

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022724\
 Data File : VX040483.D
 Acq On : 27 Feb 2024 12:44
 Operator : JC/MD
 Sample : P1578-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TT149S1-20240222

Quant Time: Feb 28 02:01:41 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	5.550	168	67523	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	120274	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	117528	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	58081	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	55535	49.174	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.340%
35) Dibromofluoromethane	5.391	113	41731	49.895	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.780%
50) Toluene-d8	8.647	98	149142	49.125	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	98.260%
62) 4-Bromofluorobenzene	11.079	95	58394	50.458	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	100.920%
Target Compounds						
					Qvalue	
16) Acetone	2.373	43	1091	2.962	ug/l #	87
19) Methyl tert-butyl Ether	3.111	73	33381	12.721	ug/l	99
44) Trichloroethene	7.129	130	4661	4.968	ug/l	95
52) Toluene	8.726	92	2991	1.397	ug/l	96
64) Tetrachloroethene	9.275	164	1034	1.253	ug/l #	82
91) 1,2-Dichlorobenzene	12.341	146	623	0.324	ug/l	88

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

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