

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022724\
 Data File : VX040491.D
 Acq On : 27 Feb 2024 15:50
 Operator : JC/MD
 Sample : P1578-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TT161S1-20240221

Quant Time: Feb 28 02:04:32 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022624W.M
 Quant Title : SW846 8260
 QLast Update : Tue Feb 27 05:01:33 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	61793	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	110299	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	106425	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	53400	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	52043	50.355	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 100.700%		
35) Dibromofluoromethane	5.391	113	38450	50.130	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 100.260%		
50) Toluene-d8	8.647	98	138126	49.611	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 99.220%		
62) 4-Bromofluorobenzene	11.079	95	52938	49.880	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 99.760%		
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.325	96	404	0.574	ug/l #	81
16) Acetone	2.379	43	1901	5.640	ug/l #	86
19) Methyl tert-butyl Ether	3.117	73	1199	0.499	ug/l	92
24) 1,1-Dichloroethane	3.617	63	4440	3.076	ug/l	98
30) Chloroform	5.105	83	660	0.431	ug/l #	60
44) Trichloroethene	7.135	130	1444	1.678	ug/l	98
64) Tetrachloroethene	9.274	164	486	0.650	ug/l #	67

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

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