

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN070819\
 Data File : VN056617.D
 Acq On : 8 Jul 2019 18:14
 Operator : JC/SP
 Sample : K3622-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OR-02-070319-B

Quant Time: Jul 09 04:46:26 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N061919W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 21 03:45:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	292548	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	515212	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	543214	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	219109	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	172488	53.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.40%	
35) Dibromofluoromethane	7.59	113	152406	47.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.84%	
50) Toluene-d8	10.09	98	649008	51.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.78%	
62) 4-Bromofluorobenzene	12.40	95	242770	58.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	116.46%	
Target Compounds						
16) Acetone	3.82	43	16270	10.016	ug/l	99
18) Methyl Acetate	4.34	43	7076	1.380	ug/l #	89
20) Methylene Chloride	4.55	84	52782	16.642	ug/l	98
43) Isopropyl Acetate	8.17	43	19597	2.852	ug/l #	88
95) Naphthalene	15.13	128	8194	5.976	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN070819\
Data File : VN056617.D
Acq On : 8 Jul 2019 18:14
Operator : JC/SP
Sample : K3622-04
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
OR-02-070319-B

Quant Time: Jul 09 04:46:26 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N061919W.M
Quant Title : SW846 8260
QLast Update : Fri Jun 21 03:45:40 2019
Response via : Initial Calibration

