

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062022\
 Data File : VX029631.D
 Acq On : 20 Jun 2022 18:23
 Operator : JC/MD
 Sample : N3339-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WC-GW-20220613

Quant Time: Jun 21 01:52:55 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jun 20 01:31:01 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	260676	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	435384	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	378396	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	160985	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	148233	42.958	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	85.920%
35) Dibromofluoromethane	5.391	113	77381	27.272	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	54.540%#
50) Toluene-d8	8.652	98	494097	47.668	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	95.340%
62) 4-Bromofluorobenzene	11.085	95	174932	44.752	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	89.500%
Target Compounds						
					Qvalue	
16) Acetone	2.385	43	40581	29.505	ug/l	100
17) Carbon Disulfide	2.513	76	1564	0.228	ug/l #	95
20) Methylene Chloride	2.794	84	3798	1.240	ug/l	97
30) Chloroform	5.104	83	40382	7.522	ug/l	97
52) Toluene	8.720	92	22993	3.154	ug/l	100
67) Ethyl Benzene	10.195	91	14181	1.062	ug/l	99
68) m/p-Xylenes	10.305	106	20842	4.001	ug/l	93
69) o-Xylene	10.646	106	13843	2.690	ug/l	96
84) 1,2,4-Trimethylbenzene	11.756	105	14337	1.414	ug/l	94
95) Naphthalene	13.780	128	24340	2.155	ug/l	99

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

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