

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN011725\
 Data File : VN085495.D
 Acq On : 17 Jan 2025 15:36
 Operator : JC\MD
 Sample : Q1115-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 366

Quant Time: Jan 17 22:37:06 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011425W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jan 15 02:16:08 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	171685	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	338097	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	310797	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	116685	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	154623	55.793	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	111.580%	
35) Dibromofluoromethane	8.165	113	114448	48.793	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	97.580%	
50) Toluene-d8	10.565	98	425716	51.083	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	102.160%	
62) 4-Bromofluorobenzene	12.847	95	131193	46.020	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	=	92.040%	
Target Compounds						
					Qvalue	
16) Acetone	4.424	43	132400	147.859	ug/l	98
18) Methyl Acetate	5.036	43	3335	1.225	ug/l #	82
20) Methylene Chloride	5.277	84	2435	1.098	ug/l #	74
25) 2-Butanone	7.488	43	14246	10.811	ug/l	93
43) Isopropyl Acetate	8.688	43	233385	43.835	ug/l #	77
51) 4-Methyl-2-Pentanone	10.441	43	15731	5.092	ug/l #	83

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

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