

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072523\
 Data File : VN078647.D
 Acq On : 25 Jul 2023 15:12
 Operator : JC\MD
 Sample : 03717-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-2

Quant Time: Jul 26 04:47:27 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071023W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 10 16:13:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.229	168	420745	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	741515	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	710624	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	247921	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	258179	45.772	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	91.540%		
35) Dibromofluoromethane	8.176	113	266113	52.867	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	105.740%		
50) Toluene-d8	10.570	98	903127	47.494	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	94.980%		
62) 4-Bromofluorobenzene	12.853	95	302301	51.612	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	103.220%		
Target Compounds					Qvalue	
16) Acetone	4.441	43	86088	48.996	ug/l	91
20) Methylene Chloride	5.288	84	27638	4.368	ug/l	97
25) 2-Butanone	7.494	43	47499	17.174	ug/l #	81
37) Ethyl Acetate	7.577	43	13920	2.297	ug/l #	96
43) Isopropyl Acetate	8.700	43	351939	35.282	ug/l #	80
51) 4-Methyl-2-Pentanone	10.453	43	150397	25.348	ug/l	91

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

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