

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052423\
 Data File : VX035838.D
 Acq On : 24 May 2023 16:01
 Operator : JC/MD
 Sample : 02911-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TT17411-20230522

Quant Time: May 25 05:32:41 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051123W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 11 13:57:54 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	173318	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	306275	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	268414	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	125772	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	130399	43.127	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	86.260%	
35) Dibromofluoromethane	5.385	113	98506	47.691	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	95.380%	
50) Toluene-d8	8.647	98	335135	43.920	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	87.840%#	
62) 4-Bromofluorobenzene	11.079	95	132662	43.179	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	86.360%	
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.325	96	2234	1.157	ug/l	91
16) Acetone	2.404	43	2937	2.260	ug/l	99
19) Methyl tert-butyl Ether	3.111	73	1873	0.244	ug/l	93
24) 1,1-Dichloroethane	3.611	63	15960	3.686	ug/l	98
30) Chloroform	5.086	83	1358	0.306	ug/l	85
32) 1,1,1-Trichloroethane	5.379	97	1302	0.337	ug/l #	47
44) Trichloroethene	7.123	130	1897	0.820	ug/l	96
64) Tetrachloroethene	9.275	164	1293	0.716	ug/l	93

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

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