

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061818\
 Data File : VN049379.D
 Acq On : 19 Jun 2018 7:32
 Operator : MD\SY
 Sample : J3431-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SW-1RR

Quant Time: Jun 19 08:53:45 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N061618W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 16 01:02:01 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	1407121	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	2078153	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1738746	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	541494	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	846592	45.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.72%	
35) Dibromofluoromethane	7.59	113	768046	45.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.36%	
50) Toluene-d8	10.09	98	2906001	45.39	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.78%	
62) 4-Bromofluorobenzene	12.41	95	770829	36.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	72.82%	
Target Compounds						
16) Acetone	3.82	43	132089	29.149	ug/l	96
18) Methyl Acetate	4.33	43	48185	3.487	ug/l	98
20) Methylene Chloride	4.55	84	49874	2.635	ug/l	89
25) 2-Butanone	6.85	43	54059	8.884	ug/l	96
31) Cyclohexane	7.66	56	293690	10.501	ug/l #	83
39) Methylcyclohexane	9.08	83	155293	6.074	ug/l	95

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN061818\
Data File : VN049379.D
Acq On : 19 Jun 2018 7:32
Operator : MD\SY
Sample : J3431-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SW-1RR

Quant Time: Jun 19 08:53:45 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N061618W.M
Quant Title : SW846 8260
QLast Update : Sat Jun 16 01:02:01 2018
Response via : Initial Calibration

