

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111320\
 Data File : VN064645.D
 Acq On : 13 Nov 2020 21:11
 Operator : JC/MD
 Sample : L4719-12
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-3A-S

Quant Time: Nov 17 03:26:55 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 04 12:09:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	213367	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	461484	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	496887	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	171916	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	218327	58.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	117.48%	
35) Dibromofluoromethane	7.55	113	160077	50.95	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.90%	
50) Toluene-d8	10.06	98	613150	50.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.98%	
62) 4-Bromofluorobenzene	12.38	95	225500	49.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.20%	
Target Compounds						
16) Acetone	3.79	43	13503	9.538	ug/l	98
18) Methyl Acetate	4.29	43	3341	1.200	ug/l #	77
20) Methylene Chloride	4.50	84	11268	4.012	ug/l #	82
43) Isopropyl Acetate	8.13	43	87252	11.043	ug/l	99

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

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