

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080823\
 Data File : VX036868.D
 Acq On : 08 Aug 2023 16:51
 Operator : JC/MD
 Sample : 03903-18
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-08W

Quant Time: Aug 09 04:16:58 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 21 12:14:12 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	122013	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	223114	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	197795	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	84460	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	85303	47.250	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	94.500%
35) Dibromofluoromethane	5.391	113	71936	46.908	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.820%
50) Toluene-d8	8.646	98	262857	48.334	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	96.660%
62) 4-Bromofluorobenzene	11.079	95	94524	46.069	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	92.140%
Target Compounds						
					Qvalue	
16) Acetone	2.385	43	5004	6.680	ug/l	92
30) Chloroform	5.092	83	45723	15.605	ug/l	98
44) Trichloroethene	7.128	130	1694	0.975	ug/l	76
47) Bromodichloromethane	7.823	83	1469	0.576	ug/l	87
64) Tetrachloroethene	9.274	164	10549	7.150	ug/l	91
68) m/p-Xylenes	10.305	106	1186	0.399	ug/l	97
84) 1,2,4-Trimethylbenzene	11.756	105	1856	0.317	ug/l	95

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

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