

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060222\
 Data File : VX029178.D
 Acq On : 03 Jun 2022 09:01
 Operator : JC/MD
 Sample : N3164-18 50X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 40S-20220531

Quant Time: Jun 03 09:43:59 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053122W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 01 04:45:42 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	305585	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	574860	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	553680	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	249732	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	201934	58.601	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	117.200%
35) Dibromofluoromethane	5.391	113	190430	49.872	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	99.740%
50) Toluene-d8	8.653	98	679533	48.659	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	97.320%
62) 4-Bromofluorobenzene	11.079	95	266128	49.298	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	98.600%
Target Compounds						
						Qvalue
3) Chloromethane	1.294	50	675	0.243	ug/l	# 86
16) Acetone	2.379	43	2362	1.741	ug/l	# 81
20) Methylene Chloride	2.794	84	817	0.258	ug/l	97
27) cis-1,2-Dichloroethene	4.507	96	1457	0.393	ug/l	79
37) Ethyl Acetate	4.721	43	9297	1.953	ug/l	# 90
44) Trichloroethene	7.135	130	6491	0.814	ug/l	92

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

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