

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010522\
 Data File : VX026356.D
 Acq On : 06 Jan 2022 05:07
 Operator : JC/MD
 Sample : N1020-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-3

Quant Time: Jan 06 12:22:09 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X123021W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 30 08:02:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	151333	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	253699	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	222217	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	101889	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	109840	50.170	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 100.340%	
35) Dibromofluoromethane	5.391	113	82364	49.936	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 99.880%	
50) Toluene-d8	8.653	98	296189	48.278	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 96.560%	
62) 4-Bromofluorobenzene	11.079	95	113804	52.108	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 104.220%	
Target Compounds						Qvalue
16) Acetone	2.386	43	19013	16.028	ug/l	97
31) Cyclohexane	5.477	56	31977	10.256	ug/l	99
39) Methylcyclohexane	7.385	83	6463	2.154	ug/l #	85
40) Benzene	6.044	78	805619	106.108	ug/l	99
52) Toluene	8.720	92	157183	33.628	ug/l	99
67) Ethyl Benzene	10.195	91	1212109	138.867	ug/l	98
68) m/p-Xylenes	10.305	106	644183	197.658	ug/l	94
69) o-Xylene	10.646	106	226697	71.092	ug/l	94
73) Isopropylbenzene	10.963	105	28101	3.548	ug/l	99
78) n-propylbenzene	11.305	91	44549	4.892	ug/l	96
80) 1,3,5-Trimethylbenzene	11.451	105	83566	12.715	ug/l	96
95) Naphthalene	13.774	128	239910	30.466	ug/l	99

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

