

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100522\
 Data File : VX031900.D
 Acq On : 05 Oct 2022 17:16
 Operator : JC/MD
 Sample : N4983-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB-RW5-TB-20220930

Quant Time: Oct 06 06:47:09 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100522W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 05 12:20:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	118479	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	168657	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	179417	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	120937	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	91117	51.060	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	102.120%
35) Dibromofluoromethane	5.385	113	86853	47.068	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.140%
50) Toluene-d8	8.653	98	323476	49.489	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	98.980%
62) 4-Bromofluorobenzene	11.079	95	110312	46.918	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	93.840%
Target Compounds						
						Qvalue
16) Acetone	2.386	43	1063	2.056	ug/l	89
30) Chloroform	5.099	83	2737	0.900	ug/l	85

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

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