

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102022\
 Data File : VX032269.D
 Acq On : 20 Oct 2022 17:37
 Operator : JC/MD
 Sample : N5185-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-01

Quant Time: Oct 21 02:14:48 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101922W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 19 16:51:09 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	150606	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	230902	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	199709	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	91635	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	95410	48.698	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	97.400%
35) Dibromofluoromethane	5.385	113	73286	45.611	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	91.220%
50) Toluene-d8	8.647	98	275458	47.625	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	95.260%
62) 4-Bromofluorobenzene	11.079	95	95698	45.184	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	90.360%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.971	59	32494	104.307	ug/l	96
19) Methyl tert-butyl Ether	3.111	73	169628	36.619	ug/l	92
22) Diisopropyl ether	3.764	45	4921	1.064	ug/l #	77
40) Benzene	6.038	78	11828	1.974	ug/l	93
67) Ethyl Benzene	10.195	91	10668	1.472	ug/l	96

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

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