

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN011625\
 Data File : VN085470.D
 Acq On : 16 Jan 2025 15:22
 Operator : JC\MD
 Sample : Q1069-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RW7B-CARBON-20250109

Quant Time: Jan 17 04:34:54 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	8.218	168	198665	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.094	114	358567	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.859	117	314710	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	128541	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	172160	53.685	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 107.360%		
35) Dibromofluoromethane	8.159	113	125899	50.611	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 101.220%		
50) Toluene-d8	10.565	98	443352	50.163	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 100.320%		
62) 4-Bromofluorobenzene	12.847	95	148587	49.146	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	= 98.300%		
Target Compounds						
						Qvalue
9) 1,1,2-Trichlorotrifluo...	4.377	101	35395	14.795	ug/l	99
16) Acetone	4.424	43	13081	12.624	ug/l	98
24) 1,1-Dichloroethane	6.565	63	21254	4.539	ug/l	99
30) Chloroform	7.965	83	9847	2.035	ug/l	100
42) 1,2-Dichloroethane	8.665	62	3778	0.956	ug/l	81
44) Trichloroethene	9.353	130	3876	1.587	ug/l	91

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

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