

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010422\
 Data File : VX026281.D
 Acq On : 04 Jan 2022 18:56
 Operator : JC/MD
 Sample : M5333-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-7

Quant Time: Jan 05 06:13:32 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X123021W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 30 08:02:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	147523	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	255628	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	224751	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	95190	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	114216	53.516	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	107.040%
35) Dibromofluoromethane	5.391	113	83380	50.171	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	100.340%
50) Toluene-d8	8.653	98	300544	48.618	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.240%
62) 4-Bromofluorobenzene	11.079	95	110061	50.014	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	100.020%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.953	59	669334	1570.195	ug/l	99
19) Methyl tert-butyl Ether	3.111	73	1015166	177.915	ug/l	99
22) Diisopropyl ether	3.764	45	25199	4.090	ug/l #	78
42) 1,2-Dichloroethane	6.092	62	1953	0.631	ug/l	85

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

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