

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081623\
 Data File : VN078907.D
 Acq On : 16 Aug 2023 16:54
 Operator : JC\MD
 Sample : 04022-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-16

Quant Time: Aug 17 01:18:47 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081523W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 16 02:42:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	147706	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	293377	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	293102	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	85945	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	129421	56.308	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 112.620%		
35) Dibromofluoromethane	8.177	113	101828	51.562	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 103.120%		
50) Toluene-d8	10.571	98	383459	52.293	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 104.580%		
62) 4-Bromofluorobenzene	12.853	95	114958	47.963	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 95.920%		
Target Compounds						
						Qvalue
16) Acetone	4.448	43	26736	28.192	ug/l	88
18) Methyl Acetate	5.036	43	5118	1.277	ug/l #	69
20) Methylene Chloride	5.283	84	15687	7.762	ug/l #	77
25) 2-Butanone	7.500	43	7752	5.098	ug/l #	74
37) Ethyl Acetate	7.577	43	4479	1.348	ug/l #	75
43) Isopropyl Acetate	8.694	43	211053	35.925	ug/l #	71
51) 4-Methyl-2-Pentanone	10.447	43	36371	10.438	ug/l #	75

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

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