

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN103119\
 Data File : VN059033.D
 Acq On : 31 Oct 2019 12:48
 Operator : JC/SP
 Sample : K5147-11
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OK-01-103019

Quant Time: Nov 01 01:31:51 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	404265	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	630977	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	579971	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	212470	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	276004	46.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.10%	
35) Dibromofluoromethane	7.57	113	203973	49.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.34%	
50) Toluene-d8	10.08	98	799226	49.63	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.26%	
62) 4-Bromofluorobenzene	12.41	95	254702	43.07	ug/l	0.00
Spiked Amount 50.000			Recovery	=	86.14%	
Target Compounds						
16) Acetone	3.80	43	41895	14.299	ug/l	98
18) Methyl Acetate	4.31	43	7008	1.052	ug/l	95
20) Methylene Chloride	4.53	84	84337	16.858	ug/l	96
43) Isopropyl Acetate	8.15	43	100333	8.888	ug/l #	92
95) Naphthalene	15.13	128	28211	4.440	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_N\DATA\VN103119\
Data File : VN059033.D
Acq On : 31 Oct 2019 12:48
Operator : JC/SP
Sample : K5147-11
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
OK-01-103019

Quant Time: Nov 01 01:31:51 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

