

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062722\
 Data File : VN073158.D
 Acq On : 27 Jun 2022 13:36
 Operator : JC\MD
 Sample : N3494-07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SP-EXC-1-4

Quant Time: Jun 28 06:37:05 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	305328	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	492136	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	454803	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.609	152	214716	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	185994	45.057	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	90.120%
35) Dibromofluoromethane	7.957	113	111526	36.664	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	73.320%#
50) Toluene-d8	10.380	98	519199	45.871	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	91.740%
62) 4-Bromofluorobenzene	12.668	95	222154	54.582	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	109.160%
Target Compounds						
					Qvalue	
16) Acetone	4.228	43	94154	41.494	ug/l	96
20) Methylene Chloride	5.016	84	56542	15.900	ug/l #	85
25) 2-Butanone	7.275	43	20941	5.816	ug/l #	85
65) Chlorobenzene	11.710	112	10078	1.135	ug/l	98

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062722\
Data File : VN073158.D
Acq On : 27 Jun 2022 13:36
Operator : JC\MD
Sample : N3494-07
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SP-EXC-1-4

Quant Time: Jun 28 06:37:05 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

