

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022824\
 Data File : VX040524.D
 Acq On : 28 Feb 2024 20:15
 Operator : JC/MD
 Sample : P1615-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TT17411-20240227

Quant Time: Feb 29 03:55:38 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	5.555	168	69984	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	120618	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	117781	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	56280	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	58009	49.558	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	99.120%
35) Dibromofluoromethane	5.385	113	41449	49.416	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.840%
50) Toluene-d8	8.646	98	149558	49.122	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	98.240%
62) 4-Bromofluorobenzene	11.079	95	58493	50.400	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	100.800%
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.324	96	957	1.200	ug/l #	84
16) Acetone	2.373	43	822	2.153	ug/l #	81
19) Methyl tert-butyl Ether	3.123	73	869	0.320	ug/l #	87
24) 1,1-Dichloroethane	3.611	63	5930	3.628	ug/l #	91
30) Chloroform	5.098	83	565	0.326	ug/l	84
44) Trichloroethene	7.122	130	2506	2.663	ug/l	94
52) Toluene	8.720	92	2437	1.135	ug/l	92
64) Tetrachloroethene	9.280	164	578	0.699	ug/l #	81

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

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