

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060521\
 Data File : VN067239.D
 Acq On : 5 Jun 2021 18:46
 Operator : JC/MD
 Sample : M2603-04
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FAM4-WC1

Quant Time: Jun 07 01:18:00 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060421W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 04 19:00:20 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.091	168	210700	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	376540	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.749	117	377383	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	156693	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.442	65	141554	49.286	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	98.580%		
35) Dibromofluoromethane	8.027	113	110645	47.400	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	94.800%		
50) Toluene-d8	10.446	98	504364	53.052	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	106.100%		
62) 4-Bromofluorobenzene	12.736	95	161570	44.127	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	88.260%		
Target Compounds					Qvalue	
16) Acetone	4.299	43	7202	7.267	ug/l	90
20) Methylene Chloride	5.111	84	5521	2.212	ug/l	93
43) Isopropyl Acetate	8.566	43	15541	2.438	ug/l #	82

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060521\
Data File : VN067239.D
Acq On : 5 Jun 2021 18:46
Operator : JC/MD
Sample : M2603-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FAM4-WC1

Quant Time: Jun 07 01:18:00 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060421W.M
Quant Title : SW846 8260
QLast Update : Fri Jun 04 19:00:20 2021
Response via : Initial Calibration

