

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN041420\
 Data File : VN060945.D
 Acq On : 14 Apr 2020 19:38
 Operator : JC/MD
 Sample : L2230-24
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-12-B

Quant Time: Apr 15 09:41:29 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031820W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 18 08:49:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	149643	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.56	114	254932	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	242048	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	111676	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	100836	49.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.18%	
35) Dibromofluoromethane	7.56	113	75791	49.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.34%	
50) Toluene-d8	10.08	98	319132	50.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.36%	
62) 4-Bromofluorobenzene	12.39	95	121027	52.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.90%	
Target Compounds						
16) Acetone	3.78	43	21890	30.083	ug/l	99
20) Methylene Chloride	4.52	84	2607	1.456	ug/l	91
25) 2-Butanone	6.81	43	6606	5.610	ug/l	94
43) Isopropyl Acetate	8.14	43	42576	9.679	ug/l #	90

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN041420\
Data File : VN060945.D
Acq On : 14 Apr 2020 19:38
Operator : JC/MD
Sample : L2230-24
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-12-B

Quant Time: Apr 15 09:41:29 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N031820W.M
Quant Title : SW846 8260
QLast Update : Wed Mar 18 08:49:09 2020
Response via : Initial Calibration

