

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091522\
 Data File : VN074492.D
 Acq On : 16 Sep 2022 04:29
 Operator : JC\MD
 Sample : N4630-25
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-17-4-5-GR

Quant Time: Sep 16 07:02:00 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091522W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 15 19:20:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	370139	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.904	114	717339	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	713879	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	301995	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	270812	43.121	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	86.240%
35) Dibromofluoromethane	7.951	113	221490	46.481	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	92.960%
50) Toluene-d8	10.381	98	827457	44.901	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	89.800%
62) 4-Bromofluorobenzene	12.674	95	326344	51.704	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	103.400%
Target Compounds						
					Qvalue	
16) Acetone	4.228	43	77718	19.339	ug/l	# 85
17) Carbon Disulfide	4.457	76	37005	2.768	ug/l	97
18) Methyl Acetate	4.787	43	13417	1.075	ug/l	# 76
20) Methylene Chloride	5.016	84	23106	3.794	ug/l	# 84
37) Ethyl Acetate	7.340	43	82705	6.349	ug/l	# 95
43) Isopropyl Acetate	8.492	43	338691	16.600	ug/l	# 85

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

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