

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090122\
 Data File : VX031039.D
 Acq On : 01 Sep 2022 22:53
 Operator : JC/MD
 Sample : N4297-23
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RW9-MW01S-20220821

Quant Time: Sep 02 02:57:05 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090122W.M
 Quant Title : SW846 8260
 QLast Update : Fri Sep 02 02:49:21 2022
 Response via : Initial Calibration

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

Internal Standards							
1) Pentafluorobenzene		5.562	168	223212	50.000	ug/l	0.00
34) 1,4-Difluorobenzene		6.763	114	364963	50.000	ug/l	0.00
63) Chlorobenzene-d5		10.055	117	313483	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4		12.024	152	138980	50.000	ug/l	0.00
System Monitoring Compounds							
33) 1,2-Dichloroethane-d4		5.964	65	159710	55.608	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	111.220%	
35) Dibromofluoromethane		5.391	113	152021	57.078	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	114.160%	
50) Toluene-d8		8.653	98	562940	57.430	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	114.860%#	
62) 4-Bromofluorobenzene		11.085	95	185691	53.635	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	107.280%	
Target Compounds							Qvalue
29) Tetrahydrofuran		5.019	42	28591	25.909	ug/l	# 84
44) Trichloroethene		7.135	130	1808	0.533	ug/l	89

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

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