

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061819\
 Data File : VN056366.D
 Acq On : 18 Jun 2019 15:19
 Operator : JC/SP
 Sample : K3433-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OILY-WATER

Quant Time: Jun 19 03:35:48 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N061719W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 18 04:23:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.66	168	339326	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	565310	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	569759	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	209710	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	192321	50.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.82%	
35) Dibromofluoromethane	7.59	113	173536	47.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.74%	
50) Toluene-d8	10.09	98	698670	50.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.90%	
62) 4-Bromofluorobenzene	12.40	95	246775	53.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.54%	
Target Compounds						
16) Acetone	3.82	43	58591	27.791	ug/l	98
18) Methyl Acetate	4.33	43	8872	0.654	ug/l	93
20) Methylene Chloride	4.54	84	6465	1.737	ug/l #	95
25) 2-Butanone	6.85	43	17585	6.882	ug/l	95
30) Chloroform	7.37	83	1001339	165.157	ug/l	99
52) Toluene	10.16	92	43592	4.628	ug/l	95

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN061819\
Data File : VN056366.D
Acq On : 18 Jun 2019 15:19
Operator : JC/SP
Sample : K3433-01
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
OILY-WATER

Quant Time: Jun 19 03:35:48 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N061719W.M
Quant Title : SW846 8260
QLast Update : Tue Jun 18 04:23:59 2019
Response via : Initial Calibration

