

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092519\
 Data File : VN058365.D
 Acq On : 25 Sep 2019 14:27
 Operator : JC/SP
 Sample : K4657-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BU-02-092419

Quant Time: Sep 26 03:45:45 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091819W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 19 09:27:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	488366	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	817206	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	717323	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	289235	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	367685	54.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.12%	
35) Dibromofluoromethane	7.58	113	255883	51.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.02%	
50) Toluene-d8	10.09	98	1042918	53.88	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.76%	
62) 4-Bromofluorobenzene	12.41	95	340745	47.39	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.78%	
Target Compounds						
16) Acetone	3.81	43	17756	7.690	ug/l	93
20) Methylene Chloride	4.53	84	3833232	952.271	ug/l	96
43) Isopropyl Acetate	8.15	43	121178	10.609	ug/l #	87
95) Naphthalene	15.14	128	109654	7.937	ug/l	98

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

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092519\
Data File : VN058365.D
Acq On : 25 Sep 2019 14:27
Operator : JC/SP
Sample : K4657-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
BU-02-092419

Quant Time: Sep 26 03:45:45 2019
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091819W.M
Quant Title : SW846 8260
QLast Update : Thu Sep 19 09:27:53 2019
Response via : Initial Calibration

