

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111220\
 Data File : VN064586.D
 Acq On : 12 Nov 2020 17:13
 Operator : JC/MD
 Sample : L4701-15
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TRANSFORMER-5-1-TOP

Quant Time: Nov 13 03:45:27 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	224166	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	494888	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	527287	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	185250	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	235753	60.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	120.76%	
35) Dibromofluoromethane	7.55	113	170947	50.73	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.46%	
50) Toluene-d8	10.06	98	644588	49.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.00%	
62) 4-Bromofluorobenzene	12.38	95	241972	49.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.26%	
Target Compounds						
16) Acetone	3.78	43	17488	11.758	ug/l	98
18) Methyl Acetate	4.28	43	7643	2.614	ug/l	93
20) Methylene Chloride	4.51	84	10175	3.448	ug/l	86
43) Isopropyl Acetate	8.13	43	77449	9.141	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111220\
Data File : VN064586.D
Acq On : 12 Nov 2020 17:13
Operator : JC/MD
Sample : L4701-15
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TRANSFORMER-5-1-TOP

Quant Time: Nov 13 03:45:27 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.M
Quant Title : SW846 8260
QLast Update : Wed Nov 04 12:09:03 2020
Response via : Initial Calibration

