

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062422\
 Data File : VN073093.D
 Acq On : 24 Jun 2022 19:25
 Operator : JC\MD
 Sample : N3455-16
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SP-EXC-2-3

Quant Time: Jun 24 23:41:10 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	8.022	168	281581	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	457838	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	423784	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	209304	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	171171	44.963	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	89.920%
35) Dibromofluoromethane	7.951	113	145665	51.474	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.940%
50) Toluene-d8	10.380	98	484156	45.979	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	91.960%
62) 4-Bromofluorobenzene	12.669	95	208554	55.079	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	110.160%
Target Compounds						
					Qvalue	
16) Acetone	4.234	43	70336	33.611	ug/l	90
18) Methyl Acetate	4.793	43	6251	1.122	ug/l #	85
20) Methylene Chloride	5.016	84	145643	44.845	ug/l	90
43) Isopropyl Acetate	8.498	43	31348	3.581	ug/l #	86

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062422\
Data File : VN073093.D
Acq On : 24 Jun 2022 19:25
Operator : JC\MD
Sample : N3455-16
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SP-EXC-2-3

Quant Time: Jun 24 23:41:10 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

