

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010622\
 Data File : VX026387.D
 Acq On : 06 Jan 2022 20:31
 Operator : JC/MD
 Sample : N1042-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TRE-1522

Quant Time: Jan 07 04:19:21 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	150938	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	250099	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	213736	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	99074	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	108598	49.733	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	99.460%
35) Dibromofluoromethane	5.391	113	81274	49.985	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.960%
50) Toluene-d8	8.653	98	288236	47.658	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	95.320%
62) 4-Bromofluorobenzene	11.079	95	108231	50.269	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	100.540%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	32449	27.427	ug/l	94
20) Methylene Chloride	2.794	84	11398	5.178	ug/l	95
40) Benzene	6.044	78	9124	1.219	ug/l	94
43) Isopropyl Acetate	6.342	43	13897	2.804	ug/l	95
52) Toluene	8.720	92	43045	9.342	ug/l	98
68) m/p-Xylenes	10.299	106	9751	3.111	ug/l	89
69) o-Xylene	10.640	106	3400	1.109	ug/l	91

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

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