

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX011722\
 Data File : VX026485.D
 Acq On : 17 Jan 2022 12:58
 Operator : JC/MD
 Sample : N1160-01 50X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FMI-114

Quant Time: Jan 18 00:40:55 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jan 11 14:45:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	135430	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	227336	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	198497	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	88294	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	93019	49.168	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.340%
35) Dibromofluoromethane	5.391	113	72286	49.239	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.480%
50) Toluene-d8	8.653	98	264372	49.973	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.940%
62) 4-Bromofluorobenzene	11.079	95	98429	52.375	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	104.760%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	3630	3.763	ug/l	99
25) 2-Butanone	4.568	43	2587	2.062	ug/l	95
30) Chloroform	5.099	83	2524	0.851	ug/l	91
40) Benzene	6.044	78	7222	1.127	ug/l	93
52) Toluene	8.720	92	80671	20.536	ug/l	100
67) Ethyl Benzene	10.195	91	34668	4.564	ug/l	98
68) m/p-Xylenes	10.299	106	40537	14.260	ug/l	98
69) o-Xylene	10.640	106	19411	7.077	ug/l	98
78) n-propylbenzene	11.305	91	5659	0.686	ug/l	99
80) 1,3,5-Trimethylbenzene	11.451	105	7793	1.306	ug/l	94
84) 1,2,4-Trimethylbenzene	11.756	105	29227	4.897	ug/l	100

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

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