

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN082520\
 Data File : VN063118.D
 Acq On : 25 Aug 2020 17:51
 Operator : JC/MD
 Sample : L3788-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 1794

Quant Time: Aug 26 06:51:00 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N081820W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 19 06:09:27 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	160555	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	273314	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	266029	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	119939	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	98430	51.73	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.46%	
35) Dibromofluoromethane	7.55	113	80589	45.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.50%	
50) Toluene-d8	10.06	98	345375	50.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.98%	
62) 4-Bromofluorobenzene	12.38	95	126990	52.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.22%	
Target Compounds						
3) Chloromethane	2.04	50	3308	1.503	ug/l	88
11) Tert butyl alcohol	4.73	59	822562	3051.793	ug/l	99
16) Acetone	3.78	43	3812211	6522.892	ug/l	100
18) Methyl Acetate	4.29	43	77422	40.942	ug/l	99
25) 2-Butanone	6.79	43	1195047	1182.932	ug/l	97
51) 4-Methyl-2-Pentanone	9.96	43	40978	17.339	ug/l	94
59) 2-Hexanone	10.72	43	120438	73.060	ug/l	97

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

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