

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN082318\
 Data File : VN050844.D
 Acq On : 23 Aug 2018 11:29
 Operator : MD\SY
 Sample : J4612-07
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TOTE

Quant Time: Aug 24 11:56:44 2018
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 22 05:38:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	635624	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	992694	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	818839	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	276021	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	403623	52.75	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.50%	
35) Dibromofluoromethane	7.59	113	383298	48.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.06%	
50) Toluene-d8	10.09	98	1366082	46.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.54%	
62) 4-Bromofluorobenzene	12.40	95	356791	36.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	72.72%	
Target Compounds						
16) Acetone	3.82	43	126797	64.74	ug/l	Qvalue 100
17) Carbon Disulfide	4.05	76	47049	1.97	ug/l	98
25) 2-Butanone	6.85	43	13456	4.36	ug/l	90
29) Tetrahydrofuran	7.22	42	7050	3.99	ug/l #	84

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

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN082318\
Data File : VN050844.D
Acq On : 23 Aug 2018 11:29
Operator : MD\SY
Sample : J4612-07
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TOTE

Quant Time: Aug 24 11:56:44 2018
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.M
Quant Title : SW846 8260
QLast Update : Wed Aug 22 05:38:12 2018
Response via : Initial Calibration

