

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN110520\
 Data File : VN064470.D
 Acq On : 5 Nov 2020 18:55
 Operator : JC/MD
 Sample : L4625-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CONTAINER-001-1B

Quant Time: Nov 06 03:43:29 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	241244	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	504934	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	540224	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	189844	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	238225	56.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.38%	
35) Dibromofluoromethane	7.55	113	171460	49.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.74%	
50) Toluene-d8	10.06	98	663824	49.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.92%	
62) 4-Bromofluorobenzene	12.38	95	251706	50.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.18%	
Target Compounds						
16) Acetone	3.78	43	20013	12.503	ug/l	99
18) Methyl Acetate	4.29	43	5328	1.693	ug/l #	87
20) Methylene Chloride	4.50	84	9426	2.968	ug/l	86
43) Isopropyl Acetate	8.13	43	106079	12.271	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN110520\
Data File : VN064470.D
Acq On : 5 Nov 2020 18:55
Operator : JC/MD
Sample : L4625-08
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
CONTAINER-001-1B

Quant Time: Nov 06 03:43:29 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

