

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082021\
 Data File : VX023854.D
 Acq On : 19 Aug 2021 21:43
 Operator : JC/MD
 Sample : M3483-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BB-TT-VPB-RW7-DUP-20210816

Quant Time: Aug 20 03:16:29 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 17 11:57:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	125870	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	215810	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	192367	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	66163	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	90219	51.356	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	102.720%
35) Dibromofluoromethane	5.397	113	71249	49.881	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.760%
50) Toluene-d8	8.653	98	264890	48.724	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.440%
62) 4-Bromofluorobenzene	11.085	95	80025	39.986	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	79.980%#
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	9037	11.505	ug/l #	86
25) 2-Butanone	4.587	43	2918	2.453	ug/l #	87
30) Chloroform	5.111	83	3118	1.193	ug/l #	78
44) Trichloroethene	7.135	130	2011	1.232	ug/l	96
64) Tetrachloroethene	9.275	164	612	0.398	ug/l #	77

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

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