

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN081519\
 Data File : VN057309.D
 Acq On : 15 Aug 2019 00:45
 Operator : JC/SP
 Sample : K4329-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ETGI-292

Quant Time: Aug 15 08:58:14 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N081519W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 15 08:22:24 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.66	168	1125699	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	1590494	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1330288	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	344033	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.02	65	536558	43.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.70%	
35) Dibromofluoromethane	7.58	113	477300	45.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.56%	
50) Toluene-d8	10.08	98	1844515	48.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.74%	
62) 4-Bromofluorobenzene	12.40	95	489463	36.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	72.64%	
Target Compounds						
16) Acetone	3.81	43	24537	5.780	ug/l	98
18) Methyl Acetate	4.32	43	31496	0.375	ug/l	96
20) Methylene Chloride	4.54	84	145140	12.219	ug/l	93
43) Isopropyl Acetate	8.16	43	208769	8.444	ug/l #	75

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN081519\
Data File : VN057309.D
Acq On : 15 Aug 2019 00:45
Operator : JC/SP
Sample : K4329-02
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
ETGI-292

Quant Time: Aug 15 08:58:14 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N081519W.M
Quant Title : SW846 8260
QLast Update : Thu Aug 15 08:22:24 2019
Response via : Initial Calibration

