Chemistry Homework and Assessments



"SFU Chemistry used CAPA from 1996 and in 2001 migrated to LON-CAPA.



Summary of regular an intermittent usages. Orange text -- in constant use. SFU: trimester system with different people teaching -- each with their own preferences and methods.

2017, taking stock: creating new resources, making obsolete some old resources. Organizing sequences according to topics.



3-d animated visualizations. SFU resources which used JMOL, are now revised to use JSMOL.

2758 mol or pdb files published in LON-CAPA.1883 in MSU domain.630 in NAU domain.129 in SFU domain.107 in FSU domain.



This question has been used many years, but is still considered difficult.

In part to avoid scamming of answers off the web, I made new differing versions of this problem, which were intended to be a bit easier AND have different questions and phrasing than the old one — so it is not as easy to search out.



Version of problem used in Spring 2017: Students are quick to use 'tutoring' services online. Google search for words in the problem statement find services with answers. NOT free though. \$\$



Only one student had this particular rendition of this question.

They submitted a photo of their display to a commercial homework help site.

The student was working up to near the deadline on the assignment,

but did not receive the answer in time.

Viewed Sun Jan 15, 2017 6:25PM until 11:25PM (due at 11:59PM).

Student's final view or their problem (at lower right).

Gravimetric Analysis Binary Mixture (2003 - 2016 – 130 courses)											
Antimony is alloyed with lead to i 1.011 g of a particular metallic al What was the percentage (by ma %	ncrease the rigidit loy, compounded ss) of antimony in	y of compone of only Pb an the alloy?	ents used in the constructio d Sb, can be quantitatively	n of lead stor converted in	rage batteries. to a 1.172-g mi	xture of th	e oxides Pt	oO ₂ an	d Sb ₂ O	4.	
Submit Answer Tries 0/5 Overall Assessment Statistical Data Recent Detailed Assessment Statistical Data											
Statistics calculated for number of students:	13319	Domain	Course	(Rumahu)	Section	(s)	Num Students	Part	Mean Tries	Degree of Difficulty	Deg Discri
Average number of tries till solved:	2.39	sru	Chemizi 2016-3 Lecture	e (Burnaby)		an	434	0	2.20	0.64	
Degree of difficulty:	(0.65)										
Degree of discrimination:	(0.21)										
Gravimetric Anal	ysis Bi	nary M	ixture		(2017	- 20	courses	S)			
Amounts of tin may be alloyed with co 7.8470 g of a particular bronze, comp (Use the atomic masses: Cu 63.546, S	opper to produce bro ounded of only Cu a Sn 118.71, O 15.999	onzes of varyin Ind Sn, can be 94.)	g composition and properties. quantitatively converted into a	a 9.8429-g mi	xture of the oxide	s CuO and S	SnO ₂ .				
What was the percentage (by mass) of %	of tin in the alloy?										
Overall Assessment Statist	ical Data										
Statistics calculated for number of students:	367										
Average number of tries till solved:	2.32										
Degree of difficulty:	(0.68)	Recent D	etailed Assessment Sta	atistical Da	ta						
Degree of discrimination:	(0.00)	Domain sfu (Course Chem121 2017-2 Lecture	Section(s) N all	um Students P 130	art Mean 1 0 2	Fries Degr e	ee of D	0.69	y Degree of	Discrim
ulser initiation.		sfu (Chem121 2017-1 Lecture	all	237	0 2	2.38		0.68	3	

Goggle-search finds a solution for the old problem, but not (yet) for the new problem. Average number of tries by student and the overall degree of difficulty similar.

	Bond Energies and Heat of Reaction			
	Use the given bond energy values to estimate ΔH for the following gas-phase read	tion.		
2004-2016 8 variations DoD=0.61 <tries> = 2.28</tries>	$H \rightarrow \left(\begin{array}{c} H \rightarrow H \\ H \rightarrow H \end{array} \right) \xrightarrow{H} \left(\begin{array}{c} H \rightarrow H \\ H \rightarrow H \end{array} \right) \xrightarrow{H} \left(\begin{array}{c} H \rightarrow H \\ H \rightarrow H \end{array} \right)$ (Simple energy units required for the answer.)			
2017		0 = 432	Due = 363	[Du = = 413
26 variations DoD=0.67		$D_{\text{H-H}} = 432$ $D_{\text{H-N}} = 391$ $D_{\text{C=C}} = 614$	$D_{H-O} = 467$ $D_{C=C} = 839$	$D_{\rm H-C} = 413$ $D_{\rm C-C} = 347$ $D_{\rm C-N} = 305$
<tries>=2.35</tries>		$D_{C=N} = 615$ $D_{C=O} = 745$	D _{C≡N} = 891 D _{C=O (CO₂)} = 799	$D_{C=0} = 358$ $D_{C=0} = 1072$
		$D_{C-Br} = 276$ $D_{N\equiv N} = 941$ $D_{N-O} = 201$	$D_{N-N} = 160$ $D_{O-O} = 146$ $D_{N=O} = 607$	$D_{N=N} = 418$ $D_{O=O} = 495$ $D_{Br-Br} = 193$
	Submit Answer Tries 0/5			
	Threaded View Chronological View Other Views Export			
	Anonymous 1 Reply (Tue Oct 25 06:06:17 pm 2016 (PDT))			
	Just in case if anyone wants to know:			
	Hold On, Don't Panic!			
	Okay, this question may seem very scary, but it can be broken down easily	!		
	 Look carefully at the diagram, this isn't something new; the last two q The table offers WAY more information than you actually need, worry A strategic recommendation is to work left to right, write out what the It is also a good idea to keep the molecules separate of each other un be careful as solid lines aren't the only bondi Elements adjacent to ea Note that double (or triple) bonds are NOT the same as single bonds Once again, keep in mind the positive and negative signs 	uestions are ju not bonds are, and til the sum at tl ch other are bo	st the same! d how many there ar he end - that way, m onds too; they will r	e listakes can be avoided leed to be considered
	Take a deep breath, it's okay! Not too bad, right?			

Creating new problems with more variations.

LON-CAPA student discussion/help for others seemed to work well in this instance, without resort to other online 'help' sites. Suspect that student was likely getting 'help' before posting to the in-course discussion.

Diffusion Partial I	Pressure Mole Fraction	
A	c	
Three glass bulbs, joined by closed stopcocks Bulb &: 250. mL of Kr(g) at 570. torr Bulb B: 500. mL of CO ₂ (g) at 0.500 atm Bulb C: 1.25 L of N ₂ (g) at 25.331 kPa After both stopcocks are opened and the gase Plotal = 152 kPa lpts Submit Answer Incorrect. Tries 3/5 Previou What is the partial pressure of Kr(g)? Pkr = lpts Submit Answer Tries 0/5 What is the mole fraction of CO ₂ (g)?	, have the following volumes and initial pressures of the specified gases as allowed to diffuse throughout, what will be the final total pressure? <u>s Tries</u>	
X _{CO2} = 1pts Submit Answer Tries 0/5		
For question 1: I think the total pressure of a and how do you find the pa Show transcribed image text	Get this answer with Chegg Study VIEW THIS ANSWER	
Expert Answer	OR	

Chegg ("chicken or egg") : originally started for/by students at Iowa State U.

- is an online textbook rental company, also providing:

homework help (answers) and online tutoring.

It is based in Santa Clara California.

Some students seem prompt to access this service and pay\$.

James () Diffusion F	Diffusion Partial Pressure Mole Fraction Diffusion - Partial Pressure - Mole Fraction Viewed by < 324 students <u>only</u> in Spring 2017										
Netwo using	Network-wide courses using resource: • Chem121 2016-3 Lecture (Burnaby) (sfu) • Chem121 2017-1 Lecture (sfu) • Chem121 2017-2 Lecture (sfu) • Chem121/120 2017-2 Surrey Lecture (sfu) • Testing in 2.11 (sfu)										
Overall	Assessment Statisti	ical Data	_								
Statis numb	tics calculated for er of students:	270									
Avera till sol	ge number of tries ved:	1.90									
Degre	e of difficulty:	(0.50)									
Degre	e of discrimination:	(0.00)									
Recent	Detailed Assessmen	t Statistical D	ata								
Domain	Course	Section(s)	Num Students	Part	Mean Tries	Degree of Difficulty	Degree of Discrimination		Time of com	putation	
sfu	Chem121 2017-1 Lect	ture all	270	12	2.31	0.60	0.00	Fri May	26 10:20:18	am 2017	(PDT)
sfu	Chem121 2017-1 Lect	ture all	261	13	1.63	0.42	0.00	Fri May	26 10:20:18	am 2017	(PDT)
sfu	Chem121 2017-1 Lect	ture all	252	15	1.74	0.47	0.00	Fri May	26 10:20:18	am 2017	(PDT)
You	u are correct. Compute	er's answer now	shown ab	ove.							

Spring 2017 — Problem restricted to one class of SFU students only.

271 out of 323 registered students (84%) attempted this question.

42 of whom (15%) got all or part of it wrong (5 tries allowed for each part)

Can easily identify the student who posted to CHEGG, because of the many variations of the problem.

This student used his/her 3rd try to input 152 kPa. Feb 3 2017, at 1:18PM, before posting for help.

On 5th and last try they entered the correct answer as 0.375 atm on Sunday Feb 5, 2017 at 4:56 PM. (C student.) There was only one other student with an identical variation of the problem (in a class of 323)— and that person's submission profile did not match.



Another problem of which Photo submitted to CHEGG could only have been taken March 29 – April 21 2017 Only one student in course with this combination/permutation of the problem.

In general: Attempted by 223 of 323 students (83%), 45 of whom (20%) never got it correct, despite answer(s) offered on commercial 'help' site.



Student first viewed this problem at 5:08PM Sun April 9, 2017.

Answered it on their first try at 11:07PM Sun April 9, 2017 (Due date), presumably after having received the answer from the commercial 'help' site.

C student.



Same problem using JSMOL visualization of molecules as well as the 2-dimensional line drawings.



New problem (bottom) possibly a little easier than old problem (top). Google search for postings containing the wording of the problem. Next slide



Student submitted their problem to CHEGG.

Only two students had this specific variation.

Neither of them made ANY attempt to answer this problem

(presumably because no answer was forthcoming).

One was working on the assignment over 3 days culminating on the due date.

The other only worked on the assignment on the due date sporadically from 5:41PM until 11:16PM. (Due at midnight).



Demo of new rendition of old problem.



New rather easy problem (2017). Nonetheless students still submitted it to commercial web sites for answers.



HIDE all until UNKNOWN ID has been accepted. Necessary for auto-grading. One try only -- grading occurs simultaneously with submission.

Randomized CORRELATION OF UNKNOWN ID## is generated within the program and provided to Lab technician for filling/labelling sample vials, randomly distributed to students.

NEXT, interactive "Chart" of student scores for their determined result Wt% Ca.

kaymond Johi	n Batchelo	r (Course Coordinal	tor) Chem	istry 21	5 Laboratory	2014-2 (More)	R.	New Messages Rol	es Help Logo
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Chemistry 21	5 Laborator	y 2014-2 » Chart							Chart 💽
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Droplists to select filters for display or output of Chart. Note blue "?"s which appear liberally throughout the interface, give contextual help. Displayed part scores (integers or single-character symbols) are each a LINK to the grading inter face for that particular resource and the individual student. Very COOL! When the exact score cannot be represented by an integer, an asterisk is displayed which then pops the actual score when moused-over. Next, click on the Expt B result for the student who scored 3.2/5.

View of the problem								
Due in 4 hours, 1 minute								
Exp B. TGA Results								
Enter the following data and	calculated results from your lab no	tebook:						
FIRST, CAREFULLY enter yu It must be correctly entered Subsequent results input fie Unknown Sample Answer Submitted: Your fr	ur Unknown Sample Number. as a simple integer, between 1 and ids will not become available until t # = 98 inal submission will be graded after	1 999. his has been done. the due date. Tries 1/1 Previous	Tries					
Sample	Mass	Heating Rate						
CaC ₂ O ₄ ·xH ₂ O	4.864480mg	30°C/min						
Polyethylene glycol	5.266522mg	30°C/min						
Unknown sample	5.676756mg	30°C/min						
Antacid tablet	7.668478mg	30°C/min						
Submit Answer Answer Sub	omitted: Your final submission will b	e graded after the due date. Trie	s 3/10 Previous Tries	Submitted				
2. Enter the Unknown weigh	nt percent Ca ²⁺ and its uncertainty.	#	Try	Answer				
(The submitted values will b	e entered in the table in Report I.)	1 will be gr	aded after the due date. (Try 1)	22.64 0.0003				
uncertainty is absolute, (not	relative).)	2 Answer S will be gr	aded after the due date. (Try 2)	22.64 5.36				
Wt% Ca = 22.64 ±	: 7.07	3 Answer S will be gr	ubmitted: Your final submission aded after the due date. (Try 3)	22.64 7.07				
Answer Sur	Sincea. Four mai submission will b	ine graded arter the dde date. The	S J/ 10 PIEVIOUS TITES					

Note student's reported Wt%Ca and the unreasonably large Uncertainty. Can pop-up the students "Previous Tries".

'Uncertain about uncertainty'. This student seems to remain confused about the meaning of "absolute" vs "relative" uncertainty.

Automatic score of 3.2/5 arises from granting 4/5 for accuracy of Wt% Ca, which is reduced to 3.2/5 for inappropriate uncertainty.



New problem sample using dynamically generated gnu plots for different quantum numbers.

Creating a new problem resource.

The requested file /priv/batchelo/testing/JME/demo_organicresponse_CSC.problem currently does not exist.

To create a new problem, select a template from the list below. Then click on the "Create problem" button.

Algebraic Response Problems

Formula Response using Computer Algebra System Example

Math Response using Computer Algebra System MAXIMA () Example Math Response using Computer Algebra System R () Example Math Response using Computer Algebra System and Hints () Example

OUnordered Multi-Answer Formula Response Problem Example

Chemistry Problems

Chemical Reaction Response Example Chemical Reaction Response with Hints Example Organic Material Response Example Organic Material Response with Hint Example

Free Form Problems

Create problem

Custom Response @ Example Custom Response using Computer Algebra System and Hints @ Example Custom Response with Partial Credit @ Example Extring Response @ Example String Response @ Example

String Response with Pre-Processing () Example

Handgraded Problems

Essay Response Example

Input-Dependent Problems

◯ Using Learner Answer in Multipart Numerical Problem Example ◯Using Learner Formula in Graph with Formula Response Example ◯Using Learner Formula in Graph with Math Response Example

Miscellaneous

Blank Problem Example
Click-On-Image Problem Example
Simple Formula Problem Example

Multiple Choice Problems

Matching Response Example Option Response - Concept Groups Example Option Response - Matching & Example Option Response - True/False Example Radio Button Response & Example Randomity Labelled Image with Option Response & Example Rank Response & Example

Numerical Problems

Curve Plot with Numerical Response Example Oata Plot with Numerical Response Bearple Numerical Response & Example Numerical Response with Custom Units Example Numerical Response with Pre-Processing Example One of Multiple Answers Numerical Problem Example

Many templates as shown by the offerings when initiate the creation of a new resource. Programming definitely NOT required, but can be advantageous. Can mix and match or devise your own special types of assessments. Have not even attempted ALL of these templates. Let's try something a bit different... In CSTR, create New problem "SampleOrganicResponse" -- click on "Go"

Problem Testing		
Problem Status:	; Problem Type:	🗧 🗧 Show All Foils
Feedback Mode:	; Apply style file:	Select
Language:	;) Math Rendering:	•
Change View Show Default View	Reset Submissions	
New Randomization Change Rand	om Seed To: 1305694014 Calculate answers for 20 VE	ersions. 🔞
Edit		
The image below is acetic	acid C ₂ H ₄ O ₂	
H ₃ CO OH		
Complete acetic acid.		
Submit Answer Tries 0		

This is the organicresponse TEMPLATE. Idea is to present a molecule and require the student to draw a molecule, not necessarily the same molecule though. Will EDIT the question and use a molecule which displays potential stereoisomerism. Click on "Edit" to go to the GUI editor.

Construction Space								
Functions 💊 📃 🚍	a							
Construction Space: /priv/batchelo/testing/JME/demo_organicresponse_CSC.problem Recent :								
Problem Editing 🕐								
Discard Edits and View EditXML Undo			Save and Ed	it Save and View				
Insert: Image	\$							
Text Block Delete? =	Edit Math	Greek Symbols 🔞	Other Symbols 🔞	Output Tags 🔞				
Rich formatting » The image above is phen	oxymethylpenicillinK/b>.							

Insert an image "tag" . I'm going to ask for a specific stereoisomer of penicillin V". Click "Save and Edit"

Construction Conco
Functions 😱 📃 🗮 🚢
Construction Space: /priv/batchelo/testing/JME/demo_organicresponse_CSC.problem
Recent \$
Problem Editing ()
Discard Edits and View EditXML Undo Save and Edit Save and View
(nsert:
Image Delete? 💶 🗧 Image Options 🛞
Image Url: /res/sfu/batchelo/testing/JME/Phenoxymethylpenicillin.png Select Search
Description: Penicillin V
width (pixel): height (pixel):
TeXwidth (mm): TeXheight (mm): Alignment: TeXwrap: ÷ Encrypt URL: no
O H S S S S S S S S S S S S S S S S S S
Insert:
Text Block Delete? Edit Math
Rich formatting »
The image above is phenoxymethylpenicillin .

IF one knows where the image is located they can just type it in, OR one could use "Select" or "Search" to either browse or search in repository for a published image.

Click "Save and Edit"



Notice the pencil icon just after the "Starting Molecule" field. Click on it to use applet to insert and save atomic coordinates & bonds for "starting" point drawing.

Next click the pencil icon over on the right and draw the CORRECT answer with full stereochemistry using wedge bonds.

Clicking "Insert Answer" in this case does two things:

Saves the coords and bonds needed to produce the "correct" answer displayed.

Saves and inserts the SMILES (Simple Molecular Input Line-Entry System) string which uniquely corresponds to the structure, including stereochemistry. This is required for automatic grading of the student's submission. NEXT Scroll down to SAVE and VIEW the problem you have created.



So, now we have a simple problem to draw a specified stereoisomer of the original molecule. NEXT Edit XML to show the simple code.

Construction Space » Problem Editing								
Functions 🕋 🗮 📇								
Construction Space: /priv/batchelo/testing/JME/demo_organicresponse_CSC.problem								
Problem Editing Script Functions () Greek Symbols () Other Symbols () Quitout Taos () ()								
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<pre><pre><pre>cycoblem> <web><startouttext></startouttext><h2>Penicillin V</h2><endouttext></endouttext> <script src="/adm/jmol/Jmol.js" type="text/javascript"></script> <script type="text/javascript"> jmolTnitialize("/adm/jmol"); jmolAppletColor(boxbgcolor = #002266",boxfgcolor = "#002266",progresscolor = "#002266"); jmolApplet(400, "load /res/sfu/batchelo/2015tests/jsmol/penicillinv3.mol"); </script> </web> <startouttext></startouttext>Text penicillin V<endouttext></endouttext> <corganicresponse <br="" jmeanswer="24 26 C 5.64 -5.18 C 5.64 -3.78 C 4.42 -3.08 C 3.21 -3.78 C 3.21 -5.18 C 4.42 -5.88 0 6.85
-5.88 C 8.06 -5.18 C 9.27 -5.18 N 10.49 -5.18 C 11.70 -5.88 C 12.06 -7.23 N 13.41 -6.87 C 13.05 -5.52 C 14.81 -6.94 C
15.31 -5.64 S 14.23 -4.76 0 9.27 -7.28 0 11.36 -8.44 C 15.57 -8.12 0 14.94 -9.36 0 16.97 -8.04 C 16.18 -4.54 C 16.70
-5.88 1 2 3 2 3 4 1 4 5 2 5 6 1 6 1 2 1 7 1 7 8 1 8 9 1 9 10 11 10 -1 12 13 113 14 111 14 1 11 12 15 16 1 16
17 1 14 17 -1 13 15 1 9 18 2 12 19 2 15 20 -2 20 21 2 20 21 1 16 23 1 16 24 1" options="autoes">answer="CC3(C)S[CSEN2]2(CEH](NC(=0)COCccccc)]C(=0)N2(CEH]3C(=0)O" 1d="11" molecule="24 26 C 4.91 -4.45 C 3.70 -5.55 C 5.10.97 -6.55 C 5.10.97 -6.55 C 5.10.97 -6.55 C 5.10.97 -6.55 C 5.11.33 -7.90 N 12.69 -7.54 C 12.32 -6.19 0 10.63 -9.12 0 8.55 -7.95 C 14.08 -7.62 C 14.59 -6.31 s 13.50 -5.43 C 19.72 -6.55 C 5.55 -5.55 N 7.6 -5.55 C 10.97 -6.55 C 5.55 -5.55 N 7.6 -5.55 C 10.97 -6.55 C 5.10.97 -6.55 C 5.11.33 -7.90 N 12.69 -7.54 C 12.32 -6.19 0 10.63 -9.12 0 8.55 -7.95 C 14.08 -7.62 C 14.59 -6.31 s 13.50 -5.43 C 15.25 -5.57 -5.50 C 1.52 -5.50</corganicresponse></pre></pre></pre>								

How about 3-D representation of these molecules? JSMOL Need to generate some xyz coordinates of the structure. There are a variety of ways that you might do this and diverse programs and databases. Suffice it to say that a file (MDL mol file or Brookhaven PDB file) can be readily obtained or generated for loading into JSMOL to show an interactive 3D-display. Replaced the original tags for the imported image with appropriate tags for the JSMOL application.

Now shows the entire xml code for the revised problem, shown next...



Video demo:

Still in Author's Construction/Testing environment and want to test out the problem.

Rotate molecule, so that one can visualize the required stereochemistry for the specified answer.

Click Pencil icon to open drawing applet.

Choose appropriate wedge-bonds for the displayed molecule and "Insert Answer" to produce the required SMILES string. Click Submit and see the system's response "Correct".

NEXT: Paper Exams.



Why? Numerous reasons not necessarily individually compelling but collectively so.

Easy to generate new versions of exams, as needed.

Control over dynamically generated content. (show four verisons of one problem with graph).

Numerical answers can also be recorded on custom bubble sheet & machine-graded .

Provides a permanent and cumulative association of assessment resources with statistics reflecting their effectiveness, taking advantage of the analytics within the LON-CAPA system.

We are thus able to easily maintain a semi-quantitative evaluation of every Exam question, which can be reviewed when selecting questions for a new exam.

Prepare seating-plans/room maps showing the distribution of different exam versions to minimize lines of sight.

Lay out collated exams and personalized bubble sheets in advance and seat students according to the design.

Three people could set up this room in 10-15 minutes.

Students can find their exam and seat themselves in ~10 minutes, for a room this size.

Simplifies & streamlines attendance-and-identity checking and grading issues.

Individual Student's Exam Score as displayed in *LON-CAPA*

Midterm Results

Total: Mks/Pts = 21.75/24 Multiple Choice: Mks/Pts = 16/18 Numerical: Mks/Pts = 3/3 Written: Mks/Pts = 2.75/3.00

Exam Version: 2222 Your Choices: bdbbdaaead eeaccbec Exam Answers: bdbbdaaead eeaccbec Point Values: llillillill llillill Q1: Student Ans = 1.14E+15 Exam Ans = 1.14E+15 ± 5.68E+13 Pts = 1 Score = 1 Q2: Student Ans = 9.28E-06 Exam Ans = 9.29E-06 ± 4.64E-07 Pts = 1 Score = 1 Q3: Student Ans = 0.00499 Exam Ans = 0.00499 ± 9.98E-05 Pts = 1 Score = 1

Class Average: 15.07/24

Scoring broken down and comments provided to individual students.

NEXT Analytics for exam questions (or other assessments).

Compiling statistics for 34 problems This will take some time. Gequence Statistics ®												
Sequence	#Items	Score Mean	Score STD	Score Max	Score Min	Score N	Count Mean	Count STD	Count Max	Count Min	Count N	KR-21
Final	36	21.99	5.48	36.00	6.00	337	21.99	5.48	36.00	6.00	337	0.74
Chemis Compiled Final	try 12 on Wea	2 Burnal d May 21 10	by 2014 0:42:29 pr	- 1 n 2014 (P	PDT)							

										1
P#	Title	Part	#Stdnts (plot)	Tries (plot)	tries/correct (plot)	#Wrng (plot)	%Wrng (plot)	DoDiff (plot)	DoDisc (plot)	
1	QvsKEqPos.exam	0	337	337	1.6	125.0	37.0	0.37	0.32	
2	No Buff	0	332	332	6.4	280.0	84.3	0.84	0.11	
3	acidic.or.basic.sol.SASB.exam	0	335	335	1.2	58.0	17.3	0.17	0.27	
4	Ka.rxn.def.exam	0	337	337	1.6	127.0	37.6	0.38	0.42	
5	pH.water.Kw.Tnot25.exam	0	337	337	1.6	120.0	35.6	0.36	0.39	
6	pH.wb.exam	0	337	337	1.4	88.0	26.1	0.26	0.40	
7	me/w.sign.isoT.expans.exam	0	337	337	1.9	158.0	46.8	0.47	0.14	
8	entropy.fusion.exam	0	337	337	1.1	17.0	5.0	0.05	0.08	
9	pH.change.concept.exam	0	336	336	1.2	50.0	14.8	0.15	0.21	
10	acid.str.rel.exam	0	337	337	1.7	133.0	39.4	0.39	0.33	
11	spontaneity.vs.T.exam	0	337	337	1.5	112.0	33.2	0.33	0.55	
12	me/rev.delHf0.mult.exam	0	337	337	2.0	170.0	50.4	0.50	0.54	
13	entropy.spont.exam	0	336	336	2.8	214.0	63.6	0.64	0.12	
14	ME/NO.Gf.K.exam	0	336	336	3.3	235.0	69.9	0.70	0.42	
15	delG0.Kps.thermo.exam	0	328	328	2.5	197.0	60.0	0.60	0.42	
16	std.red.pots.rxn.exam	0	337	337	1.2	55.0	16.3	0.16	0.21	
17	Selective Electrolysis	0	337	337	1.7	143.0	42.4	0.42	0.18	
18	Galvanic.cell.simple.conc.eff.exam	11	337	337	1.2	65.0	19.2	0.19	0.41	
19	Galvanic.cell.simple.conc.eff.exam	13	336	336	1.9	163.0	48.5	0.49	0.44	
20	ME/car.battery.cell.exam	0	332	332	1.5	116.0	34.9	0.35	0.21	

 KR-21 reliability index (for what it's worth). KR21= (N/(N-1))(1- M(N-M)) Ns^2 N=number of items in exam; M=mean score; s=stdev
 Mostly good degrees of discrimination.
 Difficulty level is variant.
 How do these results compare with Historical results for the same questions?

Next slide shows dynamic meta data summary of results for a single re-used question.

Overall Assessment Statist	ical Da	ta				
Statistics calculated for number of students:	2259					
Average number of tries till solved:	1.00					
Degree of difficulty:	ty:					
Degree of discrimination:		(0.40)				
Recent Detailed Assessmer	nt Stati	stical Data	1			
Course		Section(s)	Num Students	Mean Tries	Degree of Difficulty	Degree of Discrimination
Chemistry 122 Burnaby 2011-Spring		D100	454	1.00	0.35	0.39
Chemistry 122 Burnaby 2011-3		D100	235	1.00	0.34	0.46
Chemistry 122 Surrey 2012-Spring		D200	95	1.00	0.22	0.46
Showing 122 Suncy 2012 (Chemistry 122 Burnaby 2012-Spring		332	1.00	0.33	0.37
Chemistry 122 Burnaby 2012			242	1.00	0.26	0.43
Chemistry 122 Burnaby 2012 Chemistry 122 Burnaby 2012	-3	D100				
Chemistry 122 Burnaby 2012 Chemistry 122 Burnaby 2012 Chemistry 122 Burnaby 2013 Chemistry 122 Surrey 2013-5	-3 Spring	D100 D200	125	1.00	0.30	0.47
Chemistry 122 Burnaby 2012 Chemistry 122 Burnaby 2012 Chemistry 122 Burnaby 2013 Chemistry 122 Burnaby 2013	2-3 Spring 1-1	D100 D200 D100	125 307	1.00 1.00	0.30 0.22	0.47 0.26
hemistry 122 Burnaby 2012 hemistry 122 Burnaby 2012 hemistry 122 Burnaby 2013-5 hemistry 122 Burnaby 2013 hemistry 122 Surrey 2014-5	2-3 Spring 1-1 Spring	D100 D200 D100 D200	125 307 132	1.00 1.00 1.00	0.30 0.22 0.17	0.47 0.26 0.26

Example for an individual question used in 6 semesters over 4 years. Stats averaged over all students summarized at top.

Moderate difficulty Good degree of discrimination

Stats listed for 9 semester-courses -- 5 different instructors -- 2 campuses. Note that consistency is reasonable. Though, it appears to suggest that individual instructors get "better" with repeat performances. Overall averages shown to have good predictive value.



"Credits" almost anyone in the Chemistry Department and thanks to ITS and the LON-CAPA programmer-developers participating in the Communities MailLists and Bugzilla site, and attendees at the Annual meetings (held in late-May or early June every year for the past 20).