

Student number	point total	req total	extra total	R1 Euler integrator (1p)	R2 Spring system (2p)	R3 Trapezoid integrator (2p)	R4 Pendulum system (2p)	R5 Cloth system (3p)	mod	notes	RK4 (2p)	Spray system (1-3p)	Wind (2p)	Mouse drag/ poke (2-3p)	Frictionl. coll. (2p)	Particle spline editor (2-3p)	Cloth tearing (1+p)	Particle rendering (1-4p)	Implicit integr (8-10p)	GPU stuff (4+p)	Other extras (7p)	What other extras
225157	0	0	0																			
270034	5	5	0	1	2	2	0	0														
293846	0	0	0																			
295323	0	0	0																			
345642	0	0	0																			
348843	0	0	0																			
										RK4: final linear combination is unweighted (missing factors 1,2,2,1)			2									
349936	13.5	10	3.5	1	2	2	2	3			1.5		2									
350475	10	10	0	1	2	2	2	3														
352091	0	0	0																			
353980	0	0	0																			
										Empty readme. Only requirements checked - let us know if we missed something. R3: you seem to be using only f1 in the computation of the output state instead of 0.5*(f0+f1). xt is unused?												
354439	8	9	0	1	2	1	2	3	-1													
355593	0	0	0																			
356026	12	10	2	1	2	2	2	3			2											
361749	0	0	0																			
369181	12	10	2	1	2	2	2	3			2											
372660	1	1	0	1	0	0	0	0														
387370	0	0	0																			
425575	0	0	0																			
425614	0	0	0																			
426419	0	0	0																			
427489	7	7	0	1	2	2	2	0														
428022	0	0	0																			
429487	0	0	0																			
430829	1	1	0	1	0	0	0	0														
457598	0	0	0																			
460297	0	0	0																			
464772	0	0	0																			
46477D	0	0	0																			
46596K	0	0	0							R4+R5: you are using the variable current_state_ in the updates which contains the initial state of the step (but which is not modified by the integrators until the end of the step) - instead you should be using the State state, that is passed to evalF. This leads to the instability you described. The naming is a bit unfortunate here.												
474199	13	9	4	1	2	2	1.5	2.5			2					2						
474322	0	0	0																			
474458	10	10	0	1	2	2	2	3														
										R2: spring force function computes length only using x and y coords for a vec3f.												
474898	4.5	4.5	0	1	1.5	2	0	0														
475389	0	0	0																			
475813	0	0	0																			
475910	0	0	0																			
										R5: evalF is nearly unreadable and static to this particular configuration, but seems to work. RK4: cool use of the assignment operator mid arithmetic - I haven't thought about it that way before												
477329	12	10	2	1	2	2	2	3			2											
477811	12	10	2	1	2	2	2	3			2											
478328	1	1	0	1	0	0	0	0														
478470	0	0	0																			
478687	0	0	0																			

[illegible]

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549040	5	3	2	1	0	2	0	0		R2: spring force should have direction of the line connecting pos1 and pos2. Drag force has the wrong sign and is proportional to the location, not the velocity of the particle.	2											
549163	10	10	0	1	2	2	2	3		R3: looks like you are not clearing springs in reset, so the system feels a bit more rigid (-0p)												
55055P	12	10	2	1	2	2	2	3		Cool wind, very violent!	2											
552794	3	3	0	1	0	2	0	0														
552969	0	0	0																			
554598	12	10	2	1	2	2	2	3		Using std::transform, cool!	2											
563068	5	4	1	1	1	2	0	0		Your Euler-step actually takes two steps? R2: f[2] should not directly depend on the current force, but is rather the velocity-component of the current state. RK: You need no evaluate the forces after each RK-step and use that in the next part. The simple system diverges with your implementation.	1											
576149	0	0	0																			
585716	0	0	0																			
586333	0	0	0																			
586702	1	1	0	0.5	0.5	0	0	0		R1: eulerStep only increments the first element of the state, need a loop here! R2: fDrag should take in the current velocity, not the acceleration and should oppose the movement (flip signs!).												
586980	7	7	0	1	2	2	2	0														
587316	0	0	0																			
587471	17.5	9.5	8	1	2	2	2	2.5		R5: clearing the springs_ vector at reset() helps for the instability! Cloth GPU sim with RK4.	2								6			
588289	10	10	0	1	2	2	2	3														
589291	0	0	0																			
589343	0	0	0																			
589848	3	3	0	1	2	0	0	0														
590921	0	0	0																			
591904	5	5	0	1	2	2	0	0														
591946	7	7	0	1	2	2	2	0														
592929	0	0	0																			
593274	10	10	0	1	2	2	2	3														
593847	0	0	0																			
594435	0	0	0																			
595926	12	10	2	1	2	2	2	3		R4: you are not clearing the springs vector at reset(), hence the different behavior (-0p)	2											
595997	0	0	0																			
596747	0	0	0																			
597429	10	10	0	1	2	2	2	3														
597623	11	10	1	1	2	2	2	3		RK4: k1 should use half step, k3 full step. Also, k4 not necessary, final step direction given by lin. comp of f's.	1											
597937	38	10	28	1	2	2	2	3		Cloth GPU sim with RK4. The implicit integrator convergence criterion is somewhat looser than in the example. Not entirely sure where the around 3x perf drop wrt reference comes from: would need to have a closer look at performance metrics.	2	3	2		2			8	6		RKF45 (3p), Impl. opt 5 (2p)	

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652335	16	10	6	1	2	2	2	3		Moving cloth stationary points is a lot of fun!	2			3			1					
652649	16	10	6	1	2	2	2	3		Neat use of lambdas!	2		2		2							
652898	0	0	0																			
652937	10	10	0	1	2	2	2	3														
										Your implicit integrator is really just an explicit integrator since the Jacobian is the zero matrix and the solution of Ax=b is trivial.	2								2			
653127	14	10	4	1	2	2	2	3														
653596	0	0	0																			
										NB: your submission zip was flagged by Windows Defender												
653693	10	10	0	1	2	2	2	3														
653871	0	0	0																			
653907	10	10	0	1	2	2	2	3														
653910	10	10	0	1	2	2	2	3														
										Particle rendering: textured cloth with shading.												
654595	24	10	14	1	2	2	2	3			2	3	2	3	2			2				
										R5: non-square shape not significant enough to warrant extra points I'm afraid												
655057	10	10	0	1	2	2	2	3														
										R4: you are not clearing the springs vector at reset() (-0p). Flat shaded cloth mesh.								1.5				
655086	8.5	7	1.5	1	2	2	2				2											
655109	12	10	2	1	2	2	2	3														
655251	0	0	0																			
655264	0	0	0																			
655471	10	10	0	1	2	2	2	3														
655691	0	0	0																			
655853	13	10	3	1	2	2	2	3			2							1				
										R4: you are using the class member <code>current_state</code> instead of the supplied argument state. This causes the instabilities.												
656250	6.5	6.5	0	1	2	2	1.5	0														
656454	16	10	6	1	2	2	2	3			2		2		2							
656616	10	10	0	1	2	2	2	3														
657291	12	10	2	1	2	2	2	3			2											
657314	0	0	0																			
										RK4: not quite right, you are not taking the weighted average of the values k at the end. Spray system: you are not supposed to use the <code>current_state</code> variable in <code>evalF</code> , but rather the function arg state in order to construct a new state that you are free to modify as you wish. You could store the indices of the active particles e.g. in the <code>SpraySystem</code> instance and read from there.												
657327	14	10	4	1	2	2	2	3			1.5	0.5	2									
657482	0	0	0																			
657796	10	10	0	1	2	2	2	3														
657893	10	10	0	1	2	2	2	3														
659914	12	10	2	1	2	2	2	3			2											
										The collision-extra is quite unstable. Your implicit integrator is actually just an overly complicated and slow explicit integrator. You are never computing the Jacobian (which you should do analytically, not with finite differences!) and essentially just conveniently setting things up such that the result of the Ax=b equation is ~ step^2 F.												
660246	17	10	7	1	2	2	2	3			2		2		1				2			

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660877	0	0	0																			
660893	0	0	0																			
663191	0	0	0																			
663272	0	0	0																			
665380	10	10	0	1	2	2	2	3														
665898	10	10	0	1	2	2	2	3														
666172	12	10	2	1	2	2	2	3			2											
										R4+R5: you are using the variable current_state in evalF instead of the function argument state, which leads into the instability.	2											
666350	11	9	2	1	2	2	1.5	2.5			2											
666680	10	10	0	1	2	2	2	3														
667249	10	10	0	1	2	2	2	3														
67137M	0	0	0																			
67627H	15	10	5	1	2	2	2	3			2	3										
677734	0	0	0																			
678089	0	0	0																			
68933B	0	0	0																			
69247N	0	0	0																			
700436	0	0	0																			
																					Adaptive RKF45 (2p), Crank–Nicolson (4p), sparsity patterns (4p), Constraints (8p), GPU sprinkler and GPU integrators (+5 to GPU points)	
705570	69	10	59	1	2	2	2	3			2	3	2	3	2	3	3	4	10	9	18	
706566	0	0	0																			
										GPU cloth: evalF writes out of bounds - last if needs to check x < w and y < h (-0p)	2	3								5		
708784	20	10	10	1	2	2	2	3			2	3										
708904	0	0	0																			
708920	12	10	2	1	2	2	2	3							2							
708933	0	0	0																			
709291	5.5	5.5	0	1	2	2	0.5															
709628	12	10	2	1	2	2	2	3			2											
										fSpring: can be replaced with a single expression: '-k * ((p2 - p1).length() - rest_len) * (p2 - p1).normalized()' (handout section 1.2, third equation)												
710086	10	10	0	1	2	2	2	3														
710497	15.5	10	5.5	1	2	2	2	3			2	2						1.5				
										Spary: no collisions												
710743	13	10	3	1	2	2	2	3			2		1									
710976	5	5	0	1	2	2	2															
										Wind: looks very artificial, almost no variation over time												
711182	5	5	0	1	2	2	2			R3: bug in first for loop: x0[0] instead of x0[i]												
										Wind: random noise cast to int, which essentially removes all time-dependent variation												
711467	18	10	8	1	2	2	2	3			2	3	1		2							
711551	12	10	2	1	2	2	2	3			2											
711810	9	7	2	1	2	2	2				2											
711904	0	0	0																			
712550	10	10	0	1	2	2	2	3														
712686	0	0	0																			
										R4: indexing in first for-range loop wrong. R5: positions only not quite enough for points.												
712819	6	6	0	1	2	2	1	0														
712958	0	0	0																			
713672	0	0	0																			
										R4, R5: not dividing force by mass to get acc. Generally: please try to match the reference visually (makes debugging and grading easier)												
714985	13	9	4	1	2	2	1.5	2.5			2		2									

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716080	3	3	0	1		2																
										RK4: you need to call evalF for the different states, currently they are incorrectly treated as derivatives												
716462	10.5	10	0.5	1	2	2	2	3			0.5											
716718	0	0	0																			
										R5: logic seems fine, probably something with gravity or spring tension												
716860	9.5	9.5	0	1	2	2	2	2.5														
717377	0	0	0																			
717474	0	0	0																			
717513	12	10	2	1	2	2	2	3			2											
717539	0	0	0																			
718020	0	0	0																			
718871	0	0	0																			
722427	0	0	0																			
										RK4: one step too much, final dir given by sum of f0...f3 (weights also don't sum to 6)												
723691	6	5	1	1	2	2					1											
723905	0	0	0																			
728667	17.5	10	7.5	1	2	2	2	3			2									5.5		
728900	12	10	2	1	2	2	2	3			2											
729132	0	0	0																			
										R2: no spring force, evalF reads from wrong state indices												
729967	1	1	0	1	0																	
730309	0	0	0																			
732080	0	0	0																			
732255	0	0	0																			
732323	0	0	0																			
732336	0	0	0																			
732352	0	0	0																			
										Sprinkler: no collisions (even though readme claims so?)												
732459	14	10	4	1	2	2	2	3			2	2										
76509T	12	10	2	1	2	2	2	3			2											
										Wind: doesn't vary smoothly over time, looks artificial. Pretty unique tearing setup!												
765510	15.5	10	5.5	1	2	2	2	3			2		1.5				2					
766331	0	0	0																			
767042	0	0	0																			
										R2: fSpring: using sqr, not sqrt. R1, R3: not looping over points, just evolving first one. R4: broken, but overall idea seems OK												
767136	3.5	3.5	0	0.5	1.5	1	0.5															
768504	0	0	0																			
769396	12	10	2	1	2	2	2	3			2											
										RK4: need to evaluate forces (evalF) at the intermediate steps rather than just sum the the contributions.												
77388B	10.5	10	0.5	1	2	2	2	3			0.5											
779124	0	0	0																			
										R1: taking two steps instead of one, you need just s0[i] + step*f0[i]												
780058	0.5	0.5	0	0.5																		
780346	10	10	0	1	2	2	2	3														
782917	0	0	0																			
783563	0	0	0																			
783709	0	0	0																			
786667	0	0	0																			
78708M	0	0	0																			
787543	0	0	0																			
787640	0	0	0																			
788380	0	0	0																			
788678	0	0	0																			
791982	0	0	0																			
795700	0	0	0																			
795755	0	0	0																			

[illegible]