

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091518\
 Data File : VN051262.D
 Acq On : 16 Sep 2018 8:16
 Operator : MD\SY
 Sample : J4936-38
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 132-1-VOLT

Quant Time: Sep 17 05:59:30 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091018W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 11 02:27:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	397199	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	613787	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	496274	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	141160	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	8.03	65	220092	49.79	ug/l	0.00
Spiked Amount			Recovery	=	99.58%	
35) Dibromofluoromethane	7.59	113	233894	50.71	ug/l	0.00
Spiked Amount			Recovery	=	101.42%	
50) Toluene-d8	10.09	98	798397	45.62	ug/l	0.00
Spiked Amount			Recovery	=	91.24%	
62) 4-Bromofluorobenzene	12.40	95	197638	34.14	ug/l	0.00
Spiked Amount			Recovery	=	68.28%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	3.82	43	15110	8.18	ug/l	99
18) Methyl Acetate	4.34	43	11209	3.17	ug/l	96
20) Methylene Chloride	4.55	84	1846529	378.18	ug/l	99
30) Chloroform	7.38	83	10857	1.34	ug/l	95
43) Isopropyl Acetate	8.17	43	134495	19.53	ug/l #	84

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091518\
Data File : VN051262.D
Acq On : 16 Sep 2018 8:16
Operator : MD\SY
Sample : J4936-38
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
132-1-VOLT

Quant Time: Sep 17 05:59:30 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091018W.M
Quant Title : SW846 8260
QLast Update : Tue Sep 11 02:27:09 2018
Response via : Initial Calibration

