

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN081920\
 Data File : VN063019.D
 Acq On : 19 Aug 2020 17:57
 Operator : JC/MD
 Sample : L3496-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 AU-06-081720

Quant Time: Aug 20 01:48:07 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	160541	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	270604	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	298286	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	106724	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	104706	55.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.06%	
35) Dibromofluoromethane	7.55	113	91510	51.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.80%	
50) Toluene-d8	10.06	98	362201	53.48	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.96%	
62) 4-Bromofluorobenzene	12.38	95	122905	51.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.86%	
Target Compounds						
16) Acetone	3.79	43	7882	9.511	ug/l	96
18) Methyl Acetate	4.29	43	2779	1.470	ug/l #	90
20) Methylene Chloride	4.50	84	7251	3.069	ug/l	99
43) Isopropyl Acetate	8.13	43	35699	9.643	ug/l #	90

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

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