

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN102919\
 Data File : VN059007.D
 Acq On : 29 Oct 2019 13:55
 Operator : JC/SP
 Sample : K5576-11
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CSA-2-1-8

Quant Time: Oct 30 04:06:57 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N101819W.M
 Quant Title : SW846 8260
 QLast Update : Sat Oct 19 00:16:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	386584	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	605730	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	530870	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	183674	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	263748	46.52	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.04%	
35) Dibromofluoromethane	7.57	113	195103	48.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.98%	
50) Toluene-d8	10.08	98	759933	49.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.32%	
62) 4-Bromofluorobenzene	12.40	95	232849	41.02	ug/l	0.00
Spiked Amount 50.000			Recovery	=	82.04%	
Target Compounds						
16) Acetone	3.80	43	19182	6.847	ug/l	98
18) Methyl Acetate	4.30	43	13142	2.063	ug/l #	69
20) Methylene Chloride	4.53	84	8547	1.787	ug/l	92
43) Isopropyl Acetate	8.15	43	80913	7.466	ug/l #	92

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN102919\
Data File : VN059007.D
Acq On : 29 Oct 2019 13:55
Operator : JC/SP
Sample : K5576-11
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
CSA-2-1-8

Quant Time: Oct 30 04:06:57 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N101819W.M
Quant Title : SW846 8260
QLast Update : Sat Oct 19 00:16:20 2019
Response via : Initial Calibration

