

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN081319\
 Data File : VN057272.D
 Acq On : 14 Aug 2019 1:53
 Operator : JC/SP
 Sample : K4266-22
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 R1-R2-A4-(4.3)

Quant Time: Aug 14 07:54:35 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N081319W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 14 06:25:03 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.66	168	1485462	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	2056656	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1752577	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	458983	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.02	65	693144	46.26	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.52%	
35) Dibromofluoromethane	7.58	113	625099	47.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.20%	
50) Toluene-d8	10.09	98	2410423	50.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.64%	
62) 4-Bromofluorobenzene	12.40	95	652741	39.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	79.48%	
Target Compounds						
16) Acetone	3.81	43	29024	5.649	ug/l	96
18) Methyl Acetate	4.32	43	50946	2.747	ug/l #	88
20) Methylene Chloride	4.55	84	46966	3.130	ug/l	98
43) Isopropyl Acetate	8.16	43	324254	13.623	ug/l #	76

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN081319\
Data File : VN057272.D
Acq On : 14 Aug 2019 1:53
Operator : JC/SP
Sample : K4266-22
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 38 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
R1-R2-A4-(4.3)

Quant Time: Aug 14 07:54:35 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N081319W.M
Quant Title : SW846 8260
QLast Update : Wed Aug 14 06:25:03 2019
Response via : Initial Calibration

