

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090122\
 Data File : VX031045.D
 Acq On : 02 Sep 2022 01:13
 Operator : JC/MD
 Sample : N4298-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE104D2-20220822

Quant Time: Sep 02 02:58:03 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	217714	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	346268	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	301964	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	131615	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	156494	55.865	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	111.720%
35) Dibromofluoromethane	5.391	113	149508	59.165	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	118.340%
50) Toluene-d8	8.653	98	550075	59.147	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	118.300%#
62) 4-Bromofluorobenzene	11.085	95	180898	55.072	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	110.140%
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.319	96	1162	0.459	ug/l	98
27) cis-1,2-Dichloroethene	4.489	96	48269	15.303	ug/l	84
30) Chloroform	5.105	83	4422	0.884	ug/l	95
42) 1,2-Dichloroethane	6.111	62	2669	0.796	ug/l	90
44) Trichloroethene	7.129	130	414273	128.797	ug/l	98

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

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