

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092018\
 Data File : VN051348.D
 Acq On : 20 Sep 2018 18:55
 Operator : MD\SY
 Sample : J4982-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OILY-WATER

Quant Time: Sep 21 07:08:35 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092018W.M
 Quant Title : SW846 8260
 QLast Update : Fri Sep 21 03:27:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	822077	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1178773	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1007047	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	393045	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	441972	49.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.20%	
35) Dibromofluoromethane	7.59	113	444186	46.63	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.26%	
50) Toluene-d8	10.09	98	1616060	44.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.26%	
62) 4-Bromofluorobenzene	12.40	95	451084	37.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	75.72%	
Target Compounds						
					Qvalue	
16) Acetone	3.82	43	3938	2.151	ug/l #	83
17) Carbon Disulfide	4.06	76	9844	0.425	ug/l	96
18) Methyl Acetate	4.34	43	2681	0.426	ug/l #	77
20) Methylene Chloride	4.55	84	2956	0.323	ug/l	91
42) 1,2-Dichloroethane	8.12	62	5896	0.537	ug/l	86

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092018\
Data File : VN051348.D
Acq On : 20 Sep 2018 18:55
Operator : MD\SY
Sample : J4982-01
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
OILY-WATER

Quant Time: Sep 21 07:08:35 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N092018W.M
Quant Title : SW846 8260
QLast Update : Fri Sep 21 03:27:45 2018
Response via : Initial Calibration

