

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN011525\
 Data File : VN085460.D
 Acq On : 15 Jan 2025 14:37
 Operator : JC\MD
 Sample : Q1069-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RW7B-CARBON-20250109

Quant Time: Jan 16 00:59:06 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	181061	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	354107	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	288035	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	108344	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	162792	55.699	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	111.400%	
35) Dibromofluoromethane	8.159	113	122198	49.742	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	99.480%	
50) Toluene-d8	10.565	98	413472	47.371	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	94.740%	
62) 4-Bromofluorobenzene	12.847	95	122853	41.146	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	=	82.300%	
Target Compounds						
						Qvalue
9) 1,1,2-Trichlorotrifluo...	4.371	101	36336	16.665	ug/l	99
16) Acetone	4.430	43	13082	13.853	ug/l	100
24) 1,1-Dichloroethane	6.565	63	21243	4.978	ug/l	98
30) Chloroform	7.965	83	9663	2.191	ug/l	99
42) 1,2-Dichloroethane	8.677	62	4024	1.031	ug/l	86
44) Trichloroethene	9.353	130	3555	1.474	ug/l	86

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

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