

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062321\
 Data File : VX022856.D
 Acq On : 23 Jun 2021 21:39
 Operator : JC/MD
 Sample : M2803-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW203D1-20210618

Manual Integrations
 APPROVED

MMDadoda
 6/24/2021 5:00:11 PM

Quant Time: Jun 24 03:38:21 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 23 19:12:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.568	168	53181	50.00	ug/l	0.01
34) 1,4-Difluorobenzene	6.769	114	108966	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	96605	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	35296	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	51744	53.61	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 107.22%			
35) Dibromofluoromethane	5.397	113	36129	52.21	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 104.42%			
50) Toluene-d8	8.653	98	143104	52.77	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 105.54%			
62) 4-Bromofluorobenzene	11.085	95	47886	47.31	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 94.62%			
Target Compounds					Qvalue	
12) 1,1-Dichloroethene	2.325	96	5907	8.56	ug/l	87
16) Acetone	2.392	43	1234	2.88	ug/l #	79
19) Methyl tert-butyl Ether	3.111	73	1563	0.63	ug/l	95
24) 1,1-Dichloroethane	3.617	63	30870	21.19	ug/l	99
27) cis-1,2-Dichloroethene	4.507	96	864m	1.02	ug/l	
30) Chloroform	5.111	83	1699	1.16	ug/l	98
32) 1,1,1-Trichloroethane	5.391	97	3671	3.12	ug/l #	88
44) Trichloroethene	7.135	130	5631	7.11	ug/l	93
64) Tetrachloroethene	9.281	164	1023	1.51	ug/l #	74

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

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