

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100820\
 Data File : VN064022.D
 Acq On : 9 Oct 2020 5:40
 Operator : JC/MD
 Sample : L4291-10
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAV-B4-100620-DP

Quant Time: Oct 09 12:12:57 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 01 01:46:38 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	238768	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	461948	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	478567	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	182808	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	204190	51.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.52%	
35) Dibromofluoromethane	7.55	113	153440	47.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.74%	
50) Toluene-d8	10.06	98	594877	48.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.92%	
62) 4-Bromofluorobenzene	12.38	95	221221	48.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.86%	
Target Compounds						
16) Acetone	3.79	43	15186	12.506	ug/l	99
20) Methylene Chloride	4.51	84	10371	3.820	ug/l	91
43) Isopropyl Acetate	8.13	43	90371	11.783	ug/l	99
51) 4-Methyl-2-Pentanone	9.95	43	9432	2.096	ug/l	91

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN100820\
Data File : VN064022.D
Acq On : 9 Oct 2020 5:40
Operator : JC/MD
Sample : L4291-10
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PAV-B4-100620-DP

Quant Time: Oct 09 12:12:57 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260
QLast Update : Thu Oct 01 01:46:38 2020
Response via : Initial Calibration

