

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100219\
 Data File : VN058487.D
 Acq On : 2 Oct 2019 19:58
 Operator : JC/SP
 Sample : K5095-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 COMPOSITE-2

Quant Time: Oct 03 05:56:09 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
 Quant Title : SW846 8260
 QLast Update : Mon Sep 30 08:07:38 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	354789	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	601759	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	515147	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	190900	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	270506	52.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.34%	
35) Dibromofluoromethane	7.58	113	186808	50.65	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.30%	
50) Toluene-d8	10.08	98	756938	52.39	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.78%	
62) 4-Bromofluorobenzene	12.41	95	234926	44.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.94%	
Target Compounds						
16) Acetone	3.80	43	14468	8.186	ug/l	94
20) Methylene Chloride	4.53	84	954872	234.112	ug/l	98
43) Isopropyl Acetate	8.16	43	208690	22.749	ug/l #	81
84) 1,2,4-Trimethylbenzene	13.05	105	16210	1.370	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN100219\
Data File : VN058487.D
Acq On : 2 Oct 2019 19:58
Operator : JC/SP
Sample : K5095-02
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
COMPOSITE-2

Quant Time: Oct 03 05:56:09 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
Quant Title : SW846 8260
QLast Update : Mon Sep 30 08:07:38 2019
Response via : Initial Calibration

