

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042222\
 Data File : VX028290.D
 Acq On : 22 Apr 2022 19:32
 Operator : JC/MD
 Sample : N2511-12 10X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 1591

Quant Time: Apr 23 06:45:11 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041922W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 19 13:38:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	414704	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	708172	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	685453	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	322822	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	300302	55.537	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	111.080%
35) Dibromofluoromethane	5.385	113	272957	54.961	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	109.920%
50) Toluene-d8	8.653	98	1013983	53.158	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	106.320%
62) 4-Bromofluorobenzene	11.079	95	369670	50.686	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	101.380%
Target Compounds						
						Qvalue
16) Acetone	2.380	43	6631	2.634	ug/l	93
18) Methyl Acetate	2.703	43	2287	0.476	ug/l	100
40) Benzene	6.044	78	11464	0.605	ug/l	99
52) Toluene	8.720	92	227180	17.244	ug/l	98

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042222\
Data File : VX028290.D
Acq On : 22 Apr 2022 19:32
Operator : JC/MD
Sample : N2511-12 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
1591

Quant Time: Apr 23 06:45:11 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041922W.M
Quant Title : SW846 8260
QLast Update : Tue Apr 19 13:38:04 2022
Response via : Initial Calibration

