%
Target Compounds						
						Qvalue
12) 1,1-Dichloroethene	2.318	96	1027	0.987	ug/l	99
16) Acetone	2.379	43	1026	2.032	ug/l	# 67
24) 1,1-Dichloroethane	3.617	63	5579	2.867	ug/l	98
30) Chloroform	5.104	83	547	0.238	ug/l	96
44) Trichloroethene	7.122	130	880	0.652	ug/l	86
64) Tetrachloroethene	9.274	164	498	0.507	ug/l	87

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

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