

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050224\
 Data File : VX041359.D
 Acq On : 02 May 2024 17:19
 Operator : JC/MD
 Sample : P2352-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-18B-112-050124

Quant Time: May 03 01:42:12 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X043024W.M
 Quant Title : SW846 8260
 QLast Update : Wed May 01 02:38:17 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	169244	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	215757	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	200845	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	112031	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	82891	44.989	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	89.980%
35) Dibromofluoromethane	5.385	113	74180	50.734	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.460%
50) Toluene-d8	8.647	98	233628	50.484	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.960%
62) 4-Bromofluorobenzene	11.079	95	92313	51.530	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	103.060%

Target Compounds						Qvalue
12) 1,1-Dichloroethene	2.313	96	25154	19.179	ug/l	87
16) Acetone	2.392	43	5794	3.281	ug/l	99
18) Methyl Acetate	2.709	43	4899	2.599	ug/l	97
24) 1,1-Dichloroethane	3.611	63	8581	3.253	ug/l	# 95
27) cis-1,2-Dichloroethene	4.495	96	12391	7.140	ug/l	81
30) Chloroform	5.105	83	3593	1.309	ug/l	90
40) Benzene	6.044	78	3427	0.636	ug/l	96
42) 1,2-Dichloroethane	6.092	62	6388	3.234	ug/l	100
44) Trichloroethene	7.129	130	15198	10.679	ug/l	89
52) Toluene	8.726	92	2062	0.625	ug/l	90

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

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