

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012224\
 Data File : VX039885.D
 Acq On : 22 Jan 2024 15:45
 Operator : JC/MD
 Sample : P1186-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB-191-GW-278-280

Quant Time: Jan 23 00:18:18 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011024W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jan 11 03:51:26 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	78234	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	144514	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.049	117	139815	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	71768	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	62109	47.392	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	94.780%
35) Dibromofluoromethane	5.385	113	48935	50.274	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	100.540%
50) Toluene-d8	8.647	98	181004	49.305	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	98.620%
62) 4-Bromofluorobenzene	11.079	95	71252	49.633	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	99.260%
Target Compounds						
					Qvalue	
16) Acetone	2.374	43	1493	2.026	ug/l #	89
17) Carbon Disulfide	2.514	76	718	0.265	ug/l #	80
19) Methyl tert-butyl Ether	3.111	73	12705	3.801	ug/l	98
22) Diisopropyl ether	3.770	45	1830	0.517	ug/l #	56
93) 1,2,4-Trichlorobenzene	13.591	180	364	0.247	ug/l #	66

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

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