

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN110520\
 Data File : VN064471.D
 Acq On : 5 Nov 2020 19:19
 Operator : JC/MD
 Sample : L4625-09
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CONTAINER-001-2B

Quant Time: Nov 06 03:45:38 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	245745	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	517543	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	553319	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	205548	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	244010	57.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	114.02%	
35) Dibromofluoromethane	7.55	113	174186	49.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.86%	
50) Toluene-d8	10.06	98	671105	49.28	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.56%	
62) 4-Bromofluorobenzene	12.38	95	265737	51.59	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.18%	
Target Compounds						
16) Acetone	3.78	43	21911	13.438	ug/l	90
18) Methyl Acetate	4.30	43	5262	1.642	ug/l #	85
20) Methylene Chloride	4.50	84	9225	2.852	ug/l #	88
43) Isopropyl Acetate	8.13	43	103229	11.650	ug/l	100

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN110520\
Data File : VN064471.D
Acq On : 5 Nov 2020 19:19
Operator : JC/MD
Sample : L4625-09
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
CONTAINER-001-2B

Quant Time: Nov 06 03:45:38 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

