

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100720\
 Data File : VN063978.D
 Acq On : 8 Oct 2020 6:53
 Operator : JC/MD
 Sample : L4276-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW4D

Quant Time: Oct 08 07:15:37 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 01 01:46:38 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	239226	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	431003	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	440051	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	181877	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	201283	50.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.86%	
35) Dibromofluoromethane	7.55	113	147937	49.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.94%	
50) Toluene-d8	10.06	98	551241	48.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.24%	
62) 4-Bromofluorobenzene	12.38	95	208629	48.95	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.90%	
Target Compounds						
16) Acetone	3.78	43	29938	24.608	ug/l	92
25) 2-Butanone	6.80	43	10729	6.220	ug/l	93
27) cis-1,2-Dichloroethene	6.78	96	10227	3.296	ug/l	97
40) Benzene	8.00	78	7854	0.630	ug/l #	86
44) Trichloroethene	8.80	130	3604	1.103	ug/l	83
64) Tetrachloroethene	10.60	164	46286	13.595	ug/l	95

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

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