

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN081919\
 Data File : VN057393.D
 Acq On : 19 Aug 2019 12:18
 Operator : JC/SP
 Sample : K4339-14
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 T8-A1-(0)

Manual Integrations
 APPROVED

MMDadoda
 8/21/2019 9:45:16 AM

Quant Time: Aug 20 07:57:25 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N081519W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 15 11:33:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.66	168	1080187	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	1511684	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1272339	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	329717	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.02	65	493779	42.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	84.10%	
35) Dibromofluoromethane	7.58	113	454737	45.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.78%	
50) Toluene-d8	10.09	98	1748763	48.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.50%	
62) 4-Bromofluorobenzene	12.40	95	462587	36.12	ug/l	0.00
Spiked Amount 50.000			Recovery	=	72.24%	
Target Compounds						
16) Acetone	3.82	43	16862m	4.139	ug/l	Qvalue
20) Methylene Chloride	4.55	84	22664	1.989	ug/l #	84
43) Isopropyl Acetate	8.16	43	207458	9.462	ug/l #	85

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN081919\
Data File : VN057393.D
Acq On : 19 Aug 2019 12:18
Operator : JC/SP
Sample : K4339-14
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
T8-A1-(0)

Manual Integrations
APPROVED

MMDadoda
8/21/2019 9:45:16 AM

Quant Time: Aug 20 07:57:25 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N081519W.M
Quant Title : SW846 8260
QLast Update : Thu Aug 15 11:33:14 2019
Response via : Initial Calibration

