

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090222\
 Data File : VX031063.D
 Acq On : 02 Sep 2022 13:41
 Operator : JC/MD
 Sample : N4298-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EB01-20220822

Quant Time: Sep 03 00:49:25 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	5.556	168	209668	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	342551	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	303142	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	138974	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	131723	48.826	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	97.660%
35) Dibromofluoromethane	5.391	113	127065	50.830	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.660%
50) Toluene-d8	8.653	98	461382	50.149	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	100.300%
62) 4-Bromofluorobenzene	11.086	95	159953	49.224	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	98.440%
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	3859	3.705	ug/l #	83
30) Chloroform	5.099	83	28325	5.881	ug/l	94
44) Trichloroethene	7.129	130	974	0.306	ug/l	69
47) Bromodichloromethane	7.830	83	3110	0.863	ug/l #	84

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

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