

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN040720\
 Data File : VN060887.D
 Acq On : 7 Apr 2020 19:05
 Operator : JC/MD
 Sample : L2171-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FRAC-COMP

Quant Time: Apr 08 08:39:18 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	197335	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	329600	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	304722	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	137104	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	127249	46.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.94%	
35) Dibromofluoromethane	7.56	113	97320	48.83	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.66%	
50) Toluene-d8	10.08	98	400844	49.24	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.48%	
62) 4-Bromofluorobenzene	12.40	95	148194	49.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.34%	
Target Compounds						
16) Acetone	3.79	43	4059	4.230	ug/l	98
25) 2-Butanone	6.81	43	7495	4.827	ug/l	94
30) Chloroform	7.35	83	80523	19.408	ug/l	99
47) Bromodichloromethane	9.38	83	13068	4.092	ug/l	99
60) Dibromochloromethane	10.89	129	2337	1.011	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN040720\
Data File : VN060887.D
Acq On : 7 Apr 2020 19:05
Operator : JC/MD
Sample : L2171-01
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FRAC-COMP

Quant Time: Apr 08 08:39:18 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

