

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090222\
 Data File : VX031064.D
 Acq On : 02 Sep 2022 14:05
 Operator : JC/MD
 Sample : N4298-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FB01-20220822

Quant Time: Sep 03 00:49:36 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	214559	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	341640	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	307761	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	142525	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	132087	47.845	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	95.700%
35) Dibromofluoromethane	5.391	113	128387	51.495	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	103.000%
50) Toluene-d8	8.653	98	472791	51.526	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	103.060%
62) 4-Bromofluorobenzene	11.085	95	159185	49.118	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	98.240%
Target Compounds						
					Qvalue	
3) Chloromethane	1.288	50	309	0.845	ug/l #	85
16) Acetone	2.386	43	9168	8.602	ug/l	94
30) Chloroform	5.099	83	32225	6.538	ug/l	91
47) Bromodichloromethane	7.836	83	3823	1.064	ug/l #	75

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

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