

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN072920\
 Data File : VN062653.D
 Acq On : 29 Jul 2020 14:39
 Operator : JC/MD
 Sample : L3464-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 32537

Quant Time: Jul 30 04:07:32 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072120W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jul 21 18:48:59 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	138111	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	248860	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	254934	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	104089	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	118477	55.22	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.44%	
35) Dibromofluoromethane	7.55	113	82883	50.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.40%	
50) Toluene-d8	10.06	98	332629	51.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.00%	
62) 4-Bromofluorobenzene	12.38	95	125238	52.52	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.04%	
Target Compounds						
16) Acetone	3.78	43	32358	40.079	ug/l	99
18) Methyl Acetate	4.29	43	7560	3.699	ug/l	96
20) Methylene Chloride	4.50	84	13413	6.927	ug/l	94
25) 2-Butanone	6.79	43	8180	6.653	ug/l	95
43) Isopropyl Acetate	8.13	43	52450	11.379	ug/l #	91

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_N\DATA\VN072920\
Data File : VN062653.D
Acq On : 29 Jul 2020 14:39
Operator : JC/MD
Sample : L3464-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
32537

Quant Time: Jul 30 04:07:32 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N072120W.M
Quant Title : SW846 8260
QLast Update : Tue Jul 21 18:48:59 2020
Response via : Initial Calibration

