

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110823\
 Data File : VX038649.D
 Acq On : 08 Nov 2023 12:35
 Operator : JC/MD
 Sample : 05196-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TB-10-31-23

Quant Time: Nov 09 01:19:17 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X110723W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 08 03:19:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	280546	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	462125	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	471740	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	268406	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	160704	46.011	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	92.020%	
35) Dibromofluoromethane	5.385	113	150892	45.335	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	90.680%	
50) Toluene-d8	8.647	98	543138	45.603	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	91.200%#	
62) 4-Bromofluorobenzene	11.079	95	235735	51.050	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	102.100%	
Target Compounds						
						Qvalue
3) Chloromethane	1.301	50	2639	1.048	ug/l	95
16) Acetone	2.380	43	12513	8.507	ug/l	94
17) Carbon Disulfide	2.514	76	1473	0.212	ug/l	97
20) Methylene Chloride	2.788	84	1770	0.597	ug/l	82
25) 2-Butanone	4.580	43	3576	1.678	ug/l	97
51) 4-Methyl-2-Pentanone	8.580	43	3824	0.904	ug/l	96

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

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