

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010522\
 Data File : VX026361.D
 Acq On : 06 Jan 2022 07:02
 Operator : JC/MD
 Sample : N1020-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-7A

Quant Time: Jan 06 12:24:09 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	157175	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	264106	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	228258	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	101786	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	113026	49.707	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	99.420%
35) Dibromofluoromethane	5.391	113	85064	49.541	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.080%
50) Toluene-d8	8.653	98	304547	47.684	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	95.360%
62) 4-Bromofluorobenzene	11.079	95	113190	49.785	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	99.560%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	6967	5.655	ug/l	95
31) Cyclohexane	5.477	56	50465	15.584	ug/l	96
37) Ethyl Acetate	4.721	43	4709	1.466	ug/l	96
39) Methylcyclohexane	7.379	83	6380	2.042	ug/l #	87
40) Benzene	6.044	78	1843178	233.198	ug/l	98
52) Toluene	8.720	92	207157	42.574	ug/l	99
67) Ethyl Benzene	10.195	91	206021	22.978	ug/l	97
68) m/p-Xylenes	10.305	106	334641	99.962	ug/l	93
69) o-Xylene	10.640	106	22238	6.789	ug/l	93
73) Isopropylbenzene	10.963	105	74648	9.435	ug/l	97
78) n-propylbenzene	11.305	91	124320	13.666	ug/l	97
80) 1,3,5-Trimethylbenzene	11.451	105	29588	4.506	ug/l	95
84) 1,2,4-Trimethylbenzene	11.750	105	95438	14.547	ug/l	97
95) Naphthalene	13.774	128	246651	31.354	ug/l	100

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

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