

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110422\
 Data File : VX032600.D
 Acq On : 05 Nov 2022 03:13
 Operator : JC/MD
 Sample : N5483-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-2E

Quant Time: Nov 07 05:07:47 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 27 08:28:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	119207	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	196766	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	173113	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	79912	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	81063	48.941	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	97.880%
35) Dibromofluoromethane	5.391	113	63360	45.143	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	90.280%
50) Toluene-d8	8.647	98	236806	48.521	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.040%
62) 4-Bromofluorobenzene	11.079	95	90797	47.627	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	95.260%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	1007	1.673	ug/l	91
30) Chloroform	5.099	83	8859	3.405	ug/l	90
84) 1,2,4-Trimethylbenzene	11.756	105	1566	0.291	ug/l	86

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

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