

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062122\
 Data File : VN073014.D
 Acq On : 22 Jun 2022 09:20
 Operator : JC\MD
 Sample : N3301-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-SO-20220609

Quant Time: Jun 22 09:55:25 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 18 06:22:45 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	282018	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	455784	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	421250	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	204524	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	170079	44.607	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	89.220%
35) Dibromofluoromethane	7.957	113	146377	51.959	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	103.920%
50) Toluene-d8	10.375	98	485728	46.336	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	92.680%
62) 4-Bromofluorobenzene	12.669	95	205267	54.456	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	108.920%
Target Compounds						
					Qvalue	
16) Acetone	4.234	43	25744	12.283	ug/l	97
20) Methylene Chloride	5.022	84	40685	12.333	ug/l #	86
40) Benzene	8.392	78	4052	0.338	ug/l #	86
43) Isopropyl Acetate	8.498	43	37757	4.332	ug/l #	88

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062122\
Data File : VN073014.D
Acq On : 22 Jun 2022 09:20
Operator : JC\MD
Sample : N3301-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-SO-20220609

Quant Time: Jun 22 09:55:25 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
Quant Title : SW846 8260
QLast Update : Sat Jun 18 06:22:45 2022
Response via : Initial Calibration

