

SUPPLEMENTARY MATERIAL

Proteomic analysis of plasma proteins of high-flux haemodialysis and on-line haemodiafiltration patients reveals differences in transthyretin levels related with anaemia

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1. High-flux haemodialysis and on-line haemodiafiltration

Haemodialysis was prescribed as four-hour dialysis sessions trice weekly, with a blood flow rate between 220 and 400 ml/min. High-flux haemodialysis was performed using dialysis membranes of a blend of polyarylethersulfone and polyvinylpyrrolidone (Revaclear), and polyamide (Polyflux-210H). On-line haemodiafiltration was performed using polysulfone membranes (Helixone). Other aspects of the patient's treatment prescription did not differ between the two groups. All patients had not residual renal function and were on dialysis for more than 6 months. The treatment times and blood flow rates (individualized for each patient) remained unchanged from their usual prescription. To achieve anticoagulation, heparin was used (firstly a loading dose and then a constant infusion of heparin). Net fluid removal was also individualized. The efficacy of haemodialysis was indexed by the urea kinetic equation single-pool variable volume Kt/V according to Daugirdas [1]. Dialysis vintage years were similar for both groups. Residual renal function was negligible. Kt/V represents the fractional clearance of urea, and it is used as an index of the dialysis dose (K: rate of urea clearance by the dialyzer in ml/min, t: duration of treatment session in min, V: urea volume distribution in the patient in ml). Both groups met the recommendations from the National Kidney Foundation guidelines, Kt/V= 1.2 per haemodialysis session for patients treated thrice weekly [2].

References:

1. Daugirdas JT: Second generation logarithmic estimates of single-pool variable volume Kt/V: an analysis of error. *J Am Soc Nephrol*, 4: 1205-1213, 1993
2. National Kidney Foundation NK: I. NKF-K/DOQI Clinical Practice Guidelines for Haemodialysis Adequacy: update 2000. *Am J Kidney Dis*, 37: S7-S64, 2001

2. Baseline characteristics of the patients

The aetiologies of the ESRD in the patients from both groups were balanced between the groups and representative of the most common causes of ESRD. In the HF group, there were patients suffering from diabetes mellitus (1 patient), glomerulonephritis (2 patient), polycystic kidney disease (1 patient), nephroangiosclerosis and renal ischemia (1 patient), interstitial nephritis (2 patient), multiple myeloma (1 patient), amyloidosis (1 patient) and glomerular hyperfiltration (1 patient). The renal diseases in the HDF group were diabetes mellitus (1 patient), glomerulonephritis (1 patient), polycystic kidney disease (1 patient), nephroangiosclerosis and renal ischemia (1 patient), interstitial nephritis (2 patient), amyloidosis (1 patient), bilateral nephrectomy (1 patient) and unknown (1 patient). In the HDF group there was a patient who had undergone a bilateral nephrectomy due to a renal carcinoma; and a patient whose nephropathy was undiagnosed.

All the patients had a glomerular filtration rate (GFR) < 6 ml/ min (end-stage kidney disease, chronic kidney disease stage 5, GFR < 15 ml/min). There were two patients with diabetes mellitus (requiring the administration of insulin) in each group. The Charlson Comorbidity Index (CCI) was used to assess the base comorbidity of these patients. The CCI considers 17 comorbidities and subcategories for liver disease and diabetes mellitus (Roffman 2016). Patients from both groups received subcutaneous recombinant human erythropoietin (rhEPO) weekly. Though, the requirements of erythropoietin varied significantly between both groups. Patients were also prescribed iron (1-2 vials of 1 g, monthly), folic acid (50 mg) and vitamin D (individualized dosage) supplementation.

3. Preparation of plasma samples

The patient's blood sample was drawn after the long interdialytic period in EDTA-K₃ tubes. The samples were then centrifuged at 900g for 10 minutes and the plasma was separated. Aliquots were stored at -80°C. To study plasma proteome, purification of pooled or individual plasma samples was performed removing high abundant proteins to improve the resolution of gels and detection of low molecular weight proteins using the Pure Proteome™ Albumin/IgG Depletion kit (Millipore), as recommended by the supplier. Plasma sample corresponding to

each patient was stored separately at -80°C , and protein concentrations were determined for each sample.

To check the depletion of these proteins, both the unprocessed and the albumin/IgG-depleted plasma samples were run into SDS-PAGE, and the gel was stained with Coomassie blue for protein detection. Supplemental Figure S1 shows the effective depletion of albumin/IgGs in HF and HDF pooled plasma samples compared to unprocessed HF and HDF pooled plasma samples.

4. 1-D DIGE analysis

4.1. Fluorescence labelling. DIGE assay

Plasma samples from HF and HDF patients were analyzed by one-dimensional fluorescence difference in gel electrophoresis (1-D DIGE). Albumin/IgG-depleted HF and HDF patients' plasma samples (5 μg of protein) were made to 8.5 M urea, centrifuged and labelled on cysteine residues with Cy3 or Cy5 fluorescent dyes (GE Healthcare). After dithiothreitol addition, HF and HDF samples were combined and fluorescence-labelled proteins loaded into standard vertical slab SDS-PAGE for 1-D DIGE. SDS-PAGE was performed in 12% acrylamide (3% cross-linking) gels (1.0 mm thick). The gel was scanned using a Typhoon 9200 imager (GE Healthcare) at 450 V per channel, and the paired Cy5/Cy3 images captured at 100 $\mu\text{m}/\text{pixel}$ resolution for analysis. Cy5 was excited by the laser red-line (633 nm), and Cy3 was excited by the laser green-line (532 nm), sequentially and using appropriate emission filters (Cy5, 670 band-pass 30; Cy3, 580 band-pass 30) to minimize crosstalk. After scanning, the gel was stained with Coomassie blue.

4.2. In-gel protein digestion

Firstly, excised stained gel bands were destaining adding 200 μl of 50% ethanol (vol/vol) in 25 mM ammonium bicarbonate (AB) for 20 min, and after it was removed, added 200 μl of pure ethanol for 15 min, removed it and dried for 5 min. After, in-gel reduction and alkylation of proteins was carried out covering the bands with 200 μl 10 mM dithiothreitol (DTT) in 50 mM AB for 30 min at 56°C , and after DTT was removed, washed with 50 mM AB for 5 min and bands incubated with 55 mM iodoacetamide in 50 mM AB for 30 min in the dark. Then, iodoacetamide was removed and washed twice with 50 and 25 mM AB for 5 min. Gel pieces were dehydrated adding 200 μl of 50% acetonitrile (ACN) (vol/vol) in 25 mM AB for 20 min and removed, it was repeated again, then 200 μl of neat ACN were added for 10 min, and

after ACN removal, gel pieces were dried completely. Then, the gel pieces were saturated with trypsin adding 15 μ l of a solution containing 40 ng of trypsin (1:20 w/w) in 25 mM AB and leaving them in an ice bucket or a fridge for 45 min. After removing liquid excess, 20 μ l of 25 mM AB was added and incubated overnight at 30°C. Finally, to extract the peptide digestion products, 1 μ l of neat ACN were added (final concentration 6-15% ACN), incubated at 37°C for 15 min, and then 0.2 μ l of trifluoroacetic (TFA) was added to 0.1% (vol/vol) final concentration and incubated for 10 min at room temperature, collecting the supernatant for mass spectrometry analysis.

4.3. Protein identification by mass spectrometry

After digestion, peptide mass fingerprinting (PMF) was analyzed by matrix-assisted laser desorption/ionization time-of-flight mass spectrometry (MALDI-TOF MS) (Autoflex III TOF/TOF, Bruker-Daltonics) for protein identification, as previously described [3]. PMF spectra were obtained using an alpha-cyanocinnamic acid matrix and calibrated with a peptide calibration standard mixture (222570, Bruker Daltonics) contained nine standard peptides in a molecular mass range between \sim 700 and 3500 Da to reach a typical mass measurement accuracy $< \pm 3$ ppm. Additionally, when available and for confirmation of protein identity, peptide fragmentation was performed by MS in tandem MALDI LIFT-TOF/TOF [4]. MS data of PMFs and MS/MS data from LIFT TOF/TOF spectra were searched in the SwissProt database using the MASCOT database search algorithm (Matrix Science) for protein identification. No more than one missed tryptic cleavage was allowed and a mass accuracy of 50 ppm was used for all mass searches.

References:

3. Cid C, Garcia-Bonilla L, Camafeita E, Burda J, Salinas M, Alcazar A: Proteomic characterization of protein phosphatase 1 complexes in ischemia-reperfusion and ischemic tolerance. *Proteomics*, 7: 3207-3218, 2007
4. Suckau D, Resemann A, Schuerenberg M, Hufnagel P, Franzen J, Holle A: A novel MALDI LIFT-TOF/TOF mass spectrometer for proteomics. *Anal Bioanal Chem*, 376: 952-965, 2003

5. LC-MS/MS Analysis

5.1 Protein digestion

Plasma samples were boiled for 5 min at 100 °C in the presence of 25 mM Tris-HCl pH 6.8, 1% SDS and 50 mM DTT. This mixture (6 μ l) was diluted with denaturing buffer (8 M urea

in 100 mM Tris-HCl pH 8.5) and concentrated on FASP filters (Expedeon). After washing with denaturing buffer at 10,000 rpm for 15 min, free thiol groups were alkylated by incubation with 50 mM iodoacetamide 30 min at room temperature in the dark. Then the filters were washed with denaturing buffer followed by washing with trypsin digestion buffer (50 mM ammonium bicarbonate pH 8.8). Protein samples were digested overnight at 37°C with sequencing grade trypsin (Promega, Madison, WI, USA) at 1:40 (w/w) trypsin:protein ratio in digestion buffer. The resulting tryptic peptides from each sample were recovered by centrifugation at 10,000 rpm for 5 min after addition of 40 µl of trypsin digestion buffer, after which 50 µl of 500 mM NaCl were added and the filters centrifuged for 15 min at 10,000 rpm. Trifluoroacetic acid was added to a final concentration of 1% and the peptides were desalted on C18 Oasis HLB extraction cartridges (Waters Corporation, Milford, MA, USA) and dried-down.

5.2. Isobaric labelling of peptides using tandem mass tags (TMT)

The dried peptides were dissolved in triethylammonium bicarbonate buffer, and their concentration was determined using a Direct Detect IR spectrometer (Millipore). Equal amounts of each peptide sample were labeled with isobaric 10-plex Tandem Mass Tags (TMT, Thermo Scientific) according to the manufacturer's instructions and mixed together. An internal control was prepared by pooling the nine Control samples and was used as a reference to express relative quantification values. Samples were desalted using C18 Oasis HLB extraction cartridges (Waters) and dried-down for later LC-MS/MS analysis.

5.3. Peptide identification

Peptide samples were taken up in 0.1% formic acid and applied to an EASY-nLC 1000 nano-flow HPLC system (Thermo Fisher Scientific) coupled on-line with an orbitrap Fusion mass spectrometer (Thermo Fisher Scientific). C18-based reverse phase separation was used with a 2-cm trap column and a 50-cm analytical column (EASY-Spray, Thermo Fisher Scientific). Peptides were loaded in buffer A (0.1% formic acid (v/v)) and eluted with a 300-min linear gradient of buffer B (90% ACN, 0.1% formic acid (v/v)) at 200 nL/min flow. Mass spectra were acquired in a data-dependent manner, with an automatic switch between MS and MS/MS, using a top-speed method and 30 s dynamic exclusion. MS spectra were acquired in the 400–1500 m/z range at 120,000 resolution, while HCD MS/MS were performed at 33 normalized collision energy and analyzed with 35,000 resolution in the orbitrap.

LC-MS/MS data were analyzed with Proteome Discoverer (version 2.1, Thermo Fisher Scientific) using SEQUEST-HT (Thermo Fisher Scientific) against a Uniprot database containing all sequences from Homo sapiens (February 2020; 148,000 entries). Database search parameters were selected as follows: trypsin digestion with two maximum missed cleavage sites, precursor mass tolerance of 800 ppm, fragment mass tolerance of 20 mmu. Met oxidation was considered a variable modification, while Cys carbamidomethylation and Lys and peptide N-terminal modification of +229.163 were set as fixed modifications for TMT. The corresponding inverted protein sequences were incorporated to the database for false discovery rate (FDR) calculation. Peptide identification from MS/MS data was performed using the probability ratio method [5], and the FDR of peptide identifications was calculated using the refined method [6,7], taking 1% FDR as a threshold for peptide identification. Peptides were assigned only to the best protein proposed by the Proteome Discoverer algorithm.

References:

5. Martinez-Bartolome, S. *et al.* Properties of average score distributions of SEQUEST: the probability ratio method. *Mol Cell Proteomics* **7**, 1135-1145 (2008).
6. Navarro, P. & Vazquez, J. A refined method to calculate false discovery rates for peptide identification using decoy databases. *J Proteome Res* **8**, 1792-1796 (2009).
7. Bonzon-Kulichenko, E., Garcia-Marques, F., Trevisan-Herraz, M. & Vazquez, J. Revisiting peptide identification by high-accuracy mass spectrometry: problems associated with the use of narrow mass precursor windows. *J Proteome Res* **14**, 700-710 (2015).

5.4. Quantification at the peptide and protein levels

The quantitative information extracted from the MS/MS spectra by Proteome Discoverer was integrated from the spectrum level to the peptide level and then to the protein level on the basis of the WSPP model [8] and the systems biology triangle algorithm [9] using the SanXoT software package [10]. Briefly, the log₂-ratio of every scan was calculated using the TMT reporter ion intensities coming from samples A and B. The log₂-ratio of every peptide was then calculated as the weighted average of its scans, whereas the quantification of each protein was the weighted average of its peptides, and averaged of all the protein values [8]. Peptide quantification weights were calculated from the corresponding scan weights and the peptide variance, and protein quantification weights were calculated from the corresponding peptide weights and the protein variance. Thus, protein quantification was defined by a Zq

value as a normalized log₂-ratio at the protein level expressed in standard deviation units [8]. Outliers at the peptide and protein levels were detected at 1% FDR, and a FDR_q value < 0.05 was defined as significant [8].

References:

8. Navarro, P. *et al.* General statistical framework for quantitative proteomics by stable isotope labeling. *J Proteome Res* **13**, 1234-1247 (2014).
9. Garcia-Marques, F. *et al.* A Novel Systems-Biology Algorithm for the Analysis of Coordinated Protein Responses Using Quantitative Proteomics. *Mol Cell Proteomics* **15**, 1740-1760 (2016).
10. Trevisan-Herraz, M. *et al.* SanXoT: a modular and versatile package for the quantitative analysis of high-throughput proteomics experiments. *Bioinformatics* **35**, 1594-1596 (2019).

6. Quantification of TTR peptides by mass spectrometry. Calibration experiments

Calibration experiments were previously carried out to test the accuracy of the quantification using known concentrations of reference peptides bradykinin (765.85 Da), bombesin (1619.85 Da) and adrenocorticotrophic hormone (ACTH, 2465.67 Da), in addition to angiotensin-II peptide (1045.54 Da). Peptide relative intensities were obtained from MS spectra in the same conditions as for peptide mass fingerprinting analysis by MALDI-TOF MS (see above). Reference peptides were quantified with respect to the angiotensin-II peptide as a ratio of the relative intensities of peptide/angiotensin-II. The reference peptide concentrations used were 2-100 fmol, with 4 fmol of angiotensin-II, and 4-400 fmol, with 40 fmol of angiotensin-II.

To quantify transthyretin (TTR) peptides, the angiotensin-II peptide was added as an internal standard to obtain a relative intensity of the TTR peptides. The accuracy of MS peptide quantification was confirmed with calibration experiments carried out with increasing concentrations of reference peptides (molecular mass range from ~ 765 to 2465 Da) (Figure S4). The calibration experiments showed a strict correlation between peptide concentration and relative intensity for each peptide (Figure S4). TTR peptides with 1366.75, 1394.62, 1416.77, 1522.71, 2451.20 and 3140.51 Da were identified by MASCOT database search algorithm (Matrix Science, <http://www.matrixscience.com>), in addition to 833.40 Da TTR peptide, identified by FindPept tool (ExpASY, <https://web.expasy.org/findpept/>) (Figure 4A). Accordingly, TTR peptides from MALDI-TOF MS spectra were quantified in the mass range of the calibration experiments (TTR peptides from 833.40 to 2451.20 Da) as relative intensities (ratio TTR peptide intensity/angiotensin-II intensity).

7. Detection of the allele distribution of haptoglobin α in HF and HDF patients

Haptoglobin is a glycoprotein with α and β chains. In human, haptoglobin exists in two allelic forms designated Haptoglobin 1 (*Hp1*) and Haptoglobin 2 (*Hp2*) that results in three known phenotypes, Hp1-1, Hp2-2 and the heterozygous phenotype Hp2-1. *Hp2* contain the allele *Hpa2* that is the product of a partial *Hpa1* gene duplication [11]. Thus, the product of *Hpa2* gives a protein with ~ 7 kDa greater than that of *Hpa1*.

In our experiments, haptoglobin $\alpha 2$ was detected as proteins 20 and 21 (Figure 3A), showing similar intensity values in both HF and HDF patient samples (Figure 3B). In these experiments, haptoglobin $\alpha 1$ was significantly detected with higher levels in HF patients than in HDF patients (Figure 1B, protein *f*, and Figure 3B). However, the allele distribution between HF and HDF groups should be tested. To these end, we analysed the expression of haptoglobin $\alpha 2$ and $\alpha 1$ in HF and HDF samples by western blot using a specific anti-haptoglobin α antibody. The results showed two proteins identified by the antibody, which according to the molecular mass standards used, would correspond to haptoglobin $\alpha 2$ and $\alpha 1$ (Figure S3). These results also showed that three HF patients (patients 3, 6, and 7) had not haptoglobin $\alpha 2$ expression –patient HF 9 showed a very slight expression of haptoglobin $\alpha 2$ – (Figure S3), having therefore the phenotype Hp1-1. On the other hand, only one HDF patient (patient 6) had not haptoglobin $\alpha 2$, being of the Hp1-1 phenotype. As expected, the phenotype Hp1-1 induces more expression of haptoglobin $\alpha 1$ than the phenotypes Hp2-1 and Hp2-2 (Figure S3, see patients HF 3, 6 and 7 and HDF 6), and therefore, the presence of more subjects with phenotype Hp1-1 among HF patients can be the cause of the higher levels of haptoglobin $\alpha 1$ found in this group.

References:

11. Yang, F., Brune, J. L., Baldwin, W. D., Barnett, D. R. & Bowman, B. H. Identification and characterization of human haptoglobin cDNA. *Proc Natl Acad Sci U S A* **80**, 5875-5879 (1983).

8. Supplementary Table S1

Table S1. Proteins identified by MALDI-TOF MS in plasma samples of HF and HDF patients.

N°	Protein	Accession no. ^a	Gene name	Theoretical mass (Da)	Score ^b	Peptides matched/ searched	% Coverage	Lift ^c (score)
1	Alpha-2-macroglobulin	P01023	<i>A2M</i>	164613	95	24/86	20	1148.62(63)
2	Ceruloplasmin	P00450	<i>CP</i>	122983	112	19/66	24	2155.05(80)
3	Inter-alpha-trypsin inhibitor heavy chain H4	Q14624	<i>ITIH4</i>	103521	48	6/26	10	2184.07(50)
7	Serotransferrin	P02787	<i>TF</i>	79294	152	25/95	33	1881.91(91)
12	Alpha-1-antitrypsin	P01009	<i>SERPINA1</i>	46737	98	10/41	27	1641.83(99)
13	Apolipoprotein A-IV	P06727	<i>APOA4</i>	45371	127	17/72	37	1352.69(76)
14	Haptoglobin β	P00738	<i>HP</i>	27265	109	14/85	33	1707.87(75)
15	Haptoglobin β	P00738	<i>HP</i>	27265	123	13/29	33	
18	Apolipoprotein A-I	P02647	<i>APOA1</i>	30759	191	23/98	71	
19	Retinol-binding protein 4	P02753	<i>RBP4</i>	23337	70	5/13	33	2693.10(65)
20	Haptoglobin α 2	P00738	<i>HP</i>	15945	62	8/52	21	
21	Haptoglobin α 2	P00738	<i>HP</i>	15945	60	6/38	16	3432.62(70)
22	Transthyretin	P02766	<i>TTR</i>	13761	64	6/65	61	2451.20(189)
23	Haptoglobin α1	P00738	<i>HP</i>	9192	64	10/81	19	1708.91(91)
24	Serum amyloid A1	P0DJ18	<i>SAA1</i>	13581	63	5/39	55	

^a, Accession number in UniProt database (<https://www.uniprot.org>). ^b, Protein identification scores > 56 were significant ($p < 0.05$) in the MASCOT database search algorithm. ^c, MALDI LIFT-TOF/TOF MS identification mode; the m/z of the fragmented parental peptide is indicated; MASCOT scores (in parenthesis) > 28 were significant ($p < 0.05$). Proteins marked in bold were significant by DIGE quantification (Figure 1; $p < 0.05$, t-test) with higher levels in plasma samples of HF compared with HDF patients.

9. Supplementary Table S2

Table S2. Proteins quantified in LC-MS/MS analysis in plasma samples of HDF compared with HF patients.

Protein	Accession no. ^a	Gene name	Theoretical mass (Da)	Peptides identified	Zq value ^b	Zq value (colour code) ^c
cDNA FLJ53691, highly similar to Serotransferrin	B4E1B2		74832	4	3.04 *#	
Trypsin-1	P07477	<i>PRSS1</i>	26558	2	2.42 *	
Immunoglobulin heavy constant alpha 1	P01876	<i>IGHA1</i>	37655	14	2.09 *	
Dystroglycan	Q14118	<i>DAG1</i>	97441	2	1.77	
Immunoglobulin heavy variable 4-28	A0A0C4DH34	<i>IGHV4-28</i>	13124	3	1.57	
Immunoglobulin delta heavy chain	P0DOX3		56224	2	1.50	
Immunoglobulin heavy constant gamma 3	P01860	<i>IGHG3</i>	41287	6	1.43	
Immunoglobulin heavy variable 3-7	P01780	<i>IGHV3-7</i>	12943	2	1.42	
Rheumatoid factor G9 light chain (Fragment)	A0N5G3	<i>V-lambda-3</i>	12897	2	1.32	
IBM-A3 heavy chain variable region (Fragment)	A0A0X9UWM4		14369	2	1.28	
Uncharacterized protein DKFZp686K04218 (Fragment)	Q7Z379	<i>DKFZp686K04218</i>	51620	2	1.23	
Vitronectin	D9ZGG2	<i>VTN</i>	54306	5	1.18	
Immunoglobulin heavy constant gamma 4	P01861	<i>IGHG4</i>	35941	7	1.11	
Ig heavy chain variable region (Fragment)	A0A068LKQ8		12449	2	1.05	
IBM-B3 light chain variable region (Fragment)	A0A109PW74		11399	2	1.03	
Immunoglobulin kappa constant	P01834	<i>IGKC</i>	11765	8	1.01	
V1-22 protein (Fragment)	Q5NV88	<i>V1-22</i>	10643	2	1.00	
Alpha-2-macroglobulin	P01023	<i>A2M</i>	163291	70	0.94	
Immunoglobulin heavy constant gamma 1	P01857	<i>IGHG1</i>	36106	17	0.93	
Properdin	P27918	<i>CFP</i>	51276	3	0.87	
B cell receptor heavy chain variable region (Fragment)	A0A1C9J6R2		14524	2	0.87	
Immunoglobulin lambda-like polypeptide 5	B9A064	<i>IGLL5</i>	23063	4	0.87	
Immunoglobulin lambda constant 2	P0DOY2	<i>IGLC2</i>	11294	2	0.85	
Cartilage oligomeric matrix protein	P49747	<i>COMP</i>	82860	4	0.79	
Immunoglobulin lambda constant 7	A0M8Q6	<i>IGLC7</i>	11254	4	0.72	
Immunoglobulin heavy variable 3-49	A0A0A0MS15	<i>IGHV3-49</i>	13056	3	0.69	
Immunoglobulin heavy constant alpha 2	P01877	<i>IGHA2</i>	36591	2	0.67	
Serum amyloid A protein	D3DQX7	<i>SAA1</i>	13562	2	0.65	
Carboxypeptidase N subunit 2	P22792	<i>CPN2</i>	60557	9	0.62	
Immunoglobulin lambda-like polypeptide 1	P15814	<i>IGLL1</i>	22963	4	0.61	
MS-F1 light chain variable region (Fragment)	A0A0X9V9B3		12108	2	0.60	
Immunoglobulin kappa variable 3D-20	A0A0C4DH25	<i>IGKV3D-20</i>	12515	2	0.59	
Beta-Ala-His dipeptidase	Q96KN2	<i>CNDP1</i>	56706	7	0.56	
Immunoglobulin alpha-2 heavy chain	P0DOX2		48934	4	0.54	
MS-F1 heavy chain variable region (Fragment)	A0A125U0V1		13509	2	0.51	
Immunoglobulin kappa variable 4-1	P06312	<i>IGKV4-1</i>	13380	2	0.51	
Ig heavy chain variable region (Fragment)	A0A2U8J8Y4	<i>IgH</i>	12010	2	0.46	
Pregnancy zone protein	P20742	<i>PZP</i>	163863	16	0.44	
IBM-A1 light chain variable region (Fragment)	A0A0X9USN3		11626	2	0.44	
Myosin-reactive immunoglobulin light chain variable region (Fragment)	Q9UL83		11834	2	0.42	
Immunoglobulin kappa variable 3D-7	A0A0C4DH55	<i>IGKV3D-7</i>	13148	2	0.42	
IBM-B2 heavy chain variable region (Fragment)	A0A125QYY9		13569	2	0.39	
Fibulin-1	P23142	<i>FBLN1</i>	77214	9	0.38	
Afamin	P43652	<i>AFM</i>	69069	17	0.38	
Carboxypeptidase N catalytic chain	P15169	<i>CPN1</i>	52286	4	0.37	
Biotinidase	P43251	<i>BTD</i>	61133	4	0.36	
IgGfc-binding protein	Q9Y6R7	<i>FCGBP</i>	572017	2	0.33	
Vitamin K-dependent protein S	P07225	<i>PROS1</i>	75123	18	0.27	
Thyroxine-binding globulin	P05543	<i>SERPINA7</i>	46325	14	0.26	
Immunoglobulin kappa variable 3-7 (non-functional) (Fragment)	A0A075B6H7	<i>IGKV3-7</i>	12783	2	0.24	

Ig heavy chain variable region (Fragment)	A0A2U8J8Z6	<i>IgH</i>	10905	2	0.22
Fibronectin	P02751	<i>FN1</i>	272320	54	0.22
Apolipoprotein(a)	P08519	<i>LPA</i>	501319	10	0.21
Complement factor B	P00751	<i>CFB</i>	85533	18	0.21
Immunoglobulin kappa variable 1-6	A0A0C4DH72	<i>IGKV1-6</i>	12697	2	0.19
Coagulation factor IX	P00740	<i>F9</i>	51778	5	0.18
CP protein	A5PL27	<i>CP</i>	122205	33	0.15
Carbonic anhydrase 1 (Fragment)	E5RFE7	<i>CA1</i>	21402	3	0.13
Apolipoprotein C-I	P02654	<i>APOC1</i>	9332	5	0.12
Hyaluronan-binding protein 2	Q14520	<i>HABP2</i>	62672	10	0.12
Immunoglobulin kappa variable 3D-15	A0A087WSY6	<i>IGKV3D-15</i>	12534	3	0.11
von Willebrand factor	P04275	<i>VWF</i>	309265	16	0.11
Gelsolin	A0A0A0MS51	<i>GSN</i>	82526	14	0.11
Immunoglobulin J chain	P01591	<i>JCHAIN</i>	18099	4	0.09
Hepatocyte growth factor activator	Q04756	<i>HGFAC</i>	70682	7	0.09
Complement C3	P01024	<i>C3</i>	187148	100	0.09
SPARC-like protein 1	Q14515	<i>SPARCL1</i>	75208	2	0.09
Complement component C8 gamma chain	P07360	<i>C8G</i>	22277	7	0.07
IGK@ protein	Q6PIL8	<i>IGK@</i>	25834	2	0.05
Selenoprotein P	P49908	<i>SELENOP</i>	43174	2	0.05
Prothrombin	P00734	<i>F2</i>	70037	31	0.05
Apolipoprotein F	Q13790	<i>APOF</i>	35399	3	0.05
cDNA FLJ51266, highly similar to Vitronectin	B7Z553		23605	3	0.05
Angiotensinogen (Fragment)	A7L3A3	<i>AGT</i>	9407	4	0.04
N-acetylmuramoyl-L-alanine amidase	Q96PD5	<i>PGLYRP2</i>	62217	8	0.02
Immunoglobulin heavy variable 3-9	P01782	<i>IGHV3-9</i>	12945	3	0.01
Immunoglobulin heavy constant delta	P01880	<i>IGHD</i>	42353	4	0.00
Immunoglobulin heavy variable 3-15	A0A0B4J1V0	<i>IGHV3-15</i>	12926	2	-0.01
Cystatin-C	P01034	<i>CST3</i>	15799	6	-0.02
Immunoglobulin lambda variable 3-10	A0A075B6K4	<i>IGLV3-10</i>	12441	2	-0.02
Myosin-reactive immunoglobulin kappa chain variable region (Fragment)	Q9UL86		11928	3	-0.05
MS-C1 heavy chain variable region (Fragment)	A0A125U0U7		13097	2	-0.05
Immunoglobulin heavy constant gamma 2	P01859	<i>IGHG2</i>	35901	13	-0.06
Transforming growth factor-beta-induced protein ig-h3	Q15582	<i>TGFBI</i>	74681	3	-0.06
Angiotensinogen (Serpine peptidase inhibitor, clade A, member 8)	B0ZBE2	<i>AGT</i>	53154	5	-0.07
EGF containing fibulin-like extracellular matrix protein 1 isoform 2	A0A0S2Z3V1	<i>EFEMP1</i>	45723	5	-0.07
Coagulation factor X	P00742	<i>F10</i>	54732	5	-0.07
C4b-binding protein alpha chain	P04003	<i>C4BPA</i>	67033	26	-0.08
Peroxiredoxin-2	P32119	<i>PRDX2</i>	21892	4	-0.09
Ig heavy chain variable region (Fragment)	A0A2U8J9B3	<i>IgH</i>	12311	2	-0.10
Ig heavy chain variable region (Fragment)	A0A2U8J8R6	<i>IgH</i>	10922	2	-0.14
Serum albumin	P02768	<i>ALB</i>	69367	76	-0.16
Plasma kallikrein	P03952	<i>KLKB1</i>	71370	12	-0.16
Pigment epithelium-derived factor	P36955	<i>SERPINF1</i>	46312	11	-0.17
Sulfhydryl oxidase 1	O00391	<i>QSOX1</i>	82578	2	-0.18
Apolipoprotein C-IV	P55056	<i>APOC4</i>	14553	2	-0.18
Ig heavy chain variable region (Fragment)	A0A2U8J8Q6	<i>IgH</i>	10379	2	-0.19
Fibrinogen alpha chain	P02671	<i>FGA</i>	94973	41	-0.20
Platelet glycoprotein Ib alpha chain	P07359	<i>GP1BA</i>	71540	2	-0.20
Coagulation factor V	P12259	<i>F5</i>	251703	12	-0.21
Chromogranin-A	P10645	<i>CHGA</i>	50688	9	-0.22
Adiponectin A	A0A024RAB9	<i>C1QB</i>	26722	5	-0.22
Immunoglobulin heavy variable 3-21	A0A0B4J1V1	<i>IGHV3-21</i>	12840	3	-0.23
Apolipoprotein D	P05090	<i>APOD</i>	21276	7	-0.23
Fetuin-B	Q9UGM5	<i>FETUB</i>	42055	5	-0.24
Immunoglobulin heavy variable 4-4	A0A075B6R2	<i>IGHV4-4</i>	12848	4	-0.25
Inter-alpha-trypsin inhibitor heavy chain H4	Q14624	<i>ITIHA</i>	103357	26	-0.26
Complement factor H	P08603	<i>CFH</i>	139096	54	-0.26
Heavy chain of factor I (Fragment)	Q6LAM1		35883	2	-0.26
Immunoglobulin heavy variable 3-72	A0A0B4J1Y9	<i>IGHV3-72</i>	13203	3	-0.28
Immunoglobulin kappa variable 3D-11	A0A0A0MRZ8	<i>IGKV3D-11</i>	12625	2	-0.28
C-reactive protein	P02741	<i>CRP</i>	25039	6	-0.29
Monocyte differentiation antigen CD14	P08571	<i>CD14</i>	40076	4	-0.29
Actin, cytoplasmic 1	P60709	<i>ACTB</i>	41737	5	-0.30

Apolipoprotein B-100	P04114	<i>APOB</i>	515605	178	-0.31
Complement factor D	P00746	<i>CFD</i>	27033	9	-0.31
Complement component C8 alpha chain	P07357	<i>C8A</i>	65163	15	-0.31
Complement component C6	P13671	<i>C6</i>	104786	17	-0.31
Lipopolysaccharide-binding protein	P18428	<i>LBP</i>	53384	5	-0.33
Glutathione peroxidase 3	P22352	<i>GPX3</i>	25552	5	-0.33
Fibrinogen beta chain	P02675	<i>FGB</i>	55928	43	-0.34
Complement factor I	P05156	<i>CFI</i>	65750	17	-0.34
Inter-alpha-trypsin inhibitor heavy chain H1	P19827	<i>ITIH1</i>	101389	18	-0.34
Apolipoprotein A-I	P02647	<i>APOA1</i>	30778	38	-0.35
Inter-alpha-trypsin inhibitor heavy chain H2	P19823	<i>ITIH2</i>	106463	21	-0.37
Alpha-1B-glycoprotein	P04217	<i>A1BG</i>	54254	10	-0.37
Immunoglobulin kappa light chain	P0DOX7		23379	4	-0.38
Complement component C8 beta chain	P07358	<i>C8B</i>	67047	13	-0.38
Serotransferrin	P02787	<i>TF</i>	77064	56	-0.38
Complement C5	P01031	<i>C5</i>	188305	38	-0.39
Fibrinogen gamma chain	P02679	<i>FGG</i>	51512	31	-0.39
GCT-A1 light chain variable region (Fragment)	A0A0X9TDD0	<i>2 SV</i>	12436	2	-0.39
Alpha-2-HS-glycoprotein	P02765	<i>AHSG</i>	39341	10	-0.41
Cystatin-M	Q15828	<i>CST6</i>	16511	2	-0.42
Platelet basic protein	P02775	<i>PPBP</i>	13894	2	-0.43
Fibronectin 1, isoform CRA_n	A0A024R462	<i>FN1</i>	259211	2	-0.43
Immunoglobulin heavy variable 3-74	A0A0B4J1X5	<i>IGHV3-74</i>	12840	4	-0.43
Plasma serine protease inhibitor	P05154	<i>SERPINA5</i>	45675	5	-0.44
Vasorin	Q6EMK4	<i>VASN</i>	71713	2	-0.46
Mannan-binding lectin serine protease 1	P48740	<i>MASP1</i>	79247	5	-0.47
Carbonic anhydrase 1 (Fragment)	E5RFL2	<i>CA1</i>	13256	2	-0.48
Lectin galactoside-binding soluble 3 binding protein isoform 1	A0A0S2Z3Y1	<i>LGALS3BP</i>	65331	7	-0.49
Heparin cofactor 2	P05546	<i>SERPIND1</i>	57071	15	-0.49
Kininogen-1	P01042	<i>KNG1</i>	71957	28	-0.52
Prenylcysteine oxidase 1	Q9UHG3	<i>PCYOX1</i>	56640	3	-0.54
Basement membrane-specific heparan sulfate proteoglycan core protein	P98160	<i>HSPG2</i>	468830	2	-0.57
Mutant hemoglobin alpha 2 globin chain	A0A0K2BMD8	<i>HBA2</i>	15258	6	-0.60
Adiponectin B	A0A024RAA7	<i>C1QC</i>	25774	2	-0.60
IBM-B1 light chain variable region (Fragment)	A0A125QYY8		12167	2	-0.63
Complement factor H-related protein 1	Q03591	<i>CFHR1</i>	37651	3	-0.63
Inter-alpha-trypsin inhibitor heavy chain H3	Q06033	<i>ITIH3</i>	99849	15	-0.63
Clusterin	P10909	<i>CLU</i>	52495	18	-0.65
Complement C1s subcomponent	P09871	<i>C1S</i>	76684	16	-0.66
Coagulation factor XIII B chain	P05160	<i>F13B</i>	75511	15	-0.67
Vitamin K-dependent protein C	P04070	<i>PROC</i>	52071	3	-0.69
N90-VRC38.01 light chain variable region (Fragment)	A0A1W6IYJ9		12352	4	-0.69
Plasminogen	P00747	<i>PLG</i>	90569	32	-0.69
Serum amyloid P-component	P02743	<i>APCS</i>	25387	7	-0.69
C4b-binding protein beta chain	P20851	<i>C4BPB</i>	28357	6	-0.71
Alpha-1-antitrypsin	P01009	<i>SERPINA1</i>	46737	38	-0.71
Testicular tissue protein Li 61	A0A140VJI7		60674	7	-0.72
Epididymis secretory sperm binding protein	A0A384N669		38429	10	-0.72
Carboxypeptidase B2	Q961Y4	<i>CPB2</i>	48424	7	-0.72
Coagulation factor XIII A chain	P00488	<i>F13A1</i>	83267	4	-0.74
Kallistatin	P29622	<i>SERPINA4</i>	48542	12	-0.74
Insulin-like growth factor II	P01344	<i>IGF2</i>	20140	2	-0.76
Alpha-2-antiplasmin	P08697	<i>SERPINF2</i>	54566	12	-0.77
Corticosteroid-binding globulin	P08185	<i>SERPINA6</i>	45141	7	-0.78
Immunoglobulin kappa variable 1-27	A0A075B6S5	<i>IGKV1-27</i>	12712	2	-0.79
Phosphatidylinositol-glycan-specific phospholipase D	P80108	<i>GPLD1</i>	92336	7	-0.82
Vitamin D-binding protein	P02774	<i>GC</i>	52918	29	-0.84
Coagulation factor XII	P00748	<i>F12</i>	67792	8	-0.84
Complement C2	P06681	<i>C2</i>	83268	16	-0.85
Hemoglobin subunit delta	P02042	<i>HBD</i>	16055	9	-0.86
Apolipoprotein L1	O14791	<i>APOL1</i>	43974	7	-0.86
Phospholipid transfer protein	P55058	<i>PLTP</i>	54739	3	-0.86
Hemoglobin subunit beta	P68871	<i>HBB</i>	15998	8	-0.86
Vitamin D binding protein (Fragment)	A0A1B1CYC5	<i>Gc</i>	3718	2	-0.86
Insulin-like growth factor binding protein 4, isoform	A0A024R1U8	<i>IGFBP4</i>	27934	2	-0.86

CRA_a					
Serpin peptidase inhibitor clade G member 1 isoform 4	A0A0S2Z333	<i>SERPING1</i>	19647	10	-0.87
Serum paraoxonase/arylesterase 1	P27169	<i>PON1</i>	39731	7	-0.88
HCG40889, isoform CRA_b	A0A024R962	<i>hCG_40889</i>	139070	2	-0.88
Mannan-binding lectin serine protease 2	O00187	<i>MASP2</i>	75702	2	-0.88
Protein AMBP	P02760	<i>AMBP</i>	38999	18	-0.90
Immunoglobulin lambda variable 2-8	P01709	<i>IGLV2-8</i>	12382	2	-0.95
Histidine-rich glycoprotein	P04196	<i>HRG</i>	59578	12	-0.96
Retinol-binding protein 4	P02753	<i>RBP4</i>	23010	8	-0.96
Apolipoprotein C-II isoform 1	A0A024R0T9	<i>APOC4-APOC2</i>	11284	4	-0.97
Complement C1r subcomponent	P00736	<i>C1R</i>	80119	12	-0.98
Serum albumin	Q56G89		69084	2	-0.99
Apolipoprotein M	O95445	<i>APOM</i>	21253	4	-1.01
Complement component C7	P10643	<i>C7</i>	93518	17	-1.04
Serpin peptidase inhibitor, clade A (Alpha-1 antiprotease, antitrypsin) member 3, isoform CRA_c	A0A024R6P0	<i>SERPINA3</i>	47651	22	-1.12
Hepatocyte growth factor-like protein	P26927	<i>MST1</i>	80320	5	-1.12
Attractin	O75882	<i>ATRN</i>	158537	10	-1.13
Serpin peptidase inhibitor clade G member 1	A0A348GSH7	<i>SERPING1</i>	49757	8	-1.15
Serum amyloid A-1 protein	P0DJ18	<i>SAA1</i>	13532	7	-1.17
Apolipoprotein A-II	P02652	<i>APOA2</i>	11175	7	-1.17
Complement C1r subcomponent-like protein	Q9NZP8	<i>C1RL</i>	53498	2	-1.20
Protein Z-dependent protease inhibitor	Q9UK55	<i>SERPINA10</i>	50707	2	-1.21
Alpha-1-acid glycoprotein 2	P19652	<i>ORM2</i>	23603	9	-1.22
Antithrombin-III	P01008	<i>SERPINC1</i>	52602	29	-1.22
Prostaglandin-H2 D-isomerase	P41222	<i>PTGDS</i>	21029	5	-1.24
Beta-2-microglobulin	P61769	<i>B2M</i>	13715	4	-1.27
Complement C4-A	P0C0L4	<i>C4A</i>	192785	65	-1.30
Complement factor H-related protein 2	P36980	<i>CFHR2</i>	30651	4	-1.36
Protein S100-A9	P06702	<i>S100A9</i>	13242	6	-1.38
Complement C4-B	P0C0L5	<i>C4B</i>	192751	4	-1.42
Insulin-like growth factor-binding protein 3	P17936	<i>IGFBP3</i>	31674	4	-1.44
Complement component C9	P02748	<i>C9</i>	63173	16	-1.46
Hemopexin	P02790	<i>HPX</i>	51676	24	-1.47
Immunoglobulin kappa variable 3-20	P01619	<i>IGKV3-20</i>	12557	2	-1.51
CD44 antigen	P16070	<i>CD44</i>	81538	2	-1.56
Immunoglobulin heavy variable 5-51	A0A0C4DH38	<i>IGHV5-51</i>	12675	4	-1.58
Alpha-1-acid glycoprotein	V9HWF6	<i>HEL-S-153w</i>	23512	6	-1.61
Serum amyloid A-4 protein	P35542	<i>SAA4</i>	14747	4	-1.62
Insulin-like growth factor-binding protein complex acid labile subunit	P35858	<i>IGFALS</i>	66035	10	-1.63
Insulin-like growth factor-binding protein 2	P18065	<i>IGFBP2</i>	34814	4	-1.63
CD5 antigen-like	O43866	<i>CD5L</i>	38088	11	-1.65
Haptoglobin	P00738	<i>HP</i>	45205	30	-1.71
Ficolin-3	O75636	<i>FCN3</i>	32903	5	-1.74
Apolipoprotein A-IV	P06727	<i>APOA4</i>	45372	33	-1.81
Apolipoprotein H (Beta-2-glycoprotein I), isoform CRA_a	A0A384NKM6	<i>APOH</i>	38298	12	-1.82
Leucine-rich alpha-2-glycoprotein	P02750	<i>LRG1</i>	38178	5	-1.86
Immunoglobulin heavy constant mu	P01871	<i>IGHM</i>	49440	15	-1.91
Transthyretin	P02766	<i>TTR</i>	15887	7	-1.97
Uncharacterized protein DKFZp686G11190	Q6MZQ6	<i>DKFZp686G11190</i>	52043	3	-2.06 *
Apolipoprotein E	P02649	<i>APOE</i>	36154	17	-2.15 *
Full-length cDNA clone CS0DD006YL02	Q86TT1		41273	2	-2.30 *
Testicular tissue protein Li 227	A0A140VK00		34259	17	-2.30 *
Apolipoprotein C-III	A3KPE2	<i>APOC3</i>	10852	5	-2.48 *
Haptoglobin-related protein	P00739	<i>HPR</i>	39030	7	-3.19 *#
Epididymis secretory sperm binding protein	A0A384MDQ7		46723	2	-4.45 *#

^a, Accession number in UniProt database (<https://www.uniprot.org>). ^b, Protein quantification values (Zq) are normalized log₂-ratios expressed in standard deviation units: Zq > 0 and Zq < 0 indicate increased or decreased, respectively, protein abundance in HDF compared with HF patients. ^c, Zq values in a colour scale; red and blue represent the increased or decreased protein levels, respectively, in HDF compared with HF. * Zq ≥ 2, and ≤ -2, were significant (p ≤ 0.05); #, significant FDRq value (< 0.05).

10. Supplementary Figure S1

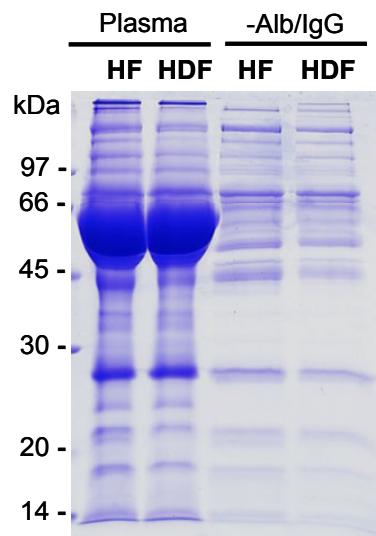


Figure S1. Plasma samples from HF and HDF patients were depleted from albumin and IgG. Plasma samples from patients on high-flux haemodialysis (HF) or on-line haemodiafiltration (HDF) were pooled, albumin and IgG depleted and ran into sodium dodecyl sulphate-polyacrylamide gel electrophoresis (SDS-PAGE), and the gel was stained with Coomassie blue for protein detection. The numbers on the left indicate the apparent molecular mass in kDa. The figure shows plasma samples from HF and HDF patients before and after albumin and IgG depletion (-Alb/IgG) in a whole stained gel.

11. Supplementary Figure S2

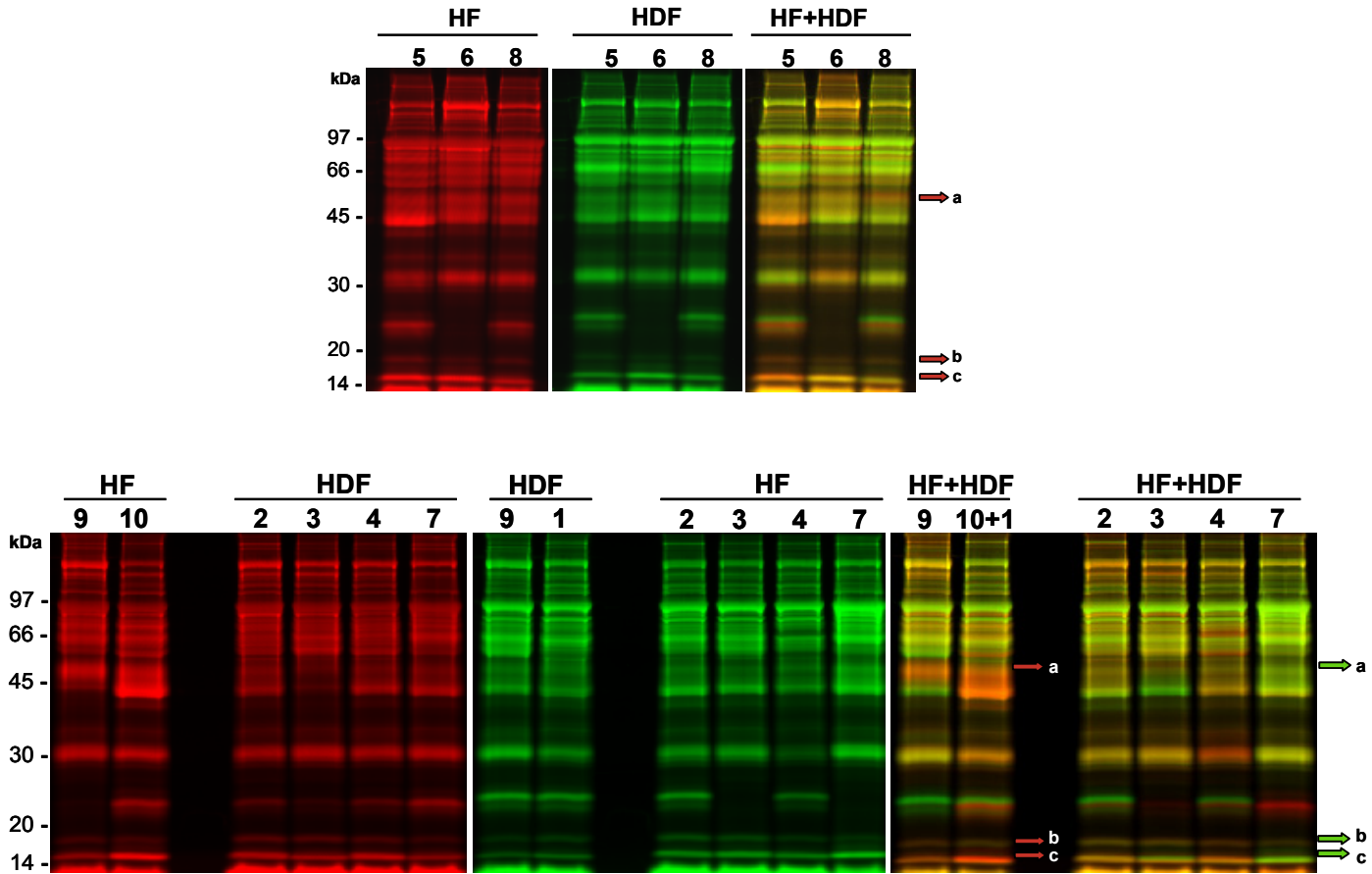
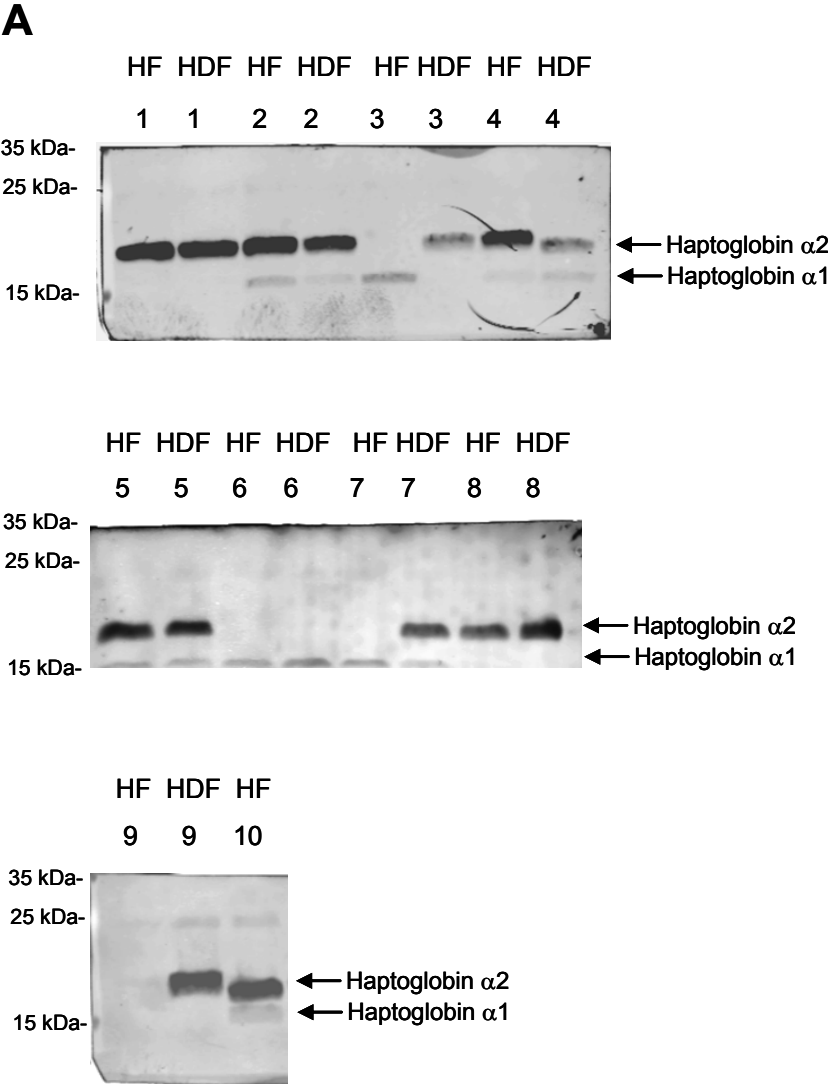


Figure S2. Differential protein detection of individual plasma samples of HF combined with HDF patients by fluorescence in gel electrophoresis (DIGE). The images show the combinations of individual paired and Cy-labelled samples of nine HF and HDF patients performed in the same way as described in Figure 1A. The scanned gel displays fluorescence-labelled proteins; proteins in the HF or HDF samples were visualised in red (Cy5-labelled) or green (Cy3-labelled). Proteins present in both HF and HDF samples were visualised in yellow due to the merge of the red and green labels (HF+HDF). The numbers on the left indicate the apparent molecular mass. The differentially detected proteins quantified in Figure 1B, *a* to *c*, are indicated in by arrows in the highest labelling colour code (red or green, HF patients). The fluorescence image corresponds to whole gels and adjacent empty lanes are not shown.

12. Supplementary Figure S3



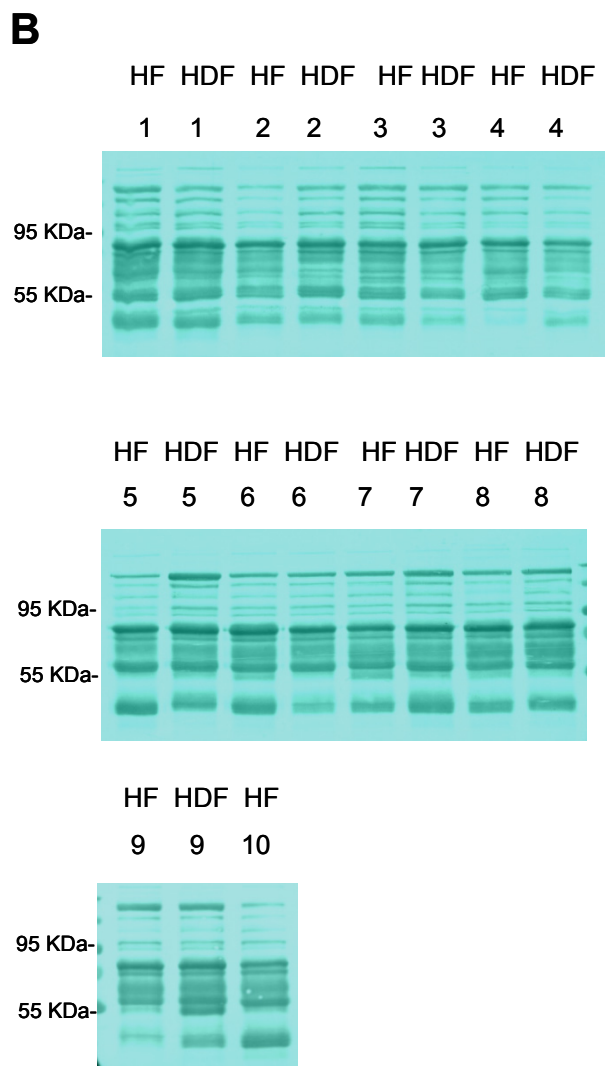


Figure S3. Haptoglobin α detection by western blot in HF and HDF plasma samples. (A) Individual albumin- and IgG-depleted plasma samples (5 μ g of protein) from 10 and 9 HF and HDF patients, respectively, were analysed independently by western blotting as describe in Methods using a mouse monoclonal anti-haptoglobin α antibody (sc-376893, Santa Cruz Biotechnology). In the images, arrows show the haptoglobin α 2 and α 1 detected. The numbers on the left indicate the apparent molecular mass from standards. (B) Blotted proteins were staining with Fast Green as loading control of the analyzed samples. . The upper piece of the PVDF membranes containing most proteins was used for staining. Images show whole blots.

11. Supplementary Figure S4

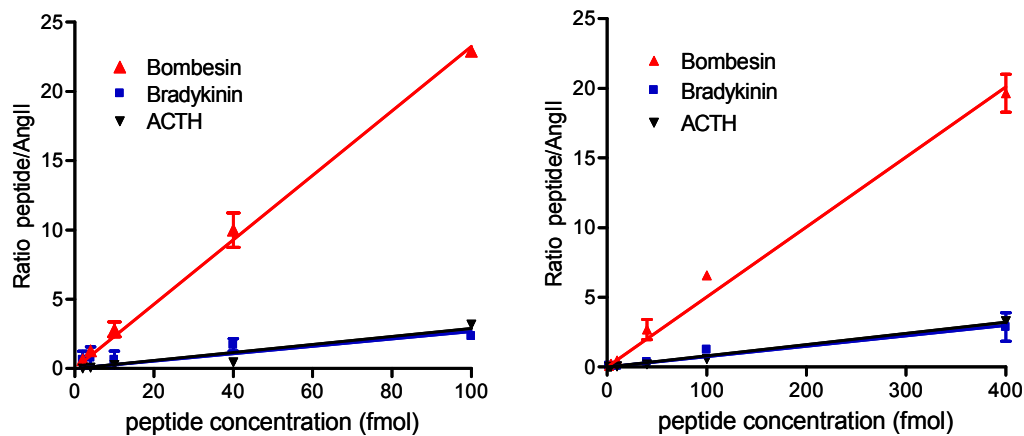


Figure S4. Quantification of the peptide intensities in the mass spectra from MALDI-TOF MS. Calibration experiments showing the linear correlation between peptide concentration and the ratio peptide intensity/angiotensin-II peptide intensity as relative intensity. Peptides used were bradykinin (765.85 Da), bombesin (1619.85 Da), and adrenocorticotrophic hormone (ACTH, 2465.67 Da). Peptide intensities were obtained from mass spectra by MALDI-TOF MS analysis, in the same conditions as mass fingerprinting analysis and using angiotensin-II peptide (1045.53 Da) as internal standard. Peptide concentrations used were 2-100 fmol, and 4 fmol of angiotensin-II (left graph), or 4-400 fmol, with 40 fmol of angiotensin-II (right graph). The correlation between peptide concentration and relative intensity for each peptide was very significant, $p < 0.0005$ (Pearson test).

13. Supplementary Figure S5

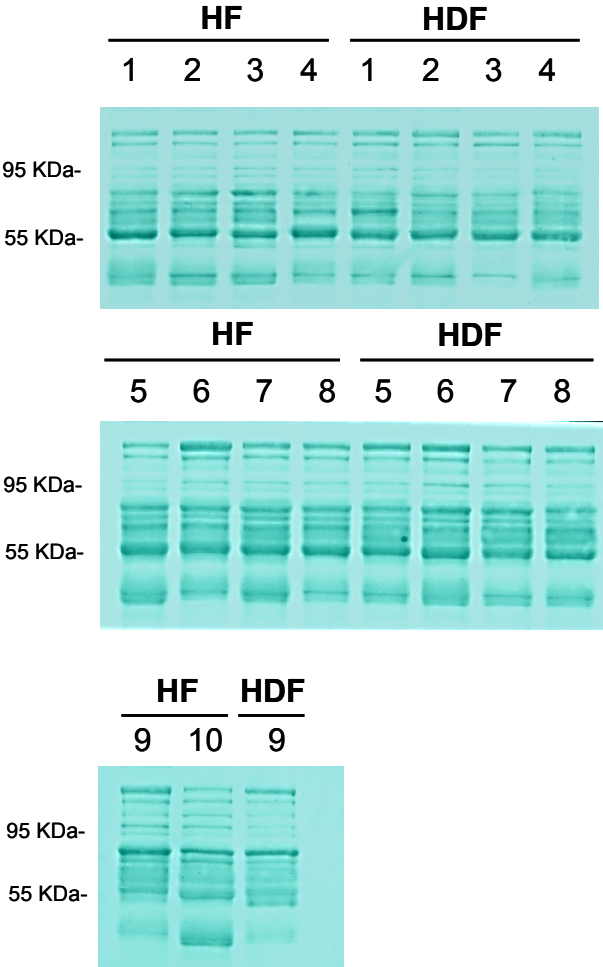


Figure S5. Loading control in western blot experiments for anti-TTR antibody detection. Blotted proteins were staining with Fast Green in the PVDF membranes as loading control of analyzed samples from HF and HDF patients. The upper piece of the membranes containing most proteins was used for staining and normalization dividing the TTR level of the western blot by the total level of stained proteins of the same loading lane. Images show stained whole blots.

15. Supplementary Figure S6

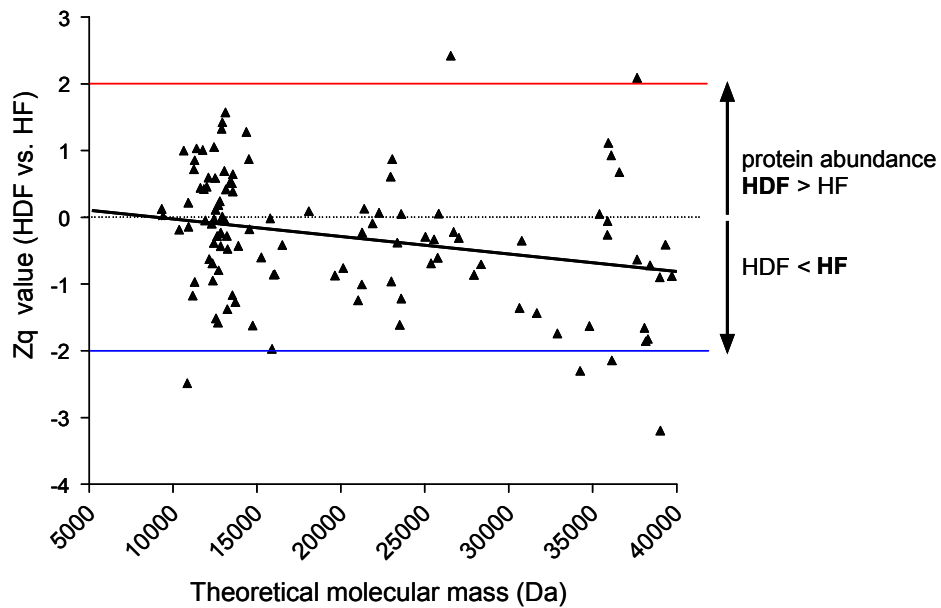
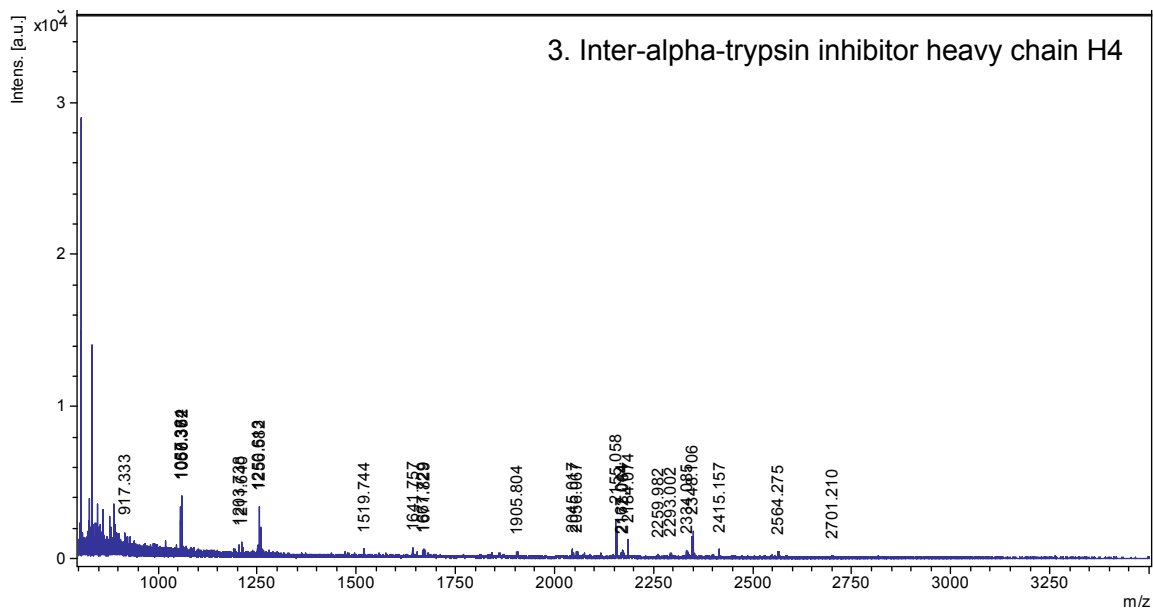
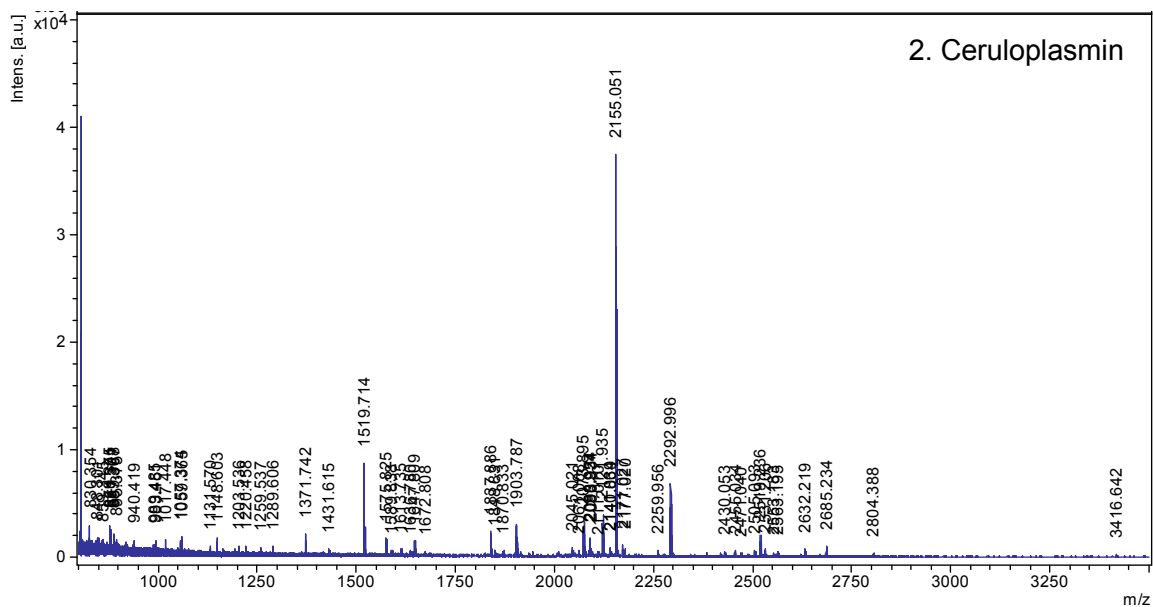
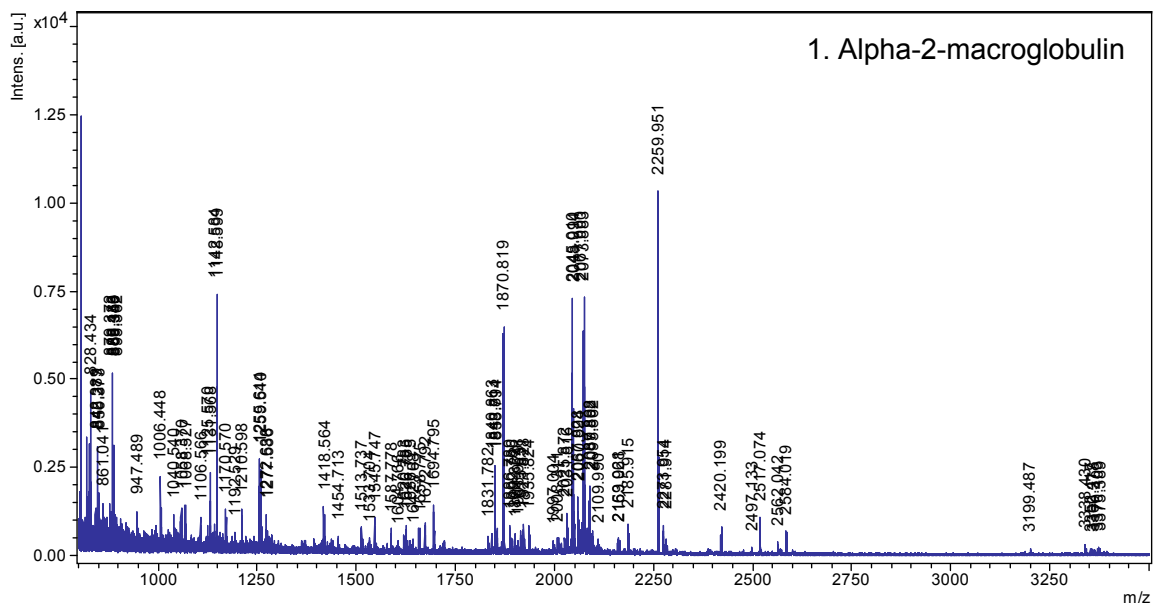
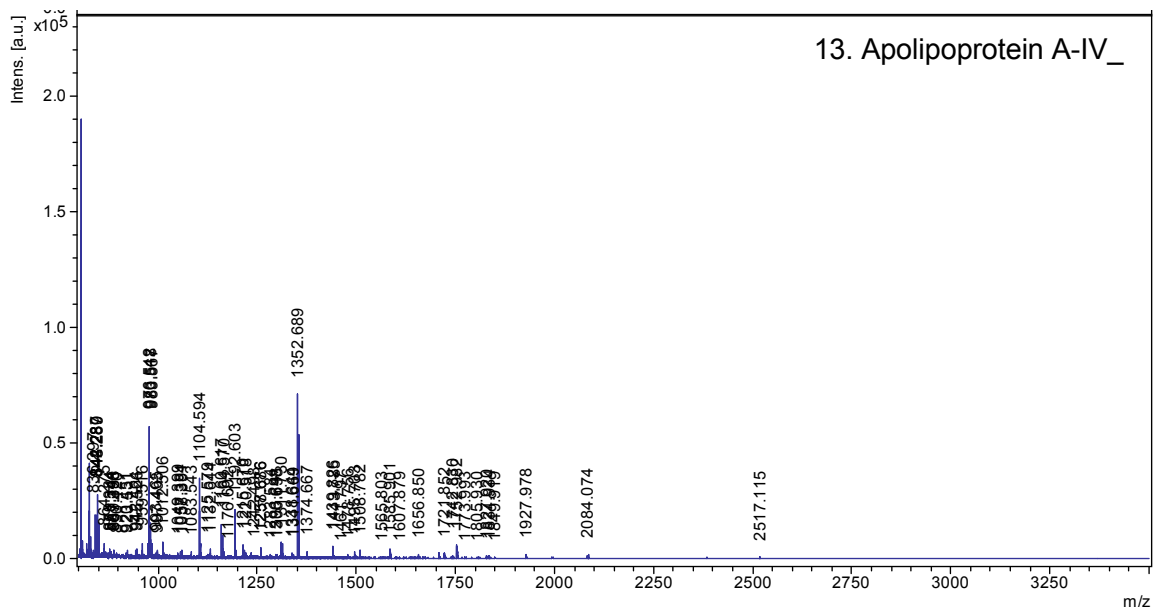
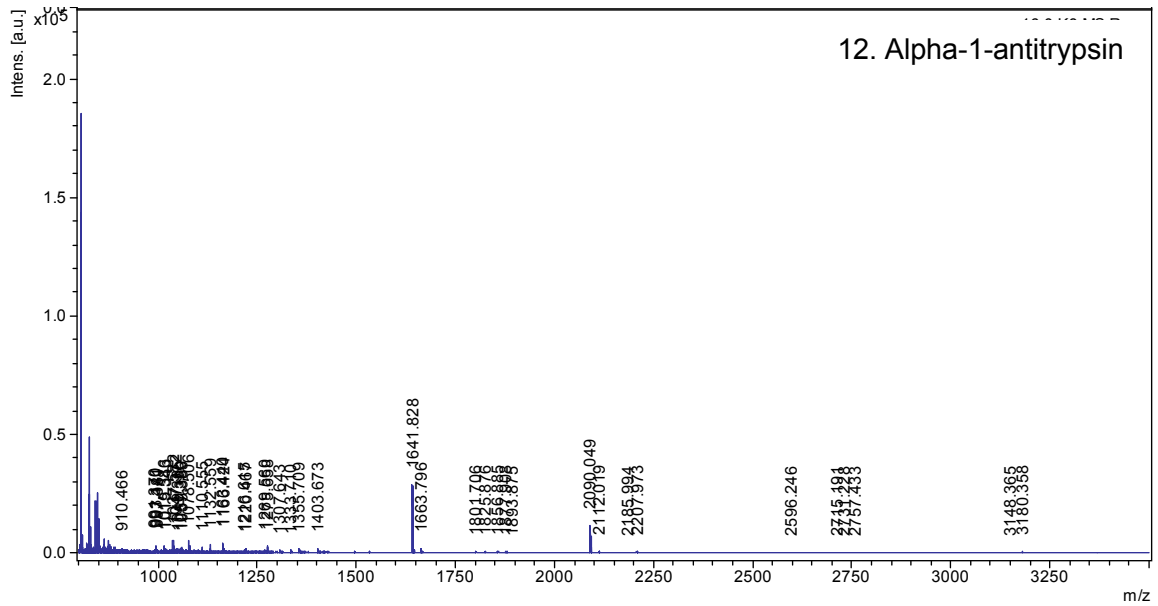
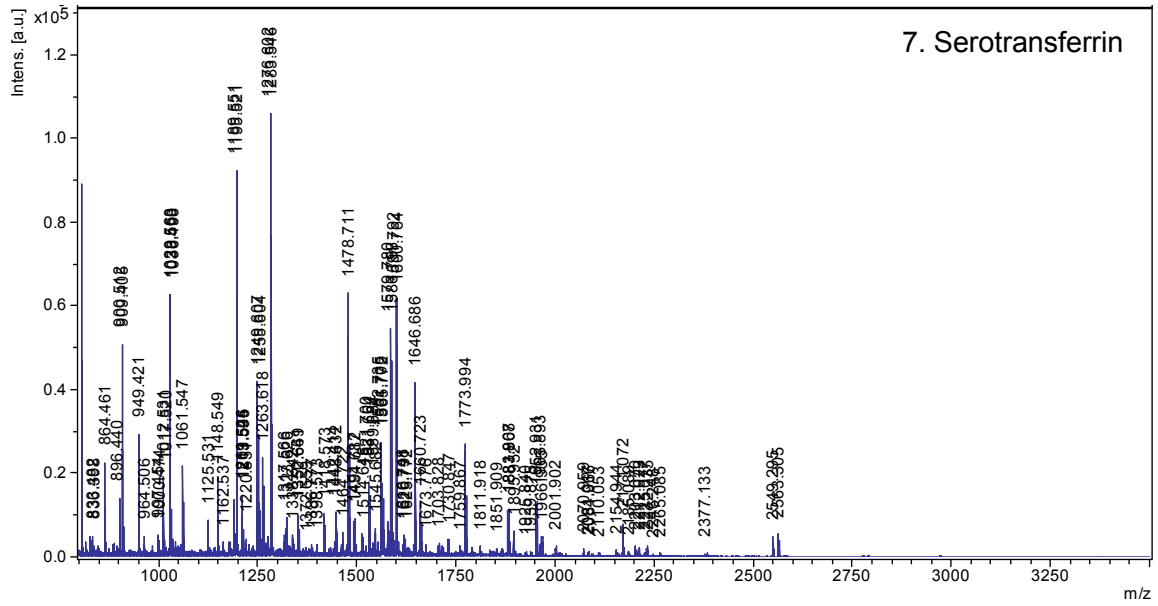
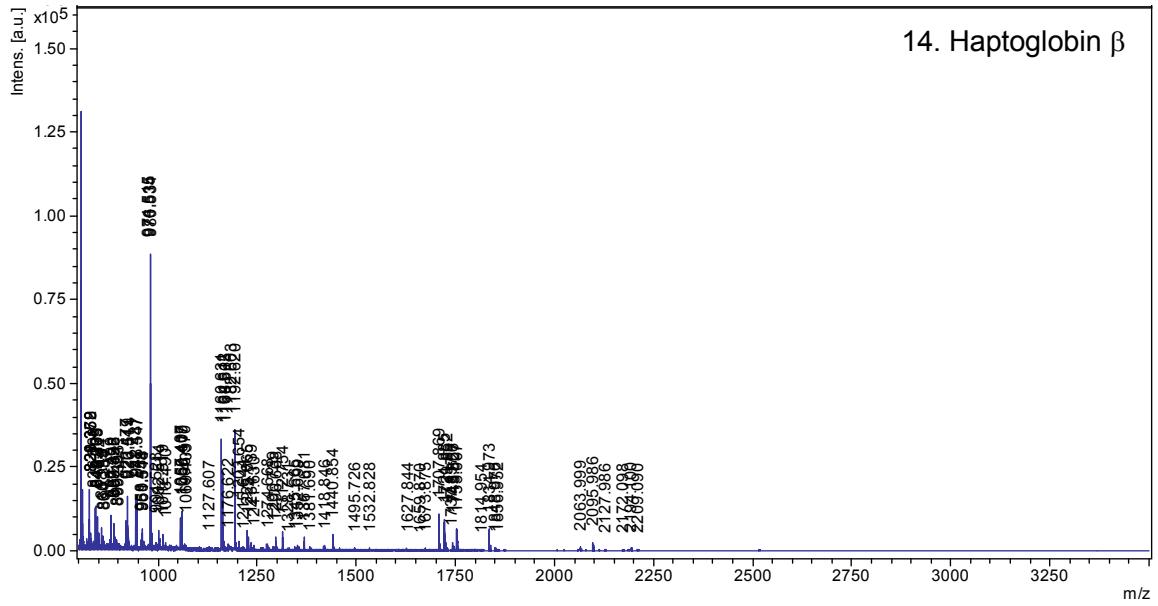


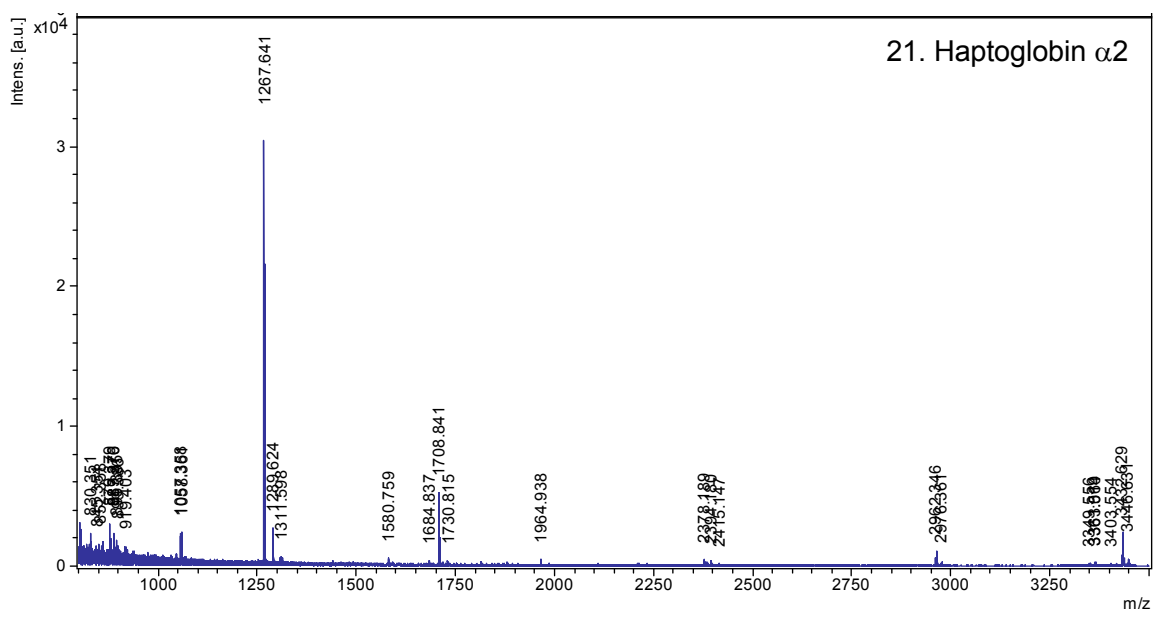
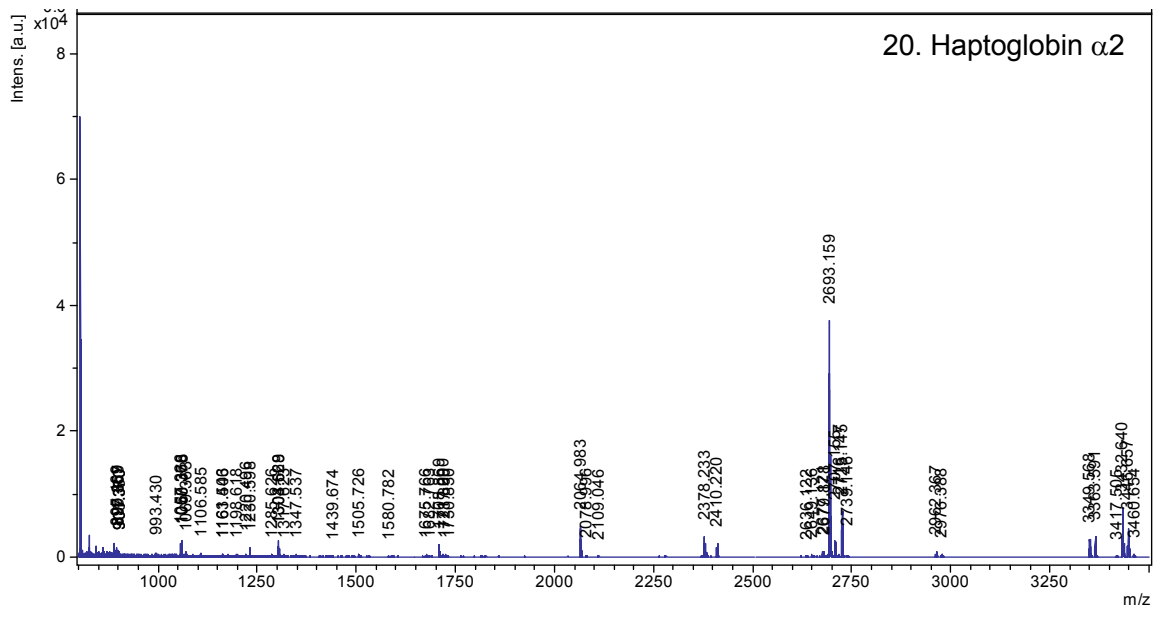
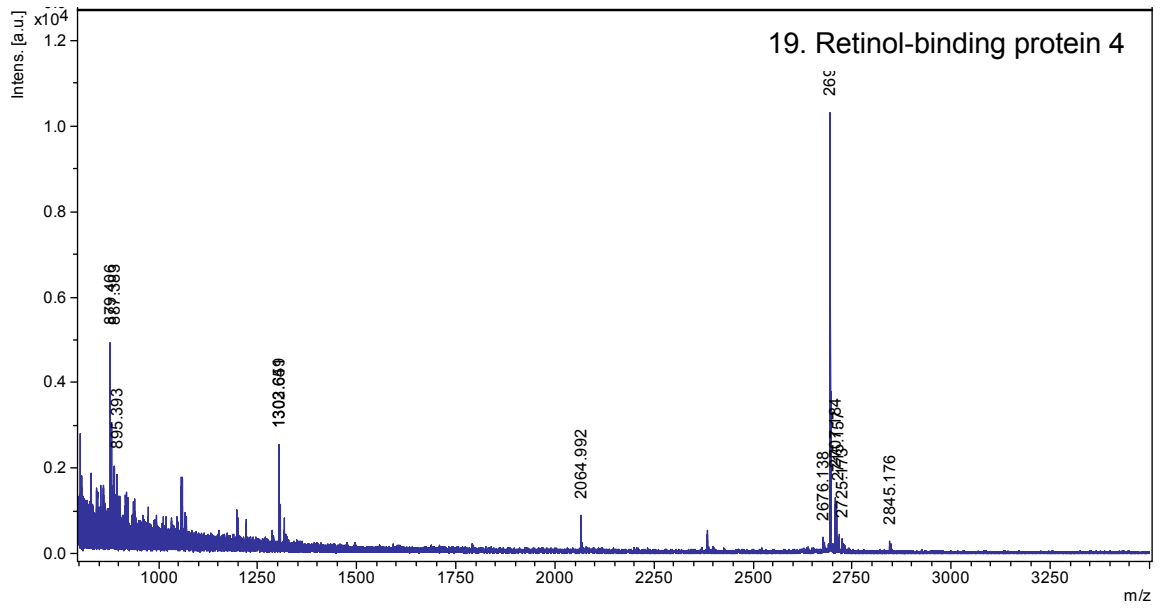
Figure S6. Correlation between the molecular mass (≤ 40 kDa) and the quantification value (Z_q) in the proteins identified in plasma samples of HDF compared with HF patients by LC-MS/MS analysis. Protein quantification values (Z_q) are the normalized \log_2 -ratios expressed in standard deviation units: $Z_q > 0$ and $Z_q < 0$ indicate increased or decreased, respectively, protein abundance in HDF compared with HF patients. $Z_q \geq 2$, and ≤ -2 , were significant ($p \leq 0.05$). Red and blue lines show the values $Z_q=2$ and $Z_q=-2$, respectively. The continuous black line shows the linear regression and correlation was significant ($p < 0.0043$, Pearson test).

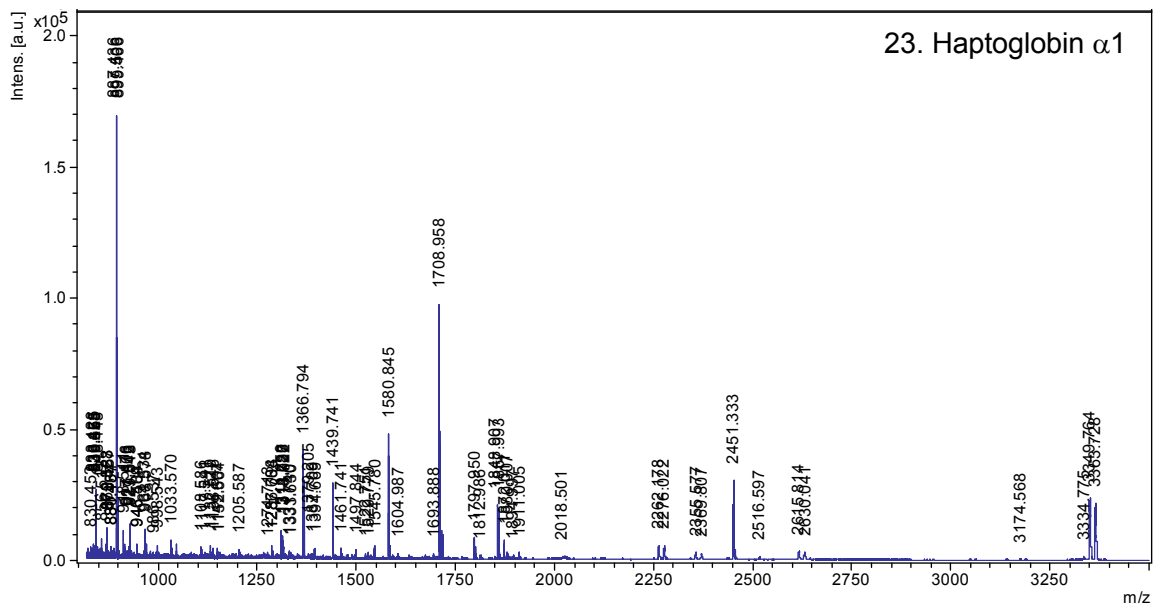
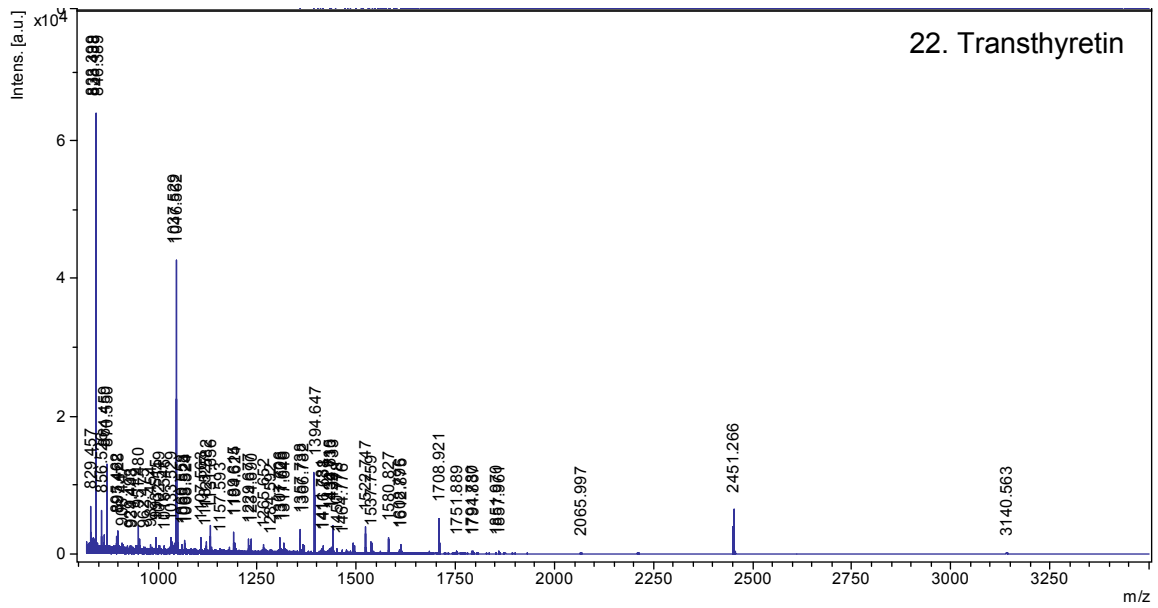
16. MALDI-TOF MS spectra and Mascot searches











Protein View: A2MG_HUMAN

Alpha-2-macroglobulin OS=Homo sapiens OX=9606 GN=A2M PE=1 SV=3

Database: SwissProt
Score: 95
Expect: 6e-06
Monoisotopic mass (M_r): 164613
Calculated pI: 6.03
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of A2MG_HUMAN against nr.](#)

Search parameters

Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)
Mass values searched: 86
Mass values matched: 24

Protein sequence coverage: 20%

Matched peptides shown in **bold red**.

1	MGKNKLLHPS	LVLLLLLVLLP	TDASVSGKPQ	YMVLVPSLLH	TETTEKGCVL
51	LSYLNETHVT	SASLESVRGN	RSLFTDLEAE	NDVLHCVAFA	VPKSSSNEEV
101	MFLTIVQVKG	TQEFKKRTTV	MVKNEDSLVF	VQTDKSIYKP	GQTVKFRVVS
151	MDENFHPLNE	LIPLVYIQDP	KGNR IAQWQS	FQLEGGLKQF	SFPLSSEPFQ
201	GSYKVVVQKK	SGGR TEHPFT	VEEFVLPKFE	VQVTVPKIIT	ILEEEMNVSV
251	CGLYTYGKPV	PGHVTVSICR	KYSDASDCHG	EDSQAFCEKF	SGQLNSHGCF
301	YQQVKTKVFQ	LKRKEYEMKL	HTEAQIQEEG	TVVELTGRQS	SEITRTITKL
351	SFVKVDSHFR	QGIPFFGQVR	LVDGKGVPIP	NKVIFIRGNE	ANYYSNATTD
401	EHGLVQFSIN	TTNVMGTSLT	VRVNYKDR SP	CYGYQWVSEE	HEEAHHTAYL
451	VFSPSKS FEVH	LEPMSHELPC	GHTQTVQAHY	ILNGGTLGL	KKLSFYYLIM
501	AKGGIVRTGT	HGLLVKQEDM	KGHFSISIPV	KSDIAPVARL	LIYAVLPTGD
551	VIGDSAKYDV	ENCLANK VDL	SFSPSQSLPA	SHAHLRVTAA	PQSVCALRAV
601	DQSVLLMKPD	AELSASSVYN	LLPEKDLTGF	PGPLNDQDNE	DCINRHNVI
651	NGITYTPVSS	TNEKDMYSFL	EDMGLKAFTN	SKIRKPK MCP	QLQQYEMHGP
701	EGLRVGFYES	DVMGRGHARL	VHVEEPTET	VRKYFPETWI	WDLVVVNSAG
751	VAEVGVTVPD	TITEWKAGAF	CLSEDAGLGI	SSTASL RAFQ	PPFVELTMPY
801	SVIRGEAFTL	KATVLNLYLPK	CIRVSVQLEA	SPAFLAVPVE	KEQAPHCICA
851	NGRQTVSWAV	TPKSLGNVNF	TVSAEALESQ	ELCGTEVPSV	PEHGRKDTVI
901	KPLLVEPEGL	EKETTFNSLL	CPSGGEVSEE	LSLKL PPNVV	EESARASVSV
951	LGDILGSAMQ	NTQNLLQMPY	GCGEQNMVLF	APNIYVLDYL	NETQQLTPEI
1001	KSK AIGYLNT	GYQRQLNYKH	YDGSYSTFGE	RYGRNQGNTW	LTAFLVLTFA
1051	QARAYIFIDE	AHITQALIWL	SQRQKDNCF	RSSGSLNNA	IKGGVEDEV
1101	LSAYITIALL	EIPLTVTHPV	VRNALFCLES	AWKTAQEGDH	GSHVYTKALL
1151	AYAFALAGNQ	DKRKEVLKSL	NEEAVKK DNS	VHWERPQKPK	APVGHFYEPQ

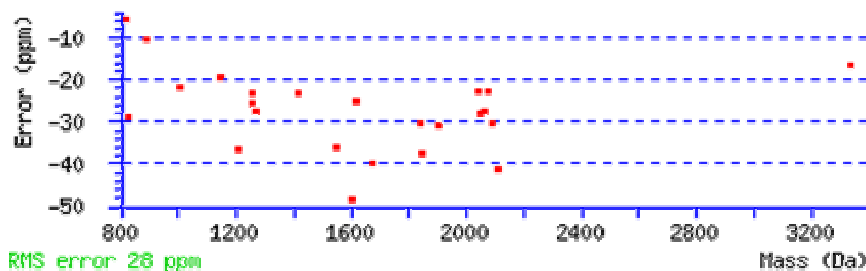
1201 APSAEVEMTS YVLLAYLTAQ PAPTSEDLTS ATNIVKWITK QQNAQGGFSS
 1251 TQDTVVALHA LSKYGAATFT RTGKAAQVTI QSSGTFSSKF QVDNNRLLL
 1301 QQVSLPELPG EYSMKVTGEG CVYLQTSKY NILPEKEEFP FALGVQTLPO
 1351 TCDEPKAHTS FQISLSVSYT GSRASANMAI VDVKMVSGFI PLKPTVKMLE
 1401 RSNHVS RTEV SSNHVLIYLD KVSNQTLSEF FTVLQDVPVR DLKPAIVKVY
 1451 DYYETDEFAI AEYNAPCSKD LGNA

Unformatted sequence string: **1474 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Peptide
175 - 188	1604.7657	1603.7584	1603.8358	-48.3	0	R.IAQWQSFQLEGGLK.Q
189 - 204	1848.8113	1847.8040	1847.8730	-37.3	0	K.QFSFPLSSEPFQGSYK.V
215 - 228	1672.7917	1671.7844	1671.8508	-39.7	0	R.TEHPFTVEEFVLPK.F
290 - 305	1899.8224	1898.8151	1898.8734	-30.7	0	K.FSQQLNSHGCFYQVK.T
320 - 338	2109.9903	2108.9831	2109.0702	-41.3	0	K.LHTEAQIQEETVVVELTGF
339 - 345	820.4113	819.4040	819.4086	-5.59	0	R.QSSEITR.T
361 - 370	1148.5988	1147.5916	1147.6138	-19.4	0	R.QGIPFFGQVR.L
429 - 456	3338.4298	3337.4225	3337.4775	-16.5	0	R.SPCYGYQWVSEEHAAHTI
532 - 539	828.4336	827.4263	827.4501	-28.7	0	K.SDIAPVAR.L
568 - 586	2048.9940	2047.9867	2048.0439	-27.9	0	K.VDLSFSPSQSLPASHALF
587 - 598	1272.6381	1271.6308	1271.6656	-27.4	0	R.VTAAPQSVCALR.A
688 - 704	2073.8831	2072.8758	2072.9230	-22.8	0	K.MCPQLQYEMHGPEGLR.V
688 - 704	2089.8621	2088.8548	2088.9179	-30.2	0	K.MCPQLQYEMHGPEGLR.V
705 - 715	1259.5404	1258.5331	1258.5652	-25.5	0	R.VGFYESDVMGR.G
720 - 732	1545.7465	1544.7393	1544.7947	-35.9	0	R.LVHVEEPHTETVR.K
788 - 804	2045.0104	2044.0031	2044.0492	-22.6	0	R.AFQPPFVELTMPYSVIR.G
788 - 804	2060.9944	2059.9872	2060.0441	-27.6	0	R.AFQPPFVELTMPYSVIR.G
935 - 945	1210.5984	1209.5911	1209.6353	-36.5	0	K.LPPNVVEESAR.A
1004 - 1014	1255.6135	1254.6063	1254.6357	-23.4	0	K.AIGYLNTGYQR.Q
1020 - 1031	1418.5643	1417.5570	1417.5899	-23.2	0	K.HYDGSYSTFGER.Y
1178 - 1190	1620.7833	1619.7761	1619.8168	-25.2	0	K.DNSVHWERPQKPK.A
1264 - 1271	886.4327	885.4254	885.4345	-10.2	0	K.YGAATFTR.T
1290 - 1297	1006.4483	1005.4410	1005.4628	-21.6	0	K.FQVDNNR.L
1357 - 1373	1840.8629	1839.8556	1839.9115	-30.4	0	K.AHTSFQISLSVSYTGSR.S

No match to: 826.2877, 842.3894, 848.2769, 850.3727, 861.0413, 879.3721, 888.3465, 889.3590, 895.3618, 947.4890, 1040.5396, 1058.3703, 1068.5270, 1106.5661, 1125.5697, 1131.5676, 1142.5836, 1170.5696, 1192.5295, 1277.5860, 1454.7129, 1513.7372, 1533.7041, 1587.7778, 1626.7851, 1636.7750, 1642.7680, 1658.7747, 1694.7954, 1831.7819, 1853.7943, 1870.8189, 1886.7886, 1892.7857, 1905.9787, 1913.8309, 1921.8281, 1935.8236, 1997.0044, 2008.3115, 2025.8719, 2031.8163, 2057.8285, 2067.0004, 2087.8984, 2095.8619, 2159.9278, 2163.0614, 2185.9147, 2259.9510, 2273.9541, 2281.9145, 2420.1989, 2497.1328, 2517.0739, 2562.0416, 2584.0188, 3199.4870, 3352.4173, 3360.3783, 3370.4030, 3375.3890



Protein View: CERU_HUMAN

Ceruloplasmin OS=Homo sapiens OX=9606 GN=CP PE=1 SV=1

Database: SwissProt
Score: 112
Expect: 1.3e-07
Monoisotopic mass (M_r): 122983
Calculated pI: 5.44
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of CERU_HUMAN against nr.](#)

Search parameters

Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)
Mass values searched: 66
Mass values matched: 19

Protein sequence coverage: 24%

Matched peptides shown in **bold red**.

```

1  MKILILGIFL FLCSTPAWAK EKHYIIGIIE TTWDYASDHG EKKLISVDTE
51  HSNIYLQNGP DRIGRLYKKA LYLQYTDETF RTTIEKPVWL GFLGPIIKAE
101 TGDKVYVHLK NLASRPYTFH SHGITYYKEH EGAIYDNTT DFQRADDKVY
151 PGEQYTYMLL ATEEQSPGEG DGNCVTRIYH SHIDAPKDIA SGLIGPLIIC
201 KKDSL DKEKE KHIDREFVVM FSVVDENFSW YLEDNIKTYC SEPEKVDKDN
251 EDFQESNRMY SVNGYTFGSL PGLSMCAEDR VKWYLFMGMN EVDVHAAFFH
301 GQALT NKNYR IDTINLFPAT LFDAYMVAQN PGEWMLSCQN LNHLKAGLQA
351 FFQVQECNKS SSKDNIRGKH VRHYIAAEE IIWNYAPSGI DIFTKENLTA
401 PGSDSAVFFE QGTRIGGSY KKL VYREYTD ASFTNRKERG PEEEHLGILG
451 PVIWAEVGDT IRVTFHNKGA YPLSIEPIGV RFNKNNEGTY YSPNYNPQSR
501 SVPPSASHVA PTEFTYEWV VPKEVGPTNA DPVCLAKMY Y SAVDPTKDIF
551 TGLIGPMKIC KKGSLHANGR QKDVDKEFY L FPTVFDENES LLEDNIRMF
601 TTAPDQVDKE DEDFQESNKM HSMNGFMYGN QPGLTMCKGD SVVWYLF SAG
651 NEADVHGIYF SGN TYLWRGE RRDTANLFPQ TSLTLHMWPD TEGTFNVECL
701 TTDHYTGGMK QKYTVNQCR QSEDSTFYLG ERTYYIAAVE VEWDYSPQRE
751 WEKELHHLQE QNVSNAFLDK GEFYIGSKYK KVVYRQYTD TFRVPVERKA
801 EEEHLGILGP QLHADVGDKV KII FKNMATR PYSIHAHVQ TESSTVTPTL
851 PGETLTYVWK IPERSGAGTE DSACIPWAYY STVDQVKDLY SGLIGPLIVC
901 RRPYLKVFNP RKRLEFALLE LVFDENESWY LDDNIKTYSD HPEKVNKDDE
951 EFIESNKMHA INGRMFGNLQ GLTMHVGDEV NWYLMGMGNE IDLHTVHFHG
1001 HSFQYKHRGV YSSDVFDFP GTYQTMLEFP RTPGIWLLHC HVTDHIHAGM
1051 ETTYTVLQNE DTKSG

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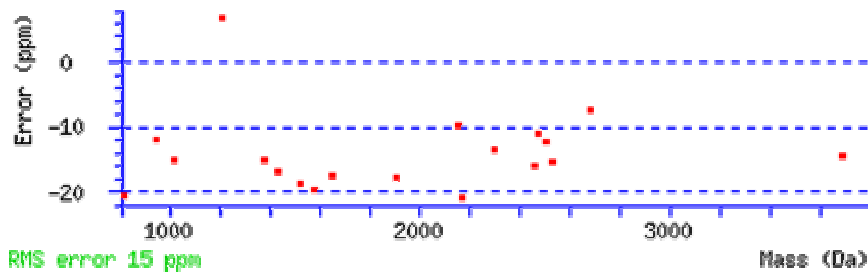
Unformatted sequence string: **1065 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass

Show matched peptides only predicted peptides also

Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Peptide
44 - 62	2171.0270	2170.0197	2170.0654	-21.1	0	K.LISVDTEHSNIYLQNGPDF
69 - 81	1647.8086	1646.8013	1646.8304	-17.7	1	K.KALYLQYTDETFR.T
70 - 81	1519.7144	1518.7072	1518.7354	-18.6	0	K.ALYLQYTDETFR.T
111 - 128	2155.0509	2154.0436	2154.0647	-9.77	0	K.NLASRPYTFHSHGITYYK.
145 - 177	3693.5792	3692.5720	3692.6247	-14.3	1	R.ADDKVYPGEQYTYMLLATE
259 - 280	2455.0337	2454.0264	2454.0654	-15.9	0	R.MYSVNGYTFGSLPGLSMCA
259 - 280	2471.0401	2470.0328	2470.0603	-11.1	0	R.MYSVNGYTFGSLPGLSMCA
427 - 436	1203.5360	1202.5287	1202.5204	6.91	0	R.EYTDASFTNR.K
469 - 481	1371.7425	1370.7352	1370.7558	-15.0	0	K.GAYPLSIEPIGVR.F
482 - 500	2292.9955	2291.9882	2292.0195	-13.6	1	R.FNKNNEGTYYSNPYNPQSF
485 - 500	1903.7868	1902.7795	1902.8132	-17.7	0	K.NNEGTYYSNPYNPQSR.S
501 - 523	2531.2065	2530.1992	2530.2380	-15.3	0	R.SVPPSASHVAPTETFTYEW
563 - 570	811.4003	810.3930	810.4096	-20.5	0	K.GSLHANGR.Q
713 - 719	940.4193	939.4120	939.4232	-11.9	0	K.YTVNQCR.R
721 - 732	1431.6146	1430.6073	1430.6314	-16.8	0	R.QSEDSTFYLGGER.T
786 - 793	1017.4482	1016.4409	1016.4563	-15.2	0	R.QYTDSTFR.V
865 - 887	2505.0933	2504.0860	2504.1166	-12.2	0	R.SGAGTEDSACIPWAYYSTV
888 - 901	1575.8252	1574.8179	1574.8490	-19.8	0	K.DLYSGLIGPLIVCR.R
1009 - 1031	2685.2344	2684.2271	2684.2469	-7.37	0	R.GVYSSDVFDFIPGTYQTLE

No match to: 826.2964, 830.3540, 848.3008, 859.3449, 879.3752, 881.3009, 887.3666, 889.3685, 895.3748, 989.4849, 993.4508, 1057.3741, 1059.3650, 1131.5696, 1148.6031, 1220.4582, 1259.5374, 1289.6064, 1589.8379, 1613.7348, 1636.7803, 1672.8079, 1837.8858, 1848.8312, 1870.8327, 2045.0215, 2061.0078, 2073.8950, 2088.9267, 2090.9538, 2093.9340, 2111.0095, 2121.9353, 2140.0393, 2141.8642, 2177.0200, 2259.9564, 2430.0527, 2519.2360, 2553.1833, 2563.1952, 2632.2186, 2804.3878, 3416.6424, 3487.5071, 3707.5985, 3867.3625



Protein View: ITIH4_HUMAN

**Inter-alpha-trypsin inhibitor heavy chain H4 OS=Homo sapiens
OX=9606 GN=ITIH4 PE=1 SV=4**

Database: SwissProt
Score: 48
Expect: 0.33
Monoisotopic mass (M_r): 103521
Calculated pI: 6.51
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of ITIH4 HUMAN against nr.](#)

Search parameters

Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)
Mass values searched: 26
Mass values matched: 6

Protein sequence coverage: 10%

Matched peptides shown in **bold red**.

```

1  MKPPRPVRTC  SKVLVLLSLL  AIHQTTTAEK  NGIDIYSLTV  DSRVSSRFAH
51  TVVTSRVVNR  ANTVQEATFQ  MELPKKAFIT  NFSMIIDGMT  YPGIIKEKAE
101 AQAQYSAAVA  KGKSAGLVKA  TGRNMEQFQV  SVSVAPNAKI  TFELVYEELL
151 KRRLGVYELL LKVRPQQLVK  HLQMDIHIFE  PQGISFLETE  STFMTNQLVD
201 ALTTWQNKTK  AHIRFKPTLS  QQQKSPEQQE  TVLDGNLIIR  YDVDRAISGG
251 SIQIENGYFV  HYFAPEGLTT  MPKNVVFVID  KSGSMSGRKI  QQTREALIKI
301 LDDLSPRDQF  NLIVFSTEAT  QWRPSLVPAS  AENVNKARSF  AAGIQALGGT
351 NINDAMLMAV  QLLDSSNQEE  RLPEGSVSLI  ILLTDGDPTV  GETNPRSIQN
401 NVREAVSGRY  SLFCLGFGFD  VSYAFLEKLA  LDNGGLARRI  HEDSDSALQL
451 QDFYQEVANP  LLTAVTFEYP  SNAVEEVTQN  NFRLLFKGSE  MVVAGKLQDR
501 GPDVLTATVS  GKLPTQNITF  QTESSVAEQE  AEFQSPKYIF  HNFMERLWAY
551 LTIQQLLEQT  VSASDADQQA  LRNQALNLSL  AYSFVTPLTS  MVVTKPDDQE
601 QSQVAEKPME  GESRNRNVHS  GSTFFKYYLQ  GAKIPKPEAS  FSPRRGWNRQ
651 AGAAGSRMNF  RPGVLSSRQL  GLPGPPDVPD  HAAYHPFRRL  AILPASAPPA
701 TSNPDAVSR  VMNMKIEETT  MTTQTPAPIQ  APSAILPLPG  QSVRLCVDP
751 RHRQGPVNLL  SDPEQGEVET  GQYEREKAGF  SWIEVTFKNP  LVVWHASPEH
801 VVVTRNRRSS  AYKWKETLFS  VMPGLKMTMD  KTGLLLLSDP  DKVTIGLLFW
851 DGRGEGRLRL  LRDTDRFSSH  VGGTLGQFYQ  EVLWGSPAAS  DDGRRTLRVQ
901 GNDHSATRER  RLDYQEGPPG  VEISCWSVEL

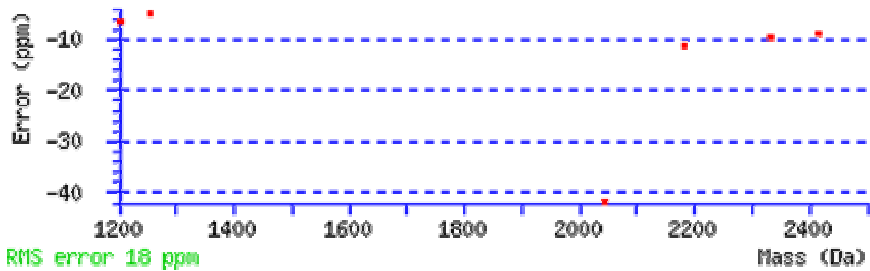
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Unformatted sequence string: **930 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Peptide
153 - 162	1203.7383	1202.7310	1202.7387	-6.36	1	R.RLGVYELLK.V
538 - 546	1256.5818	1255.5745	1255.5808	-4.99	0	K.YIFHNFMER.L
669 - 688	2184.0738	2183.0666	2183.0912	-11.3	0	R.QLGLPGPPDVPDHAAYHPE
690 - 710	2045.0168	2044.0095	2044.0953	-42.0	0	R.LAILPASAPPATSNPDPVAV
754 - 775	2415.1570	2414.1497	2414.1714	-8.98	0	R.QGPVNLLSDPEQGVETGQ
911 - 930	2334.0845	2333.0772	2333.0998	-9.67	1	R.RLDYQEGPPGVEISCWSVE

No match to: 917.3328, 1057.3640, 1058.3722, 1060.3816, 1211.6404, 1250.6130, 1519.7444, 1641.7572, 1667.7294, 1671.8289, 1905.8039, 2056.0675, 2155.0575, 2167.0439, 2172.0641, 2259.9816, 2293.0024, 2348.1058, 2564.2746, 2701.2096



Protein View: TRFE_HUMAN

Serotransferrin OS=Homo sapiens OX=9606 GN=TF PE=1 SV=3

Database: SwissProt
Score: 152
Expect: 1.3e-11
Monoisotopic mass (M_r): 79294
Calculated pI: 6.81
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of TRFE_HUMAN against nr.](#)

Search parameters

Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)
Mass values searched: 95
Mass values matched: 25

Protein sequence coverage: 33%

Matched peptides shown in **bold red**.

1 MRLAVGALLV CAVLGLCLAV PDK**TVRWCAV SEHEATK**CQS FRDHMKSVIP
 51 SDGPSVACVK **KASYLDCIRA** IAANEADAVT LDAGLVYDAY LAPNNLKPVV
 101 AEFYGSKEDP QTFYYAVAVV **KDSGFQMNQ** LRGKKSCHTG LGR**SAGWNIP**
 151 **IGLLYCDLPE** PRKPLEKAVA NFFSGSCAPC ADGTDLPQLC QLCPGCGCST
 201 LNQYFGYSGA FKCLKDGAGD VAFVKHSTIF ENLANK**ADRD QYELLCLDNT**
 251 **RKPVDEYKDC HLAQVPSHTV** VARSMGGKED LIWELLNQAQ EHFVKDK**SKE**
 301 **FQLFSSPHGK** DLLFKDSAAG FLKVPPRMDA **KMYLGYEYVT AIR**NLREGTC
 351 PEAPTDECKP VK**WCALSHHE RLKCD**EWVSN **SVGK**IECVSA ETTEDCIAKI
 401 MNGEADAMSL DGGFVYIAGK CGLVPVLAEN YNKSDNCEDT PEAGYFAIAV
 451 VKK**SASDLTW DNLK**GKKSCH TAVGRTAGWN IPMGLLYNKI NHCRFDEFFS
 501 EGCAPGSKKD SSLCKLCMGS GLNLCEPNNK **EGYYGYTGAF** RCLVEKGDVA
 551 FVKHQTVPQN TGGKNPDPWA **KNLNEKDYEL LCLDGTRKPV** **EYANCHLAR**
 601 **APNHAVVTRK** DKEACVHKIL RQQQHLLFGSN VTDCSGNFCL FRSETK**DLLF**
 651 **RDDTVCLAKL** HDRNTYEKYL **GEEYK**AVGN LRK**CSTSSLL** **EACTFR**RP

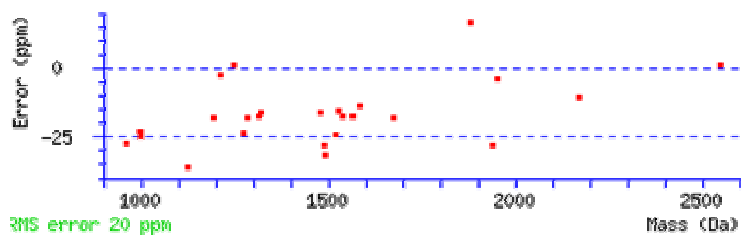
Unformatted sequence string: **698 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Peptide
24 - 37	1673.7763	1672.7690	1672.7991	-18.0	1	K.TVRWCAVSEHEATK.C
27 - 37	1317.5662	1316.5590	1316.5819	-17.4	0	R.WCAVSEHEATK.C
61 - 69	1125.5315	1124.5242	1124.5648	-36.1	1	K.KASYLDCIR.A
62 - 69	997.4542	996.4469	996.4698	-23.0	0	K.ASYLDCIR.A
122 - 132	1323.6261	1322.6188	1322.6401	-16.1	1	K.KDSGFQMNQLR.G
123 - 132	1211.5441	1210.5368	1210.5401	-2.70	0	K.DSGFQMNQLR.G + Oxidation (M)
144 - 162	2171.0721	2170.0649	2170.0881	-10.7	0	R.SAGWNIPIGLLYCDLPEPR.K
237 - 251	1881.9075	1880.9003	1880.8687	16.8	1	K.ADRDQYELLCLDNTR.K
240 - 251	1539.6842	1538.6769	1538.7035	-17.3	0	R.DQYELLCLDNTR.K
252 - 273	2549.2953	2548.2881	2548.2856	0.95	1	R.KPVDEYKDCHLAQVPSHTVVAR.S
298 - 310	1491.7170	1490.7097	1490.7518	-28.2	1	K.SKEFQLFSSPHGK.D
300 - 310	1276.6019	1275.5946	1275.6248	-23.7	0	K.EFQLFSSPHGK.D
328 - 343	1939.8748	1938.8675	1938.9219	-28.1	1	R.MDAKMYLGYEYVTAIR.N + Oxidation (M)
332 - 343	1478.7112	1477.7039	1477.7275	-16.0	0	K.MYLGYEYVTAIR.N
332 - 343	1494.6823	1493.6750	1493.7224	-31.7	0	K.MYLGYEYVTAIR.N + Oxidation (M)
363 - 371	1195.5207	1194.5134	1194.5352	-18.3	0	K.WCALSHHER.I
372 - 384	1521.6995	1520.6922	1520.7293	-24.4	1	R.LKCDEWVNSVVGK.I
454 - 464	1249.6074	1248.6001	1248.5986	1.22	0	K.SASDLTWDNLK.G

Start - End	Observed	Mr(expt)	Mr(calc)	ppm	M	Peptide
531 - 541	1283.5464	1282.5391	1282.5618	-17.7	0	K.EGYGYTGAFR.C
572 - 587	1952.9306	1951.9233	1951.9309	-3.90	1	K.NLNEKDYELLCLDGTR.K
588 - 600	1586.7526	1585.7454	1585.7671	-13.7	0	R.KPVEEYANCHLAR.A
601 - 609	964.5057	963.4985	963.5250	-27.6	0	R.APNHAVVTR.K
647 - 659	1565.7719	1564.7646	1564.7919	-17.5	1	K.DLLFRDDTVCLAK.L
669 - 676	1000.4736	999.4664	999.4913	-24.9	0	K.YLGEEYVK.A
684 - 696	1531.6641	1530.6568	1530.6807	-15.6	0	K.CSTSSLLEACTFR.R

No match to: 806.3029, 833.4825, 836.3980, 864.4605, 896.4401, 900.5118, 909.4053, 949.4210, 1012.5311, 1017.5199, 1029.5627, 1033.5502, 1036.4978, 1061.5471, 1148.5492, 1162.5366, 1189.5506, 1209.5260, 1213.5953, 1220.5509, 1255.6044, 1263.6182, 1338.6423, 1350.6488, 1352.6613, 1372.6586, 1386.7772, 1398.5728, 1418.5731, 1442.4144, 1448.6320, 1464.7217, 1514.6448, 1545.6818, 1552.7351, 1561.7047, 1579.7796, 1591.7917, 1600.7638, 1616.7480, 1620.7910, 1625.7720, 1646.6860, 1660.7235, 1703.8285, 1730.8471, 1759.8666, 1773.9937, 1811.9185, 1851.9095, 1883.8668, 1895.9519, 1925.8401, 1963.8930, 1966.9351, 2001.9016, 2070.9587, 2081.0165, 2084.9664, 2110.0527, 2154.9435, 2185.0887, 2203.0696, 2212.0231, 2225.0772, 2232.2346, 2246.2482, 2265.0848, 2377.1333, 2563.3055



Protein View: A1AT_HUMAN

Alpha-1-antitrypsin OS=Homo sapiens OX=9606 GN=SERPINA1 PE=1 SV=3

Database: SwissProt
Score: 98
Expect: 3.1e-06
Monoisotopic mass (M_r): 46878
Calculated pI: 5.37
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of A1AT_HUMAN against nr.](#)

Search parameters

Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)
Mass values searched: 41
Mass values matched: 10

Protein sequence coverage: 27%

Matched peptides shown in **bold red**.

```

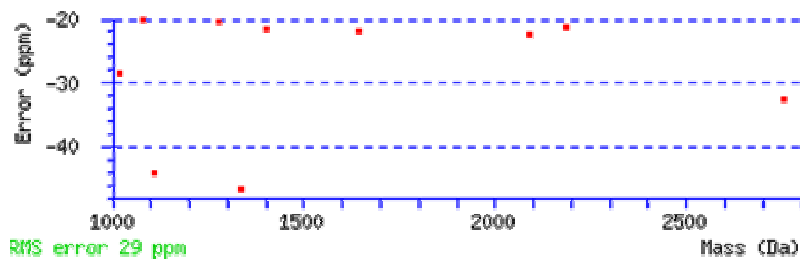
1  MPSSVSWGIL  LLAGLCLLVP  VSLAEDPQGD  AAQKTDTSHH  DQDHPTFNKI
51  TPNLAEFAFS  LYRQLAHQSN  STNIFFSPVS  IATAFAMLSL  GTKADTHDEI
101  LEGLNFNLTE  IPEAQIHEGF  QELLRTLNPQ  DSQQLTGTGN  GLFLSEGLKL
151  VDKFLEDVKK  LYHSEAFTVN  FGDTEEAKQ  INDYVEKGTQ  GKIVDLVKEL
201  DRDTVFALVN  YIFFKGKWER  PFEVKDTEEE  DFHVDQVTTV  KVPMMKRLGM
251  FNIQHCKKLS  SWVLLMKYLG  NATAIFFLPD  EGKLQHLENE  LTHDIITKFL
301  ENEDRRSASL  HLPKLSITGT  YDLKSVLGQL  GITKVFSNGA  DLSGVTEEAP
351  LKLSKAVHKA  VLTIDEKGTE  AAGAMFLEAI  PMSIPPEVKF  NKPFVFLMIE
401  QNTKSPLFMG  KVVNPTQK
  
```

Unformatted sequence string: **418 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Peptide
50 - 63	1641.8278	1640.8205	1640.8562	-21.8	0	K. ITPNLAEF AFS LYR .Q
150 - 160	1333.7102	1332.7029	1332.7653	-46.8	2	K. LVDK FLED VKK .L
161 - 179	2185.9936	2184.9864	2185.0327	-21.2	1	K. LYHSEAF TVN FGDTEEAK .Q
193 - 215	2757.4331	2756.4259	2756.5153	-32.4	2	K. IVDLV KELDRD TVFALVNYIFFK .G
199 - 215	2090.0489	2089.0416	2089.0884	-22.4	1	K. ELDRD TVFALVNYIFFK .G
216 - 225	1275.6584	1274.6511	1274.6771	-20.4	1	K. GKWER PFEVK .D
247 - 257	1403.6731	1402.6658	1402.6962	-21.7	1	K. RLGM FNIQH CK .K
299 - 306	1078.5059	1077.4986	1077.5203	-20.1	1	K. FL ENEDRR .S
315 - 324	1110.5552	1109.5479	1109.5968	-44.1	0	K. LSITGT YDLK .S
325 - 334	1015.5856	1014.5784	1014.6073	-28.6	0	K. SVL GQLGITK .V

No match to: 910.4660, 991.3701, 993.4395, 997.3814, 1019.3428, 1037.5817, 1047.3717, 1049.3650, 1057.3498, 1059.3851, 1132.5594, 1163.4403, 1166.4240, 1216.6149, 1220.4668, 1269.5595, 1307.6426, 1355.7092, 1663.7964, 1801.7064, 1825.8762, 1856.8854, 1877.9080, 1893.8750, 2112.0186, 2207.9732, 2596.2458, 2715.1913, 2731.2280, 3148.3646, 3180.3577



Protein View: APOA4_HUMAN

Apolipoprotein A-IV OS=Homo sapiens OX=9606 GN=APOA4 PE=1 SV=3

Database: SwissProt
Score: 127
Expect: 4.1e-09
Monoisotopic mass (M_r): 45371
Calculated pI: 5.28
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of APOA4 HUMAN against nr.](#)

Search parameters

Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)
Mass values searched: 72
Mass values matched: 17

Protein sequence coverage: 37%

Matched peptides shown in **bold red**.

```

1  MFLKAVVLTLL ALVAVAGARA EVSADQVATV MWDYFSQLSN NAKEAVEHLQ
51  KSELTQQLNA LFQDKLGEVN TYAGDLQKKL VPFATELHER LAKDSEKLKE
101 EIGKELEELR ARLLPHANEV SQKIGDNLRE LQQRLEPYAD QLRTQVNTQA
151 EQLRRQLTPY AQRMERVLRE NADSLQASLR PHADELKAKI DQNVEELKGR
201 LTPYADEFKV KIDQTVEELR RSLPYAQDT QEKLNHQLEG LTFQMKKNAE
251 ELKARISASA EELRQLAPL AEDVRGNLRG NTEGLQKSLA ELGGHLDQOV
301 EEFRRRVEPY GENFNKALVQ QMEQLRQKLG PHAGDVEGHL SFLEKDLRDK
351 VNSFFSTFKE KESQDKTSLSL PELEQQQEQQ QEQQQEQQVQM LAPLES
  
```

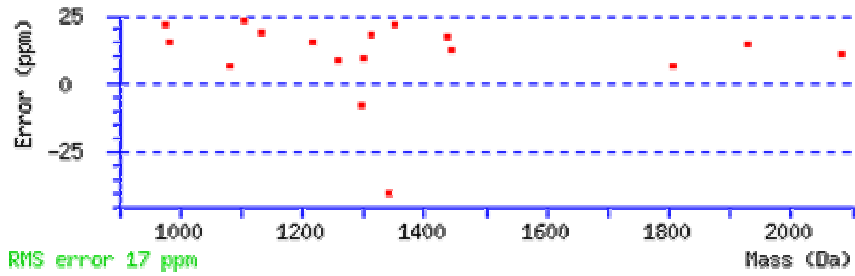
Unformatted sequence string: **396 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Peptide
79 - 90	1439.8258	1438.8185	1438.7932	17.6	1	K.KLV VPFATELHER .L
80 - 90	1311.7298	1310.7225	1310.6983	18.5	0	K.L VPFATELHER .L
124 - 134	1341.6693	1340.6620	1340.7160	-40.3	1	K.I GDNLRELQQR .L
135 - 143	1104.5936	1103.5863	1103.5611	22.8	0	R.L EPYADQLR .T
144 - 155	1443.7846	1442.7773	1442.7590	12.7	1	R.T QVNTQAEQLRR .Q
155 - 163	1132.6439	1131.6366	1131.6149	19.2	1	R.R QLTPYAQR .M
156 - 163	976.5419	975.5346	975.5138	21.4	0	R.Q LTPYAQR .M
190 - 200	1300.6979	1299.6906	1299.6783	9.53	1	K.I DQNVEELKGR .L
199 - 209	1296.6480	1295.6407	1295.6510	-7.95	1	K.G R LTPYADEFK.V
201 - 209	1083.5429	1082.5356	1082.5284	6.67	0	R.L T PYADEFK.V
212 - 221	1258.6863	1257.6790	1257.6677	8.97	1	K.I D QTVEELRR.S
267 - 275	983.5670	982.5597	982.5447	15.2	0	R.L APL AEDVR.G
288 - 304	1927.9782	1926.9709	1926.9435	14.2	0	K.S L AE LGGHLDQQVEEF R.R
288 - 305	2084.0741	2083.0669	2083.0446	10.7	1	K.S L AE LGGHLDQQVEEF RR.R
306 - 316	1352.6888	1351.6815	1351.6520	21.8	1	R.R VEPYGENFNK .A

Start - End	Observed	Mr(expt)	Mr(calc)	ppm	M	Peptide
317 - 326	1215.6702	1214.6629	1214.6441	15.5	0	K.ALVQQMEQLR.Q
329 - 345	1805.9297	1804.9225	1804.9108	6.47	0	K.LGPHAGDVEGHLSFLEK.D

No match to: 805.3172, 806.3257, 810.3573, 812.4041, 820.3092, 826.2984, 836.2974, 842.2797, 844.2874, 848.2890, 864.2745, 879.3837, 881.3443, 887.3934, 889.3875, 896.4195, 920.4506, 923.5308, 942.4642, 945.5263, 959.5158, 980.5180, 993.4652, 997.4929, 1012.5058, 1049.3988, 1057.3844, 1058.3915, 1125.5786, 1160.6169, 1164.5703, 1176.6023, 1192.6034, 1220.5182, 1242.4985, 1257.6756, 1283.5845, 1338.6445, 1374.6673, 1461.7965, 1478.7556, 1495.7034, 1508.7822, 1565.8029, 1585.9006, 1607.8789, 1656.8503, 1721.8523, 1742.8298, 1752.9820, 1773.9928, 1827.9204, 1834.9345, 1849.9194, 2517.1152



Protein View: HPT_HUMAN

Haptoglobin OS=Homo sapiens OX=9606 GN=HP PE=1 SV=1

Database: SwissProt
Score: 109
Expect: 2.6e-07
Monoisotopic mass (M_r): 45861
Calculated pI: 6.13
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of HPT_HUMAN against nr.](#)

Search parameters

Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)
Mass values searched: 85
Mass values matched: 14

Protein sequence coverage: 33%

Matched peptides shown in **bold red**.

```

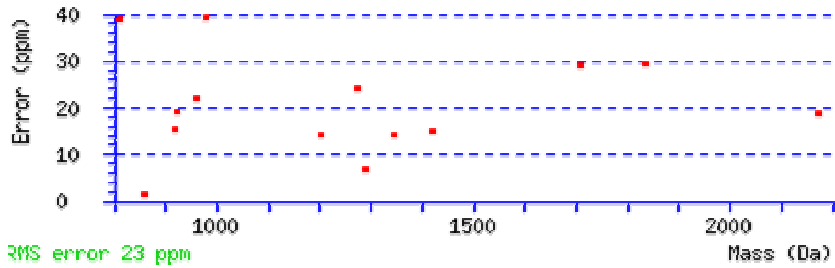
1 MSALGAVIAL LLWGQLFAVD SGNDVTDIAD DGCPKPPEIA HGYVEHSVRY
51 QCKNYKLRRT EGDGVYTLND KKQWINKAVG DKLPECEADD GCPKPPEIAH
101 GYVEHSVRYQ CKNYKLRTE GDGVYTLNNE KQWINKAVGD KLPECEAVCG
151 KPKNPANPVQ RILGGHLDAK GSFPWQAKMV SHHNLTTGAT LINEQWLLTT
201 AKNLFLNHSE NATAKDIAPT LTLVVGKKQL VEIEKVVLHP NYSQVDIGLI
251 KLKQKVSVNE RVMPICLPSK DYAEVGRVGY VSGWGRNANF KFTDHLKYVM
301 LPVADQDQCI RHYEGSTVPE KKTPKSPVGV QPILNEHTFC AGMSKYQEDT
351 CYGDAGSAFA VHDLEEDTWY ATGILSFDKS CAVAEYGVYV KVTSIQDWVQ
401 KTIAEN
  
```

Unformatted sequence string: **406 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Peptide
162 - 170	923.5485	922.5413	922.5236	19.1	0	R. ILGGHLD AK.G
171 - 178	920.4769	919.4696	919.4552	15.7	0	K. GSFPWQ AK.M
216 - 227	1290.7392	1289.7319	1289.7231	6.85	0	K. DIAPTLTLYVGK .K
216 - 228	1418.8465	1417.8392	1417.8181	14.9	1	K. DIAPTLTLYVGKK .Q
229 - 235	858.4943	857.4870	857.4858	1.37	0	K. QLVEIEK .V
254 - 261	959.5478	958.5405	958.5196	21.9	1	K. QKVS VNER.V
262 - 277	1834.9734	1833.9661	1833.9117	29.6	1	R. VMPICLPSKDYAEVGR .V
271 - 277	809.4102	808.4030	808.3715	38.9	0	K. DYAEVGR .V
278 - 286	980.5336	979.5263	979.4876	39.5	0	R. VGYSVSGWGR .N
298 - 311	1707.8687	1706.8614	1706.8120	29.0	0	K. YVMLPVADQDQCI .H
312 - 322	1274.6685	1273.6612	1273.6302	24.3	1	R. HYEGSTVPEK .T
326 - 345	2172.0981	2171.0908	2171.0504	18.6	0	K. SPVGVQPILNEHTFCAGMSK .Y
380 - 391	1345.6648	1344.6575	1344.6384	14.2	0	K. SCAVAEYGVYK .V
392 - 401	1203.6541	1202.6469	1202.6295	14.4	0	K. VTSIQDWVQK .T

No match to: 806.3576, 820.3330, 826.3276, 828.3517, 829.3691, 831.3781, 836.3084, 842.2846, 844.3085, 848.3047, 860.3589, 864.2837, 880.4994, 887.3919, 888.3954, 889.4263, 896.4505, 899.3928, 926.5110, 942.4867, 945.5466, 958.4582, 961.5135, 974.5148, 976.5354, 994.5275, 1002.5139, 1012.5193, 1018.4898, 1057.4071, 1058.4071, 1060.4188, 1066.5704, 1069.4091, 1127.6075, 1160.6310, 1164.6120, 1168.5647, 1176.6222, 1182.6026, 1192.6205, 1215.6439, 1225.6653, 1234.7393, 1241.6298, 1296.6677, 1312.7538, 1328.7313, 1352.6991, 1367.6808, 1381.6896, 1440.8543, 1495.7257, 1532.8276, 1627.8444, 1659.8698, 1673.8749, 1721.8847, 1729.8525, 1737.8567, 1742.8559, 1745.8273, 1753.0072, 1814.9542, 1848.9763, 1856.9516, 2063.9987, 2095.9860, 2127.9858, 2194.1005, 2209.0901



Protein View: HPT_HUMAN

Haptoglobin OS=Homo sapiens OX=9606 GN=HP PE=1 SV=1

Database: SwissProt
Score: 123
Expect: 1e-08
Monoisotopic mass (M_r): 45861
Calculated pI: 6.13
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of HPT_HUMAN against nr.](#)

Search parameters

Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)
Mass values searched: 29
Mass values matched: 13

Protein sequence coverage: 33%

Matched peptides shown in **bold red**.

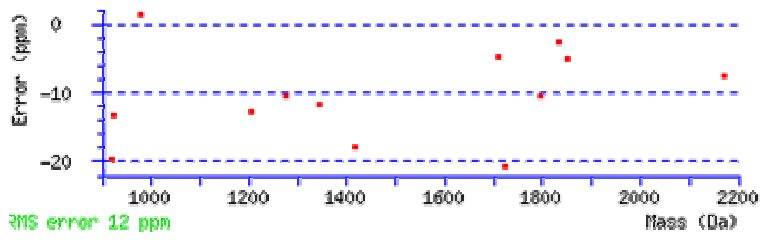
1 MSALGAVIAL LLWQQLFAVD SGNDVTDIAD DGCPKPPEIA HGYVEHSVRY
51 QCKNYYKLRT EGDGVYTLND KKQWINKAVG DKLPECEADD GCPKPPEIAH
101 GYVEHSVRYQ CKNYYKL RTE GDGVYTLNNE KQWINKAVGD KLPECEAVCG
151 KPKNPANPVQ **RILGGHLDK** **GSFPWQAKM**V SHHNLTTGAT LINEQWLLTT
201 AKNLFLNHSE NATAK**DIAPT** **LTLVVGKK**QL VEIEK**VVLHP** **NYSQVDIGLI**
251 **KLKQK**SVNE **RVMPICLPSK** **DYAEVGRVGY** **VSGWGRNANF** KFTDHLK**YVM**
301 **LPVADQDQCI** **RHYEGSTVPE** **KKTPKSPVGV** **QPILNEHTFC** **AGMSKYQEDT**
351 CYGDAGSAFA VHDLEEDTWY ATGILSFDKS **CAVAEYGVYV** **KVTSIQDWVQ**
401 **KTIAEN**

Unformatted sequence string: **406 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
Show matched peptides only predicted peptides also

Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Peptide
162 - 170	923.5187	922.5114	922.5236	-13.2	0	R.ILGGHLDK.G
171 - 178	920.4445	919.4372	919.4552	-19.5	0	K.GSFPWQAK.M
216 - 228	1418.7999	1417.7926	1417.8181	-17.9	1	K.DIAPTLTLVVGKK.Q
236 - 251	1794.9926	1793.9853	1794.0040	-10.4	0	K.VVLHPNYSQVDIGLIK.L
262 - 277	1834.9145	1833.9072	1833.9117	-2.47	1	R.VMPICLPSKDYAEVGR.V
262 - 277	1850.9047	1849.8974	1849.9066	-4.97	1	R.VMPICLPSKDYAEVGR.V + Oxidation (M)
278 - 286	980.4961	979.4888	979.4876	1.30	0	R.VGYVSGWGR.N
298 - 311	1707.8112	1706.8040	1706.8120	-4.72	0	K.YVMLPVADQDQCIR.H
298 - 311	1723.7788	1722.7715	1722.8069	-20.6	0	K.YVMLPVADQDQCIR.H + Oxidation (M)
312 - 322	1274.6241	1273.6169	1273.6302	-10.5	1	R.HYEGSTVPEKK.T
326 - 345	2172.0412	2171.0339	2171.0504	-7.59	0	K.SPVGVPILNEHTFCAGMSK.Y
380 - 391	1345.6297	1344.6224	1344.6384	-11.9	0	K.SCAVAEYGVYV.V
392 - 401	1203.6215	1202.6142	1202.6295	-12.8	0	K.VTSIQDWVQK.T

No match to: 1012.4791, 1160.5889, 1164.5811, 1192.5747, 1234.6895, 1495.6757, 1509.6839, 1685.7507, 1721.8256, 1737.8151, 1742.7969, 1752.9457, 1774.9204, 1816.9740, 1845.0189, 1874.8669



Protein View: APOA1_HUMAN

Apolipoprotein A-I OS=Homo sapiens OX=9606 GN=APOA1 PE=1 SV=1

Database: SwissProt
Score: 191
Expect: 1.6e-15
Monoisotopic mass (M_r): 30759
Calculated pI: 5.56
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of APOA1_HUMAN against nr.](#)

Search parameters

Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)
Mass values searched: 98
Mass values matched: 23

Protein sequence coverage: 71%

Matched peptides shown in **bold red**.

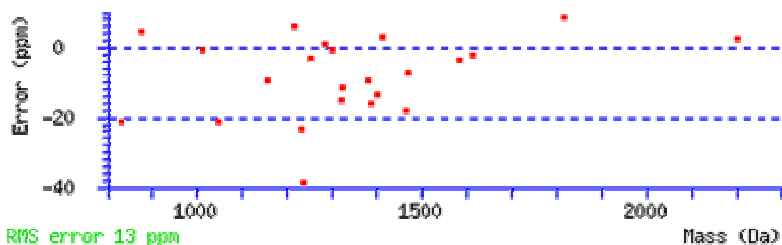
1 MKAAVLTLLAV LFLTGSQARH FWQQDEPPQS PWDR**VKDLAT VYVDVLKDSG**
 51 **RDYVSQFEFS ALGKQLNLKL LDNWDSVTST FSKLREQQLGP VTQEFWDNLE**
 101 **KETEGLRQEM SKDLEEVKAK VQPYLDDFQK KWQEEMELYR QKVEPLRAEL**
 151 **QEGARQKLHE LQEKLSPLGE EMRDRARAHV DALR**THLAPY SDELQR**LAA**
 201 **RLEALKENGG ARLAEYHAKA TEHLSTLSEK AKPALEDLRQ GLLPVLESFK**
 251 **VSFLSALEY TKKLNTQ**

Unformatted sequence string: **267 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Peptide
35 - 47	1462.8248	1461.8175	1461.8443	-18.3	1	R.VKDLATVYVDVLK.D
37 - 47	1235.6406	1234.6334	1234.6809	-38.5	0	K.DLATVYVDVLK.D
48 - 64	1815.8664	1814.8591	1814.8435	8.60	1	K.DSGRDYVSQFEFSALGK.Q
52 - 64	1400.6506	1399.6433	1399.6620	-13.3	0	R.DYVSQFEFSALGK.Q
70 - 83	1612.7816	1611.7743	1611.7781	-2.31	0	K.LLDNWDSVTSTFSK.L
84 - 101	2202.1243	2201.1170	2201.1117	2.42	1	K.LREQQLGPVTQEFWDNLEK.E
102 - 112	1323.6061	1322.5988	1322.6136	-11.2	1	K.ETEGLRQEMSK.D + Oxidation (M)
121 - 130	1252.6167	1251.6094	1251.6136	-3.28	0	K.VQPYLDDFQK.K
121 - 131	1380.7031	1379.6959	1379.7085	-9.16	1	K.VQPYLDDFQKK.W
131 - 140	1411.6716	1410.6643	1410.6601	2.93	1	K.KWQEEMELYR.Q
132 - 140	1283.5738	1282.5665	1282.5652	1.05	0	K.WQEEMELYR.Q
143 - 155	1467.7810	1466.7737	1466.7841	-7.06	1	K.VEPLRAELQEGAR.Q
148 - 155	873.4464	872.4392	872.4352	4.59	0	R.AELQEGAR.Q
165 - 173	1047.4918	1046.4845	1046.5066	-21.1	0	K.LSPLGEEEMR.D + Oxidation (M)
165 - 175	1318.6222	1317.6149	1317.6347	-15.0	1	K.LSPLGEEEMRDR.A + Oxidation (M)
185 - 195	1301.6477	1300.6404	1300.6411	-0.57	0	R.THLAPYSDELQR.Q
185 - 197	1585.8026	1584.7954	1584.8008	-3.45	1	R.THLAPYSDELQRQ.L
202 - 212	1157.6162	1156.6089	1156.6200	-9.56	1	R.LEALKENGGAR.L
213 - 219	831.4185	830.4112	830.4286	-21.0	0	R.LAEYHAK.A
220 - 230	1215.6289	1214.6216	1214.6143	6.04	0	K.ATEHLSTLSEK.A
231 - 239	1012.5782	1011.5709	1011.5713	-0.31	0	K.AKPALEDLR.Q
240 - 250	1230.6805	1229.6732	1229.7020	-23.4	0	R.QGLLPVLESFK.V
251 - 262	1386.6931	1385.6858	1385.7078	-15.9	0	K.VSFLSALEYTK.K

No match to: 806.3176, 820.3073, 826.2799, 842.2619, 855.5099, 859.3318, 861.3599, 869.3898, 879.3716, 887.3625, 889.3674, 979.5530, 993.5614, 1007.5780, 1057.3568, 1059.3701, 1158.5947, 1219.6572, 1226.5387, 1266.5546, 1268.6157, 1315.5574, 1405.6679, 1408.6805, 1484.8199, 1495.6729, 1580.8064, 1602.7801, 1632.7889, 1645.7979, 1653.7629, 1661.9146, 1675.9292, 1689.8772, 1705.8905, 1723.9390, 1733.9008, 1743.8551, 1765.8481, 1775.8486, 1797.9000, 1811.9147, 1819.9022, 1870.9907, 1878.9977, 1884.0078, 1946.0107, 1968.0051, 2037.9130, 2051.9329, 2065.0389, 2084.8869, 2094.9423, 2102.1329, 2108.9815, 2137.1176, 2141.0338, 2169.1222, 2206.1028, 2234.1261, 2303.1680, 2309.0439, 2323.0749, 2352.1872, 2380.2003, 2409.2121, 2423.2270, 2439.2305, 2450.2495, 2453.2422, 2466.2508, 2481.2564, 2494.2630, 2624.3350, 2677.3036



RMS error 13 ppm

Mass (Da)

Protein View: RET4_HUMAN

Retinol-binding protein 4 OS=Homo sapiens OX=9606 GN=RBP4 PE=1 SV=3

Database: SwissProt
Score: 70
Expect: 0.0019
Monoisotopic mass (M_r): 23337
Calculated pI: 5.76
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of RET4_HUMAN against nr.](#)

Search parameters

Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)
Mass values searched: 13
Mass values matched: 5

Protein sequence coverage: 33%

Matched peptides shown in **bold red**.

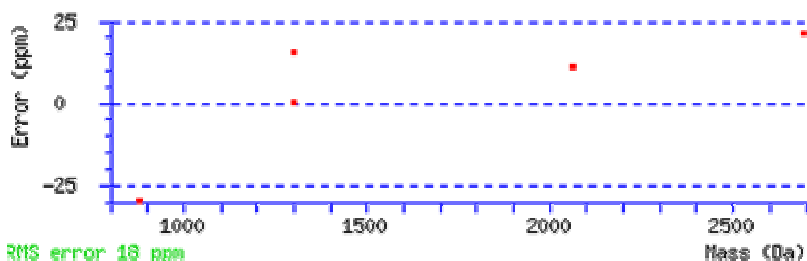
1 MKVWVALLLL AALGSGRAER DCRVSSFRVK **ENFDKAR**FSG TWYAMAKKDP
 51 EGLFLQDNIV AEFSVDETQQ MSATAKGRVR LLNNWDVCAD MVTFTDTE
 101 PAKFKMKYWG VASFLQ**GND DHWIVD**TDYD **TYAVQY**SCRL **LNL**DGTCADS
 151 **YSEVFSR**DPN GLPPEAQKIV **RQRQEEL**CLA **RQYRLIV**HNG **YCDGR**SERNL
 201 L

Unformatted sequence string: **201 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Peptide
31 - 37	879.4060	878.3987	878.4246	-29.5	1	K.ENFDKAR.F
118 - 139	2693.1786	2692.1713	2692.1136	21.4	0	K.GNDDHWIVD TDYD TYAVQY SCR.L
140 - 157	2064.9919	2063.9847	2063.9623	10.9	0	R.LLNL DGTCAD SYSFVFSR .D
172 - 181	1302.6585	1301.6513	1301.6510	0.21	1	R.QRQEEL CLAR.Q
185 - 195	1303.6413	1302.6341	1302.6139	15.5	0	R.LIVHNG YCDGR.S

No match to: 803.3802, 887.3827, 895.3929, 2676.1381, 2707.1841, 2715.1574, 2725.1726, 2845.1762



Protein View: HPT_HUMAN

Haptoglobin OS=Homo sapiens OX=9606 GN=HP PE=1 SV=1

Database: SwissProt
Score: 62
Expect: 0.011
Monoisotopic mass (M_r): 45861
Calculated pI: 6.13
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of HPT_HUMAN against nr.](#)

Search parameters

Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)
Mass values searched: 52
Mass values matched: 8

Protein sequence coverage: 21%

Matched peptides shown in **bold red**.

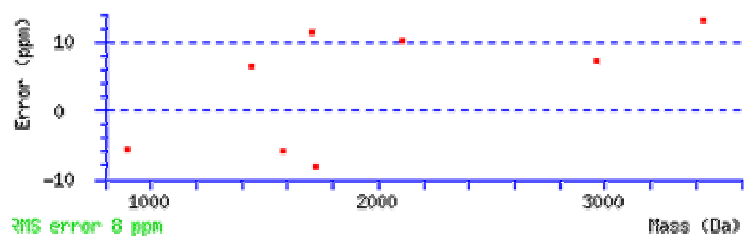
1 MSALGAVIAL LLWQQLFAVD SGNDVTDIAD DGCPKPPEIA HGYVEHSVRY
51 QCKNYYKLRT **EGDGVYTLND KKQWINKAVG DKLPECEADD GCPKPPEIAH**
101 **GYVEHSVRYQ** CKNYYKLRTE **GDGVYTLNNE KQWINKAVGD** KLPECEAVCG
151 **KPKNPANPVQ** RILGGHLDK GSFPWQAKMV SHHNLTTGAT LINEQWLLTT
201 AKNLFLNHSE NATAKDIAPT LTLYVGKKQL VEIEKVVLHP NYSQVDIGLI
251 **KLKQKVSUNE** RVMPICLPSK **DYAEVGRVGY** VSGWGRNANF **KFTDHLKYVM**
301 **LPVADQDQCI** RHYEGSTVPE **KKTPKSPVGV** QPILNEHTFC AGMSKYQEDT
351 CYGDAGSAFA VHDLEEDTWY ATGILSFDKS CAVAEGVYV KVTSIQDWVQ
401 KTIAEN

Unformatted sequence string: **406 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
Show matched peptides only predicted peptides also

Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Peptide
58 - 71	1580.7821	1579.7748	1579.7842	-5.96	1	K.LRTEGDGVYTLNDK.K
78 - 108	3432.6397	3431.6325	3431.5874	13.1	1	K.AVGDKLPECEADDGCPKPPEIAHGYVEHSVR.Y
83 - 108	2962.3669	2961.3596	2961.3385	7.12	0	K.LPECEADDGCPKPPEIAHGYVEHSVR.Y
117 - 131	1708.8695	1707.8622	1707.8428	11.4	1	K.LRTEGDGVYTLNNEK.Q
119 - 131	1439.6740	1438.6667	1438.6576	6.35	0	R.TEGDGVYTLNNEK.Q
119 - 136	2109.0464	2108.0391	2108.0174	10.3	1	R.TEGDGVYTLNNEKQWINK.A
154 - 161	895.4694	894.4621	894.4671	-5.66	0	K.NPANPVQR.I
298 - 311	1723.8001	1722.7928	1722.8069	-8.21	0	K.YVMLPVADQDQCIR.H + Oxidation (M)

No match to: 897.3565, 899.3691, 901.3695, 903.3802, 993.4297, 1057.3665, 1058.3675, 1060.3780, 1069.3631, 1106.5854, 1161.5031, 1163.4464, 1198.6184, 1220.4959, 1230.5978, 1285.6265, 1303.6282, 1304.5689, 1317.6233, 1347.5372, 1505.7258, 1675.7664, 1682.7634, 1717.2898, 1730.8297, 2064.9828, 2078.9959, 2378.2329, 2410.2200, 2636.1222, 2649.1358, 2677.1207, 2679.8777, 2693.1589, 2707.1545, 2716.1171, 2725.1448, 2739.1459, 2976.3878, 3349.5677, 3363.5906, 3417.5049, 3446.6574, 3460.6543



Protein View: HPT_HUMAN

Haptoglobin OS=Homo sapiens OX=9606 GN=HP PE=1 SV=1

Database: SwissProt
Score: 60
Expect: 0.022
Monoisotopic mass (M_r): 45861
Calculated pI: 6.13
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of HPT_HUMAN against nr.](#)

Search parameters

Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)
Mass values searched: 38
Mass values matched: 6

Protein sequence coverage: 16%

Matched peptides shown in **bold red**.

```

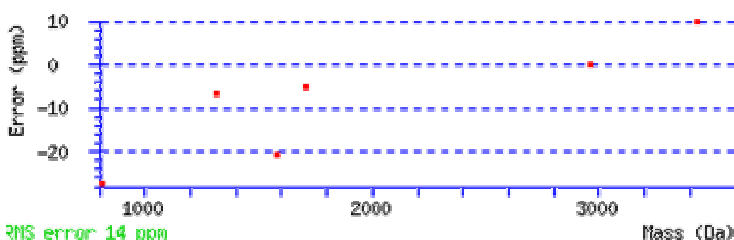
1 MSALGAVIAL LLWQQLFAVD SGNDVTDIAD DGCPKPPEIA HGYVEHSVRY
51 QCKNYYKLRT EGDGVYTLND KKQWINKAVG DKLPECEADD GCPKPPEIAH
101 GYVEHSVRYQ CKNYYKLRTE GDGVYTLNNE KQWINKAVGD KLPECEAVCG
151 KPKNPANPVQ RILGGHLDK GSFQWQAKMV SHHNLTTGAT LINEQWLLTT
201 AKNLFLNHSE NATAKDIAPT LTLVYVGGKQL VEIEKVVLP NYSQVDIGLI
251 KLKQKVSUNE RVMPICLPSK DYA EVGRVGY VSGWGRNANF KFTDHLKYVM
301 LPVADQDQCI RHYEGSTVPE KKTPKSPVGV QPILNEHTFC AGMSKYQEDT
351 CYGDAGSAFA VHDLEEDTWY ATGILSFDKS CAVA EYGVYV KVTSIQDWVQ
401 KTIAEN
  
```

Unformatted sequence string: **406 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Peptide
58 - 71	1580.7589	1579.7516	1579.7842	-20.6	1	K.LRTEGDGVYTLNDK.K
60 - 71	1311.5977	1310.5905	1310.5990	-6.54	0	R.TEGDGVYTLNDK.K
78 - 108	3432.6286	3431.6213	3431.5874	9.88	1	K.AVGDKLPECEADDGCPKPPEIAHGYVEHSVR.Y
83 - 108	2962.3461	2961.3389	2961.3385	0.11	0	K.LPECEADDGCPKPPEIAHGYVEHSVR.Y
117 - 131	1708.8414	1707.8342	1707.8428	-5.03	1	K.LRTEGDGVYTLNNEK.Q
271 - 277	809.3566	808.3493	808.3715	-27.5	0	K.DYAEVGR.V

No match to: 800.3451, 801.3705, 803.3641, 805.3823, 807.3957, 819.3379, 825.3708, 830.3505, 845.3537, 852.3576, 879.3788, 887.3705, 889.3759, 895.3855, 899.3863, 919.4034, 1057.3584, 1058.3607, 1267.6405, 1289.6242, 1684.8375, 1730.8151, 1964.9382, 2378.1887, 2394.1799, 2415.1472, 2976.3607, 3349.5558, 3361.0096, 3363.5637, 3403.5543, 3446.6305



Protein View: TTHY_HUMAN

Transthyretin OS=Homo sapiens OX=9606 GN=TTR PE=1 SV=1

Database: SwissProt
Score: 64
Expect: 0.0081
Monoisotopic mass (M_r): 15991
Calculated pI: 5.52
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of TTHY_HUMAN against nr.](#)

Search parameters

Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)
Mass values searched: 65
Mass values matched: 6

Protein sequence coverage: 61%

Matched peptides shown in **bold red**.

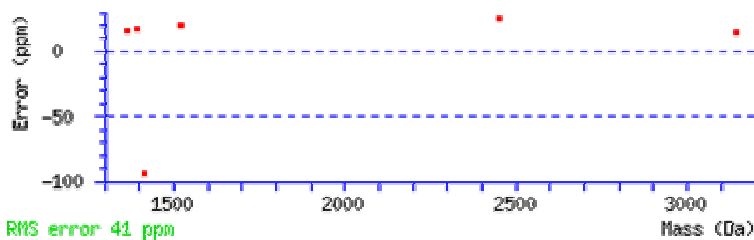
1 MASHRLLLLC LAGLVFVSEA GPTGTGESKC **PLMVKVLDAV RGSPAINVAV**
 51 **HVFRKAADDT WEPFASGKTS ESGELHGLTT EEEFVEGIYK VEIDTKSYWK**
 101 **ALGISPFHEH AEVVFTANDS GPRRYTIAAL LSPYSYSTTA VVTNPKE**

Unformatted sequence string: **147 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Peptide
30 - 41	1416.6379	1415.6306	1415.7629	-93.4	1	K.CPLMVKVLDAVR.G + Oxidation (M)
42 - 54	1366.7816	1365.7743	1365.7517	16.5	0	R.GSPAINVAVHVFR.K
55 - 68	1522.7470	1521.7397	1521.7100	19.6	1	R.KAADDTWEPFASGK.T
56 - 68	1394.6468	1393.6395	1393.6150	17.6	0	K.AADDTWEPFASGK.T
69 - 96	3140.5632	3139.5559	3139.5085	15.1	1	K.TSESGELHGLTTEEEFVEGIYKVEIDTK.S
101 - 123	2451.2658	2450.2586	2450.1979	24.8	0	K.ALGISPFHEHAEEVFTANDSGPR.R

No match to: 820.3859, 822.3767, 824.3827, 829.4568, 833.3989, 838.4381, 846.3893, 856.5263, 864.4590, 870.5501, 895.4615, 897.4285, 908.4467, 929.4930, 934.4477, 951.4799, 962.4752, 982.4537, 993.5146, 1002.5493, 1016.5464, 1033.5290, 1037.5285, 1046.5618, 1060.5584, 1065.5226, 1068.5239, 1107.5627, 1118.5271, 1121.5817, 1131.6564, 1157.5929, 1189.6149, 1194.6238, 1229.6773, 1234.6904, 1265.6520, 1284.5922, 1307.7057, 1311.6276, 1317.6457, 1357.7376, 1411.7805, 1432.6300, 1434.7149, 1439.7327, 1450.7778, 1464.7760, 1537.7588, 1580.8267, 1608.7958, 1612.8750, 1708.9214, 1751.8887, 1791.7799, 1794.8374, 1851.9701, 1857.9608, 2065.9970



Protein View: HPT_HUMAN

Haptoglobin OS=Homo sapiens GN=HP PE=1 SV=1

Database: SwissProt
 Score: 64
 Expect: 0.008
 Nominal mass (M_r): 45861
 Calculated pI: 6.13
 Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of HPT_HUMAN against nr.](#)

Search parameters

Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
 Fixed modifications: Carbamidomethyl (C)
 Variable modifications: Oxidation (M)
 Mass values searched: 81
 Mass values matched: 10

Protein sequence coverage: 19%

Matched peptides shown in **bold red**.

```

1 MSALGAVIAL LLWQQLFAVD SGNDVTDIAD DGCPKPPEIA HGYVEHSVRY
51 QCKNYYKLRTE EGDGVYTLND KKQWINKAVG DKLPECEADD GCPKPPEIAH
101 GYVEHSVRYQ CKNYYKL RTE GDGVYTLNNE KQWINKAVGD KLPECEAVCG
151 KPKNPANPVQ RILGGHLDK GSFPWQAKMV SHHNLTTGAT LINEQWLLTT
201 AKNLFLNHSE NATAKDIAPT LTLYVGKKQL VEIEKVVLHP NYSQVDIGLI
251 KLKQKVSUNE RVMPICLPSK DYAEVGRVGY VSGWGRNANF KFTDHLKYVM
301 LPVADQDQCI RHYEGSTVPE KKTPKSPVGV QPILNEHTFC AGMSKYQEDT
351 CYGDAGSAFA VHDLEEDTWY ATGILSFDKS CAVA EYGVYV KVTSIQDWVQ
401 KTIAEN
  
```

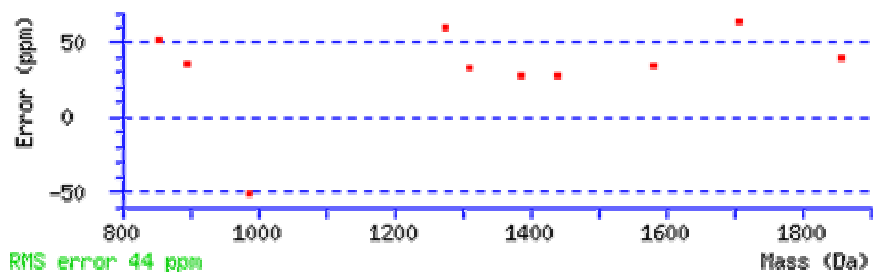
Unformatted sequence string: **406 residues** (for pasting into other applications).

Sort peptides by Residue Number Increasing Mass Decreasing Mass

Show predicted peptides also

Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Peptide
58 - 71	1580.8454	1579.8381	1579.7842	34.1	1	K.LRTEGDGVYTLNDK.K
60 - 71	1311.6498	1310.6425	1310.5990	33.2	0	R.TEGDGVYTLNDK.K
60 - 72	1439.7408	1438.7335	1438.6940	27.5	1	R.TEGDGVYTLNDKK.Q
113 - 118	856.5123	855.5050	855.4603	52.3	1	K.NYYKL.R.T
117 - 131	1708.9585	1707.9512	1707.8428	63.5	1	K.LRTEGDGVYTLNNEK.Q
137 - 153	1857.9926	1856.9853	1856.9124	39.2	1	K.AVGDKLPECEAVCGKPK.N
142 - 153	1387.7088	1386.7015	1386.6635	27.4	0	K.LPECEAVCGKPK.N
154 - 161	895.5063	894.4990	894.4671	35.7	0	K.NPANPVQR.I
228 - 235	986.5370	985.5298	985.5808	-51.8	1	K.KQLVEIEK.V
312 - 322	1274.7130	1273.7058	1273.6302	59.3	1	R.HYEGSTVPEKK.T

No match to: 820.3984, 822.4516, 824.4226, 830.4521, 833.4226, 836.4647,
838.4726, 839.3575, 848.4429, 869.3826, 870.5281, 873.4573, 880.4565, 881.4837,
887.4361, 899.4657, 913.5113, 917.4760, 921.5032, 927.5120, 929.5450, 945.4888,
946.5040, 964.5340, 968.5782, 998.5432, 1033.5704, 1108.5862, 1117.5409,
1132.5793, 1138.5461, 1149.6068, 1152.5638, 1205.5870, 1286.7040, 1287.6882,
1313.7879, 1314.7267, 1318.6225, 1331.7396, 1333.6300, 1366.7941, 1379.2050,
1394.6691, 1461.7405, 1497.8438, 1522.7508, 1529.7794, 1545.7805, 1604.9875,
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1911.0051, 2018.5006, 2262.1784, 2276.0221, 2355.5774, 2369.8066, 2451.3325,
2516.5968, 2615.8135, 2630.0413, 3174.5680, 3334.7752, 3349.7638, 3363.7276



Protein View: SAA1_HUMAN

Serum amyloid A-1 protein OS=Homo sapiens OX=9606 GN=SAA1 PE=1 SV=1

Database: SwissProt
 Score: 63
 Expect: 0.01
 Monoisotopic mass (M_r): 13581
 Calculated pI: 6.28
 Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of SAA1_HUMAN against nr.](#)

Search parameters

Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
 Fixed modifications: Carbamidomethyl (C)
 Variable modifications: Oxidation (M)
 Mass values searched: 39
 Mass values matched: 5

Protein sequence coverage: 55%

Matched peptides shown in **bold red**.

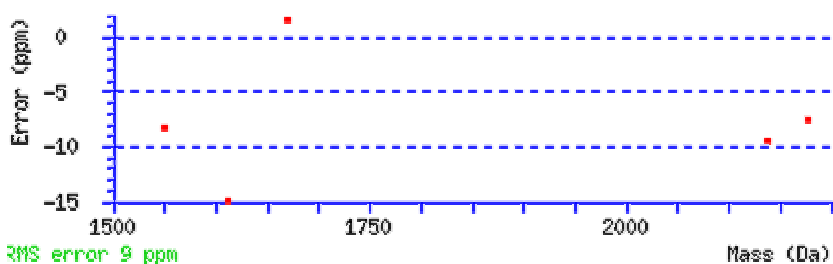
1 MKLLTGLVFC SLVLGVSSRS **FFSFLGEAFD GARDMWRAYS** DMREANYIGS
 51 **DKYFHARGNY** DAAK**RGPGGV WAAEAISDAR** ENIQ**RFFGHG AEDSLADQAA**
 101 **NEWGR**SGKDP NHFRPAGLPE KY

Unformatted sequence string: **122 residues** (for pasting into other applications).

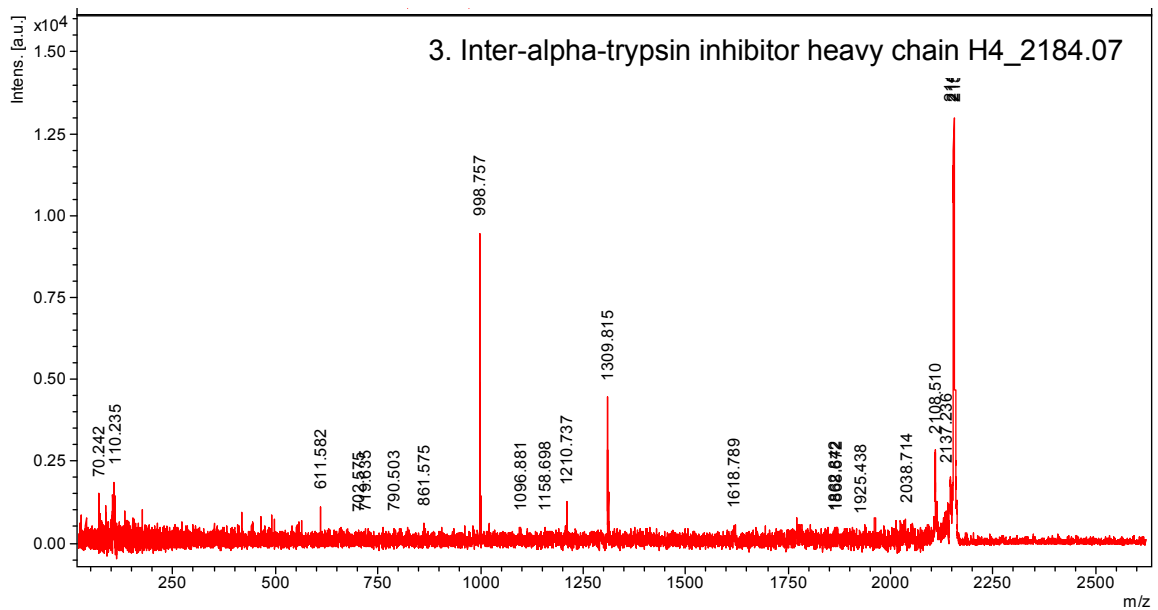
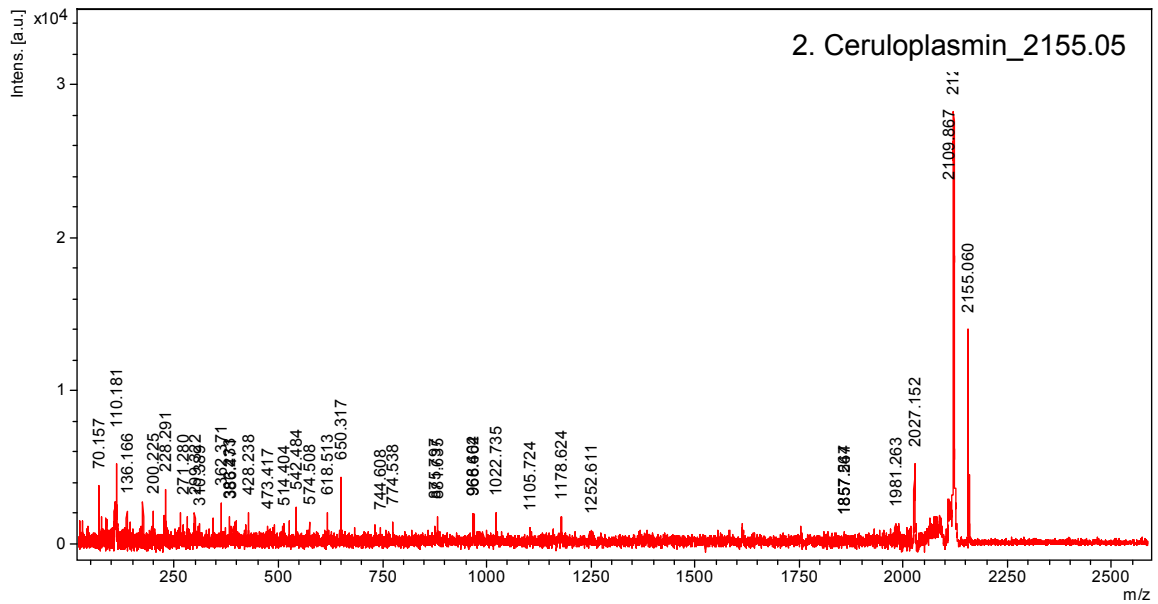
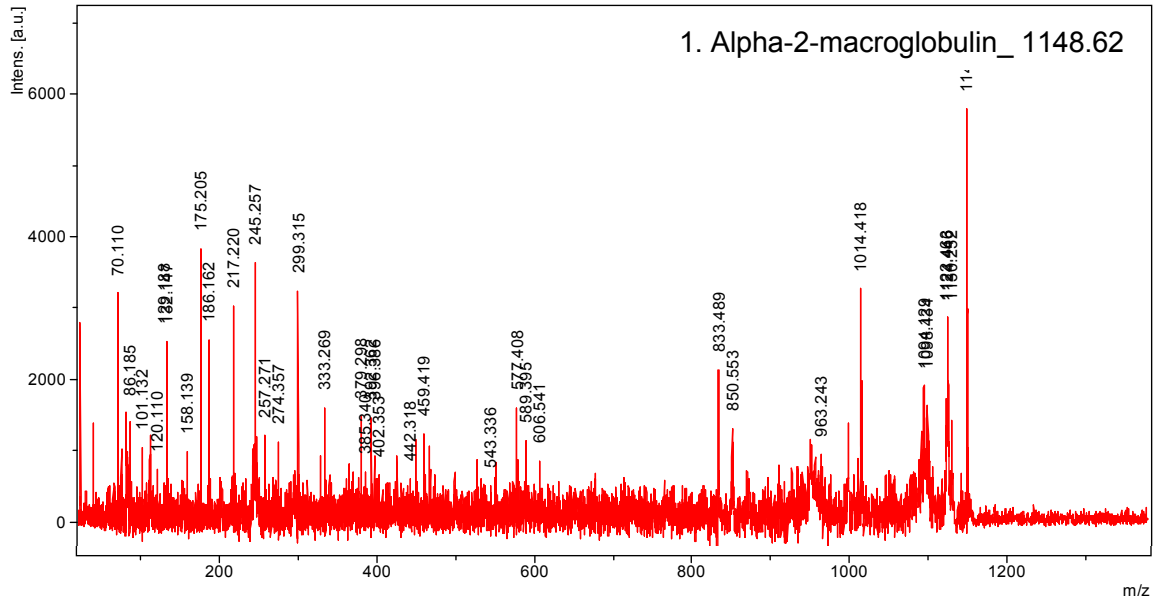
Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

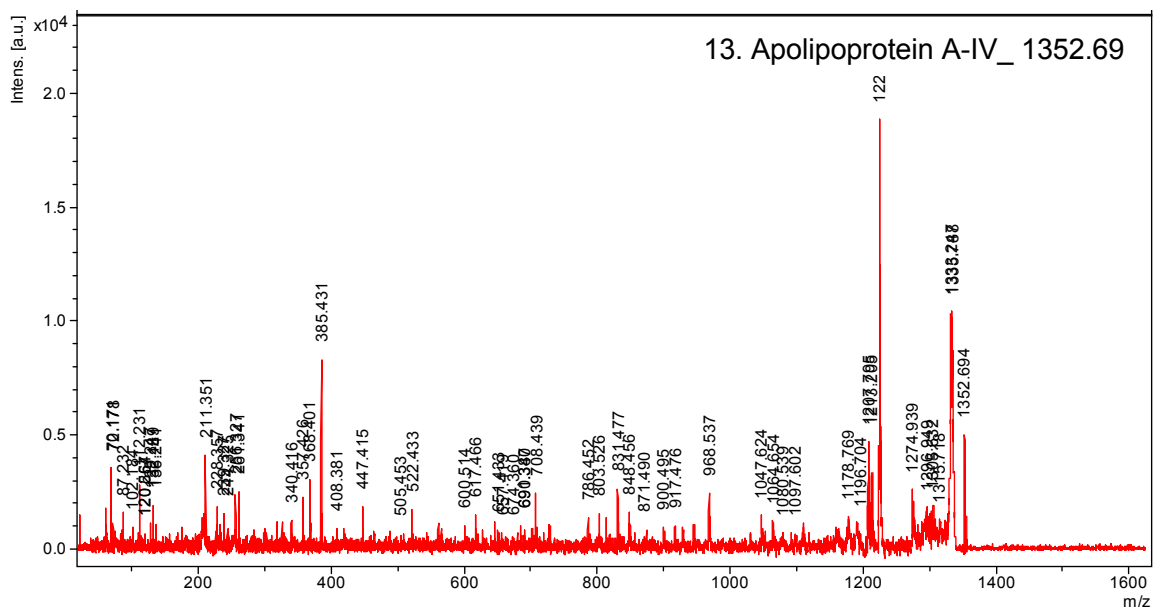
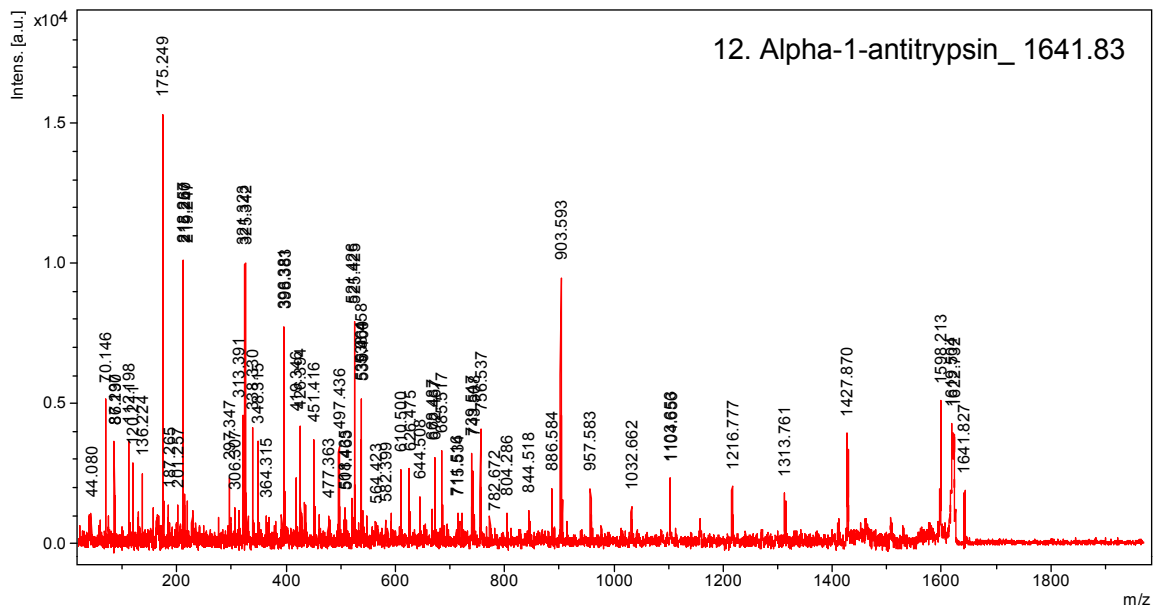
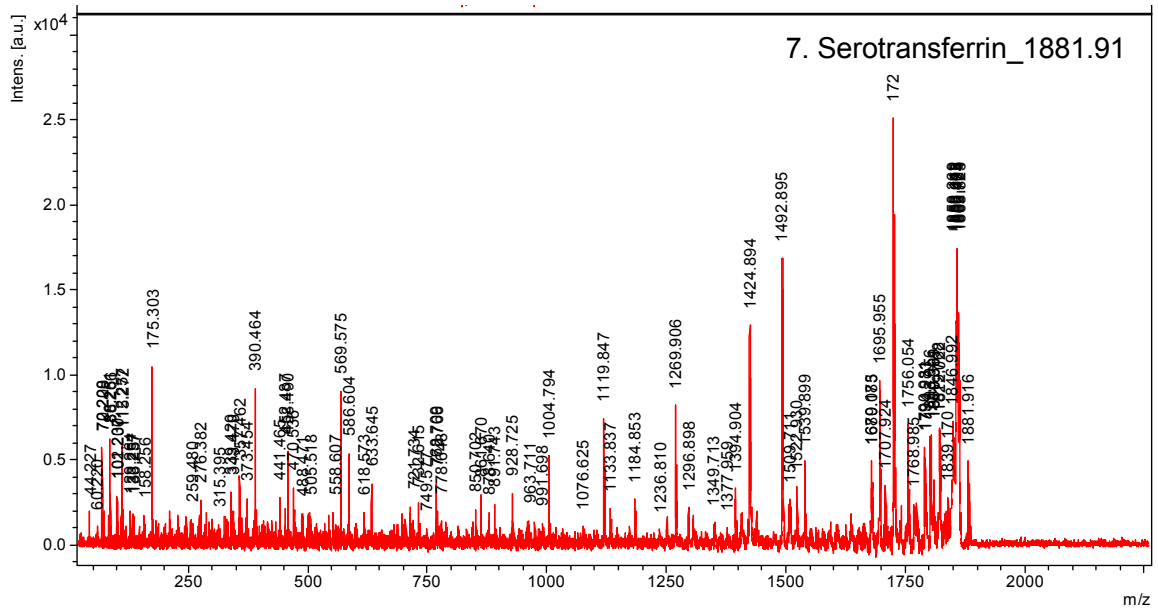
Start - End	Observed	Mr(expt)	Mr(calc)	ppm	M	Peptide
20 - 33	1550.7146	1549.7073	1549.7202	-8.28	0	R.SFFSFLGEAFDGAR.D
20 - 37	2138.9551	2137.9478	2137.9680	-9.45	1	R.SFFSFLGEAFDGARDMWR.A
44 - 57	1670.7946	1669.7874	1669.7848	1.52	1	R.EANYIGSDKYFHAR.G
65 - 80	1612.7949	1611.7876	1611.8117	-14.9	1	K.RGPGGVWAAEAISDAR.E
86 - 105	2177.9472	2176.9400	2176.9562	-7.48	0	R.FFGHGAEDSLADQAANEWGR.S

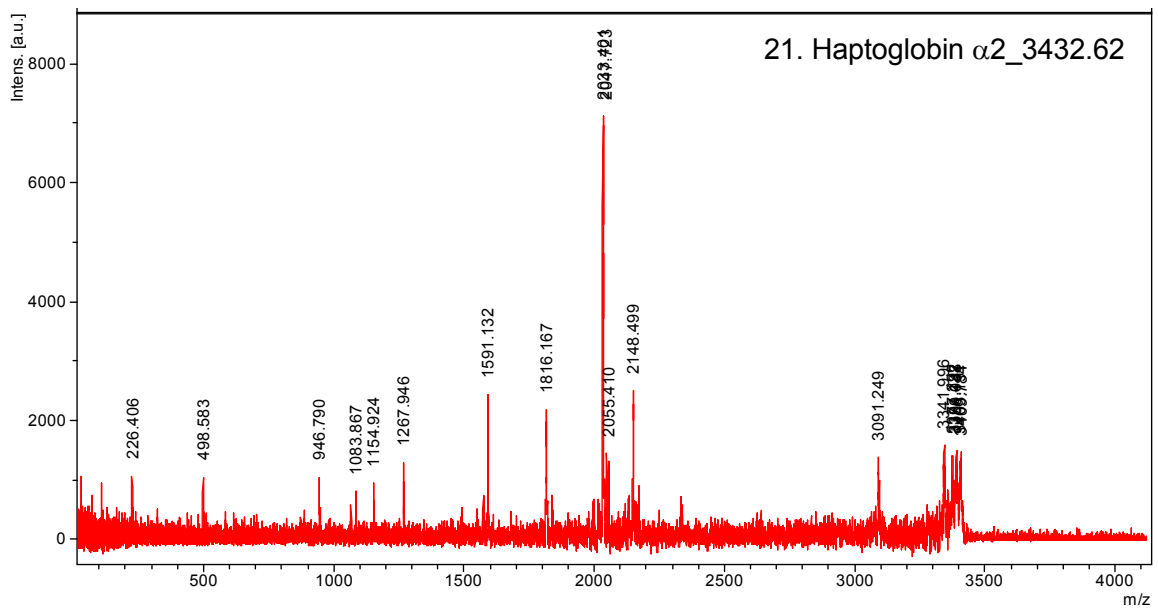
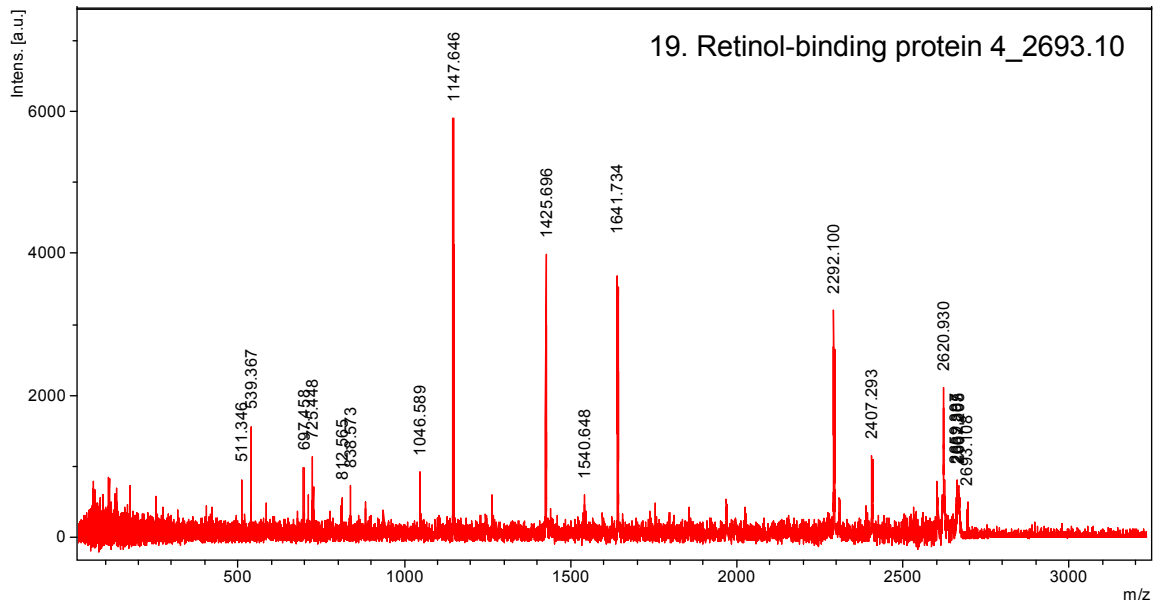
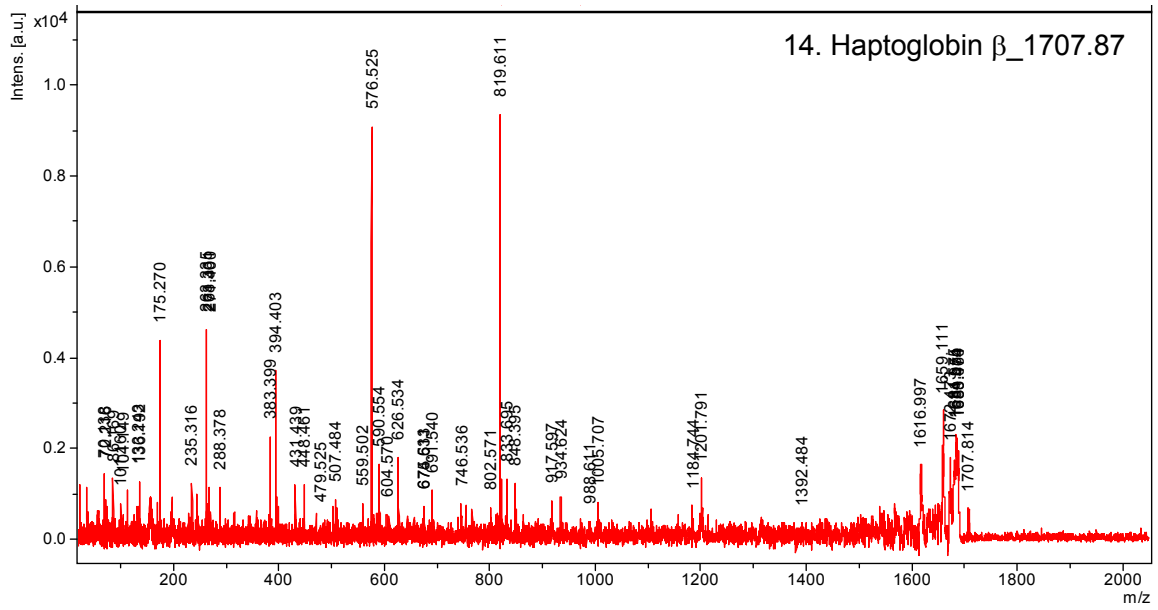
No match to: 915.4331, 919.4010, 1057.3753, 1274.7195, 1290.6976, 1400.6719, 1601.5489, 1639.8174, 1684.8545, 1706.8060, 1791.7161, 1848.8244, 1864.8034, 1892.9673, 1897.8046, 1912.9490, 1934.9070, 1940.9421, 2090.9553, 2102.2195, 2104.9846, 2116.2476, 2153.9570, 2155.9506, 2168.9625, 2185.9487, 2193.9496, 2200.9420, 2209.9841, 2318.9566, 2342.9688, 2705.1554, 2717.1025, 3312.2954

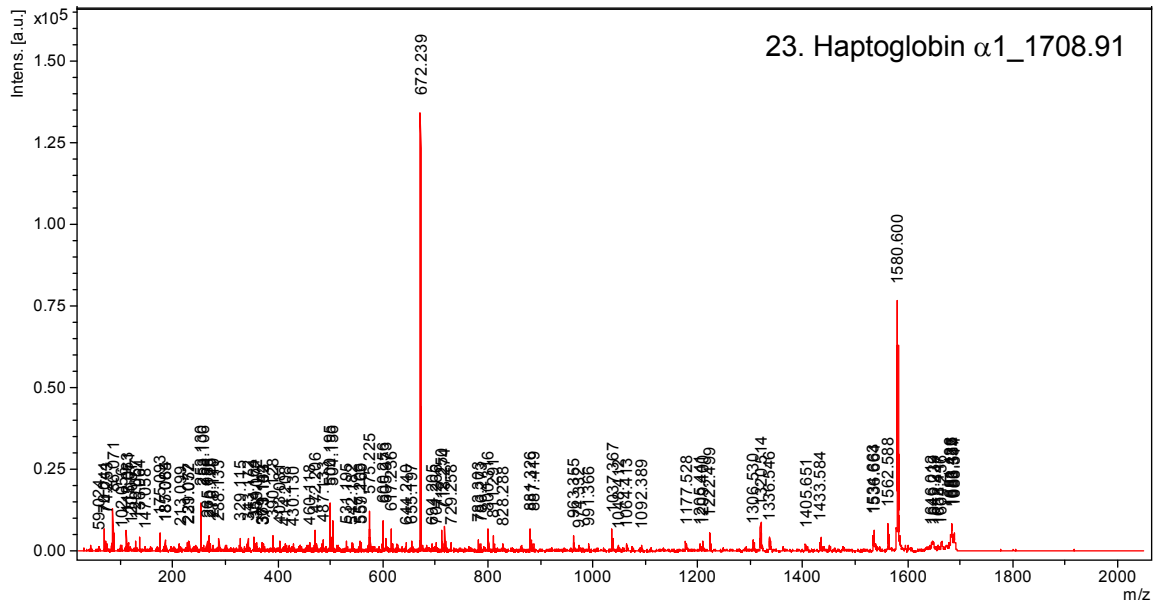
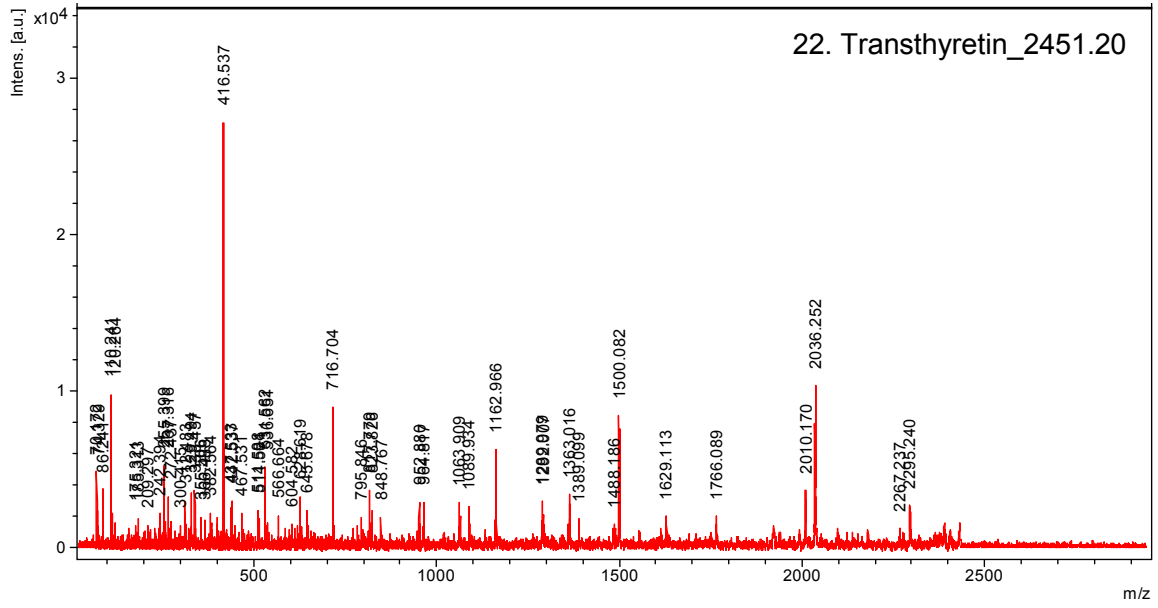


17. MALDI LIF-TOF/TOF MS spectra and Mascot searches









Protein View: A2MG_HUMAN

Alpha-2-macroglobulin OS=Homo sapiens OX=9606 GN=A2M PE=1 SV=3

Database: SwissProt
Score: 63
Monoisotopic mass (M_r): 164613
Calculated pI: 6.03
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of A2MG_HUMAN against nr.](#)

Search parameters

MS data file: DATA.TXT
Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)

Protein sequence coverage: 0%

Matched peptides shown in **bold red**.

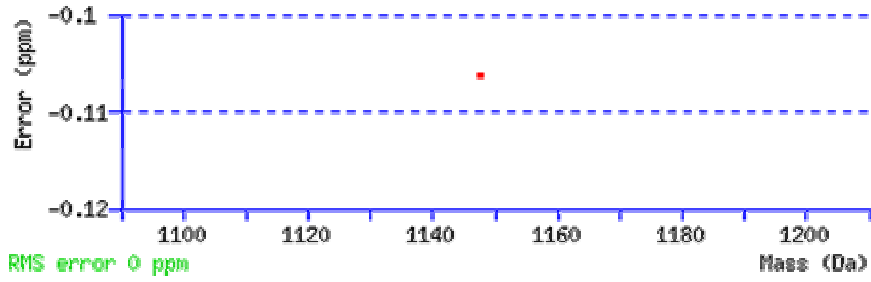
1	MGKNKLLHPS	LVLLLLLVLLP	TDASVSGKPQ	YMVLVPSLLH	TETTEKGCVL
51	LSYLNETHVT	SASLESVRGN	RSLFTDLEAE	NDVLHCVAFA	VPKSSSNEEV
101	MFLTVQVKGP	TQEFKKRTTV	MVKNEDSLVF	VQTDKSIYKP	GQTVKFRVVS
151	MDENFHPLNE	LIPLVYIQDP	KGNRIAQWQS	FQLEGGLKQF	SFPLSSEPFQ
201	GSYKVVVQKK	SGGRTEHPFT	VEEFVLPKFE	VQVTVPKIIT	ILEEEMNVSV
251	CGLYTYGKPV	PGHVTVSICR	KYSDASDCHG	EDSQAFCEKF	SGQLNSHGCF
301	YQQVKTKVFQ	LKRKEYEMKL	HTEAQIQEEG	TVVELTGRQS	SEITRTITKL
351	SFVKVDSHFR	QGIPFFGQVR	LVDGKGVPIP	NKVIFIRGNE	ANYYSNATTD
401	EHGLVQFSIN	TTNVMGTSLT	VRVNYKDRSP	CYGYQWVSEE	HEEAHHTAYL
451	VFSPSKSFVH	LEPMSHELPC	GHTQTVQAHY	ILNGGTLGL	KKLSFYYLIM
501	AKGGIVRTGT	HGLLVKQEDM	KGHFSISIPV	KSDIAPVARL	LIYAVLPTGD
551	VIGDSAKYDV	ENCLANKVDL	SFSPSQSLPA	SHAHLRVTA	PQSVCALRAV
601	DQSVLLMKPD	AELSASSVYN	LLPEKDLTGF	PGPLNDQDNE	DCINRHNVI
651	NGITYTPVSS	TNEKDMYSFL	EDMGLKAFTN	SKIRKPKMCP	QLQQYEMHGP
701	EGLRVGFYES	DVMGRGHARL	VHVEEPTTET	VRKYFPETWI	WDLVVVNSAG
751	VAEVGVTPD	TITWVKAGAF	CLSEDAGLGI	SSTASLRAFQ	PPFVELTMPY
801	SVIRGEAFTL	KATVLNLYPK	CIRVSVQLEA	SPAFLAVPVE	KEQAPHCICA
851	NGRQTVSWAV	TPKSLGNVNF	TVSAEALSEQ	ELCGTEVPSV	PEHGRKDTVI
901	KPLLVEPEGL	EKETTFNSLL	CPSGGEVSEE	LSLKLPPNVV	EESARASVSV
951	LGDILGSAMQ	NTQNLLQMPY	GCGEQNMVLF	APNIYVLDYL	NETQQLTPEI
1001	KSKAIGYLN	GYQRQLNYKH	YDGSYSTFGE	RYGRNQNTW	LTAFVLKTFA
1051	QARAYIFIDE	AHITQALIWL	SQRQKDNCGF	RSSGSLNNA	IKGGVEDEV
1101	LSAYITIAL	EIPLTVTHPV	VRNALFCLES	AWKTAQEGDH	GSHVYTKALL
1151	AYAFALAGNQ	DKRKEVLKSL	NEEAVKKNNS	VHWERPQKPK	APVGHFYEPQ
1201	APSAEVEVMS	YVLLAYLTAQ	PAPTSDELTS	ATNIVKWITK	QQNAQGGFSS
1251	TQDTVVALHA	LSKYGAATFT	RTGKAAQVTI	QSSGTFSSKF	QVDNNNRLLL

1301 QQVSLPELPG EYSMKVTGEG CVYLQTSISKY NILPEKEEFP FALGVQTLPQ
1351 TCDEPKAHTS FQISLSVSYT GSRASANMAI VDVKMVSGFI PLKPTVKMLE
1401 RSNHVSRTVEV SSNHVLIYLD KVSNQTLSEF FTVLQDVPVR DLKPAIVKVY
1451 DYYETDEFAI AEYNAPCSKD LGNA

Unformatted sequence string: **1474 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Query	Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Score	Expe
1	361 - 370	1148.6210	1147.6137	1147.6138	-0.11	0	63	1.8e-



Protein View: CERU_HUMAN

Ceruloplasmin OS=Homo sapiens OX=9606 GN=CP PE=1 SV=1

Database: SwissProt
Score: 80
Monoisotopic mass (M_r): 122983
Calculated pI: 5.44
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of CERU_HUMAN against nr.](#)

Search parameters

MS data file: DATA.TXT
Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)

Protein sequence coverage: 1%

Matched peptides shown in **bold red**.

```

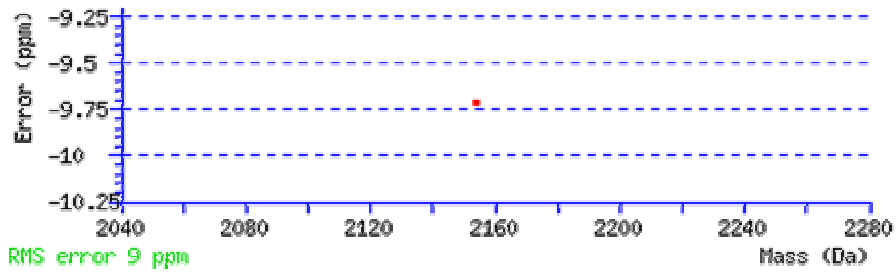
1  MKILILGIFL  FLCSTPAWAK  EKHYIIGIIE  TTWDYASDHG  EKKLISVDTE
51  HSNIYLQNGP  DRIGRLYKKA  LYLQYTDETF  RTTIEKPVWL  GFLGPIIKAE
101  TGDKVYVHLK  NLASRPYTFH SHGITYYKEH  EGAIYPDNTT  DFQRADDKVY
151  PGEQYTYMLL  ATEEQSPGEG  DGNVCVTRIH  SHIDAPKDIA  SGLIGPLIIC
201  KKDSLDKKEK  KHIDREFVVM  FSVVDENFSW  YLEDNIKTYC  SEPEKVDKDN
251  EDFQESNRMY  SVNGYTFGSL  PGLSMCAEDR  VKWYLFMGMN  EVDVHAAFFH
301  GQALTNKNYR  IDTINLFPAT  LFDAYMVAQN  PGEWMLSCQN  LNHLKAGLQA
351  FFQVQECNKS  SSKDNIRGKH  VRHYIIAAEE  IIWNYAPSGI  DIFTKENLTA
401  PGSDSAVFFE  QGTRIGGSY  KKLVIYREYTD  ASFTNRKERG  PEEEHLGILG
451  PVIWAEVGDG  IRVTFHNKGA  YPLSIEPIGV  RFNKNNEGTY  YSPNYPQSR
501  SVPPSASHVA  PTETFTYEWY  VPKEVGPTNA  DPVCLAKMY  SAVDPTKDIF
551  TGLIGPMKIC  KKGSLHANGR  QKDVDKEFY  FPTVFDENES  LLEDNIRMF
601  TTAPDQVDKE  DEDFQESNKM  HSMNGFMYGN  QPGLTMCKGD  SVVWYLFSA
651  NEADVHGIYF  SGNTYLWRGE  RRDTANLFPQ  TSLTLHMWPD  TEGTFNVECL
701  TTDHYTGGMK  QKYTVNQCR  QSEDSTFYLG  ERTYYIAAVE  VEWDYSPQRE
751  WEKELHHLQE  QNVSNAFLDK  GEFYIGSKYK  KVVYRQYTD  TFRVPVERKA
801  EEEHLGILGP  QLHADVGDKV  KIIFKNMATR  PYSIHAHVQ  TESSTVPTL
851  PGETLTYVWK  IPERSGAGTE  DSACIPWAY  STVDQVKDLY  SGLIGPLIVC
901  RRPYLKVFNP  RRRKLEFALL  LVFDENESWY  LDDNIKTYSD  HPEKVNKDDE
951  EFIESNKMHA  INGRMFGNLQ  GLTMHVGDEV  NWYLMGMGNE  IDLHTVHFHG
1001  HSFQYKHRGV  YSSDVFDIFP  GTYQTLEMFP  RTPGIWLLHC  HVTDHIHAGM
1051  ETTYTVLQNE  DTKSG

```

Unformatted sequence string: **1065 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Query	Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Score	Expe
1	111 - 128	2155.0510	2154.0437	2154.0647	-9.72	0	80	2.3e-



Protein View: ITIH4_HUMAN

Inter-alpha-trypsin inhibitor heavy chain H4 OS=Homo sapiens OX=9606 GN=ITIH4 PE=1 SV=4

Database: SwissProt
Score: 50
Monoisotopic mass (M_r): 103521
Calculated pI: 6.51
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of ITIH4_HUMAN against nr.](#)

Search parameters

MS data file: DATA.TXT
Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)

Protein sequence coverage: 2%

Matched peptides shown in **bold red**.

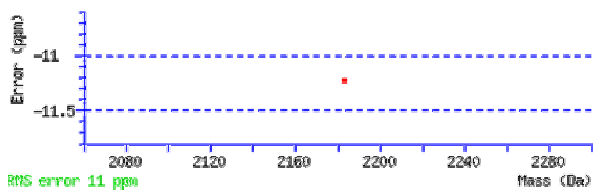
```

1 MKPPRPVRTC SKVLVLLSLL AIHQTTTAEK NGIDIYSLTV DSRVSSRFAH
51 TVVTSRVVNR ANTVQEATFQ MELPKKAFIT NFSMIIDGMT YPGIIKEKAE
101 AQAQYSAAVA KGKSAGLVKA TGRNMEQFQV SVSVAPNAKI TFEVLVYEELL
151 KRRLGVYELL LKVRPQQLVK HLQMDIHIFE PQGISFLETE STFMTNQLVD
201 ALTTWQNKTK AHIRFKPTLS QQQKSPEQQE TVLDGNLIIR YVDRAISGG
251 SIQIENGYFV HYFAPEGLTT MPKNVVFVID KSGSMSGRKI QQTREALIKI
301 LDDLSPRDQF NLIVFSTEAT QWRPSLVPAS AENVNKARSF AAGIQALGGT
351 NINDAMLMAV QLLDSSNQEE RLPESVSLI ILLTDGDPTV GETNPRSIQN
401 NVREAVSGRY SLFCLGFGFD VSYAFLEKLA LDNGGLARRI HEDSDSALQL
451 QDFYQEVANP LLTAVTFEYP SNAVEEVTQN NFRLLFKGSE MIVAGKLQDR
501 GPDVLTATVS GKLPTQNITF QTESSVAEQE AEFQSPKYIF HNFMERLWAY
551 LTIQQLEQT VSASDADQQA LRNQALNLSL AYSFVTPLTS MIVTKPDDQE
601 QSQVAEKPME GESRNRNVHS GSTFFKYLLQ GAKIPKPEAS FSPRRGWNRO
651 AGAAGSRMNF RPGVLSRQL GLPGPPDVPD HAAYHFFRRL AILPASAPPA
701 TSNPDPAVSR VMNMKIEETT MTTQTPAPIQ APSAILPLPG QSVRLCVDP
751 RHRQGPVNL SDPEQGVVET GQYEREKAGF SWIEVTFKNP LVVWHASPEH
801 VVTRNRRSS AYKWKETLFS VMPGLKMTMD KTGLLLSDP DKVTIGLLFW
851 DGRGGLRLL LRDTDRFSSH VGGTLGQFYQ EVLWGSPPAS DDGRRTLVRQ
901 GNDHSATRER RLDYQEGPPG VEISCWSVEL
    
```

Unformatted sequence string: **930 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Query	Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Score	Expect	Rank	U	Peptide
1	669 - 688	2184.0740	2183.0667	2183.0912	-11.2	0	50	0.00025	1	U	R.QLGLPGPPDVPDHAAYHFFR.R



Protein View: TRFE_HUMAN

Serotransferrin OS=Homo sapiens OX=9606 GN=TF PE=1 SV=3

Database: SwissProt
 Score: 91
 Monoisotopic mass (M_r): 79294
 Calculated pI: 6.81
 Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of TRFE_HUMAN against nr.](#)

Search parameters

MS data file: DATA.TXT
 Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
 Fixed modifications: Carbamidomethyl (C)
 Variable modifications: Oxidation (M)

Protein sequence coverage: 2%

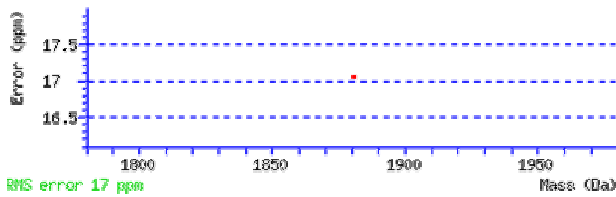
Matched peptides shown in **bold red**.

1 MRLAVGALLV CAVLGLCLAV PDKTVRWCAV SEHEATKCQS FRDHMKSVIP
 51 SDGPSVACVK KASYLDCIRA IAANEADAVT LDAGLVYDAY LAPNNLKPVV
 101 AEFYGSKEDP QTFYYAVAVV KKDSGFQMNQ LRGKKSCHTG LGRSAGWNIP
 151 IGLLYCDLPE PRKPLEKAVA NFFSGSCAPC ADGTDFFQLC QLCPCGCGST
 201 LNQYFGYSGA FKCLKDGAGD VAFVKHSTIF ENLANK**ADRD QYELLCLDNT**
 251 **R**KFPVDEYKDC HLAQVPSHTV VARSMGGKED LIWELLNQAQ EHFQKDKSKE
 301 FQLFSSPHGK DLLFKDSAAG FLKVPPRMDA KMYLGYEYVT AIRNLREGTC
 351 PEAPTDECKP VKWCALSHHE RLKCEWSVN SVGKIECVSA ETTEDCIAKI
 401 MNGEADAMSL DGGFVYIAGK CGLVPVLAEN YNKSDNCEDT PEAGYFAIAV
 451 VKKSASDLTW DNLKGGKKSCH TAVGRTAGWN IPMGLLYNKI NHCRFDEFFS
 501 EGCAPGSKKD SSLCKLCMGS GLNLCEPNNK EGYGYTGAF RCLVEKGDVA
 551 FVKHQVTPQN TGGKNPDPWA KNLNEKDYEL LCLDGRKPV EYANCHLAR
 601 APNHAVVTRK DKEACVHKIL RQQQHLFGSN VTDCSGNFCL FRSETKDLLE
 651 RDDTVCLAKL HDRNTYEKYL GEEYKAVGN LRKSTSSLL EACTFRFP

Unformatted sequence string: 698 residues (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Query	Start - End	Observed	Mr(expt)	Mr(calc)	ppm	M	Score	Expect	Rank	U	Peptide
1	237 - 251	1881.9080	1880.9007	1880.8687	17.1	1	91	2e-08	1	U	K.ADRDQYELLCLDNT.R.K



Protein View: A1AT_HUMAN

Alpha-1-antitrypsin OS=Homo sapiens OX=9606 GN=SERPINA1 PE=1 SV=3

Database: SwissProt
Score: 99
Monoisotopic mass (M_r): 46878
Calculated pI: 5.37
Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of A1AT_HUMAN against nr.](#)

Search parameters

MS data file: DATA.TXT
Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
Fixed modifications: Carbamidomethyl (C)
Variable modifications: Oxidation (M)

Protein sequence coverage: 3%

Matched peptides shown in **bold red**.

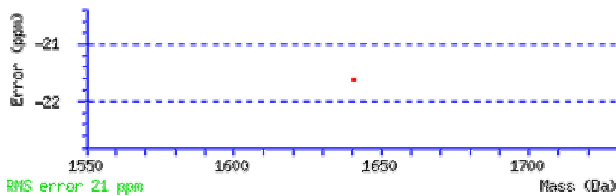
```

1  MPSSVSWGIL  LLAGLCCLVP  VSLAEDPQGD  AAQKTDTSHH  DQDHPTFNKI
51  TPNLAEFAPS  LYRQLAHQSN  STNIFFSPVS  IATAFAMLSL  GTKADTHDEI
101 LEGLNFNLTE  IPEAQIHEGF  QELLRTLNPQ  DSQLQLTTGN  GLFLSEGLKL
151 VDKFLEDVKK  LYHSEAFQVN  FGDTEEAKKQ  INDYVEKGTQ  GKIVDLVKEL
201 DRDTVFALVN  YIFFKQKWER  PFEVKDTEEE  DFHVDQVTTV  KVPMMKRLGM
251 FNIQHCKKLS  SWVLLMKYLG  NATAIFFLPD  EGKLQHLENE  LTHDIITKFL
301 ENEDRRSASL  HLPKLSITGT  YDLKSVLGQL  GITKVFSSNGA  DLSGVTEEAP
351 LKLSKAVHKA  VLTIDEKGTE  AAGAMFLEAI  PMSIPPEVKF  NKPFVFLMIE
401 QNTKSPFLMG  KVVNPTQK
    
```

Unformatted sequence string: **418 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Query	Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Score	Expect	Rank	U	Peptide
1	50 - 63	1641.8280	1640.8207	1640.8562	-21.6	0	99	3.8e-09	1	U	K.ITPNLAEFAPSLYR.Q



Protein View: APOA4_HUMAN

Apolipoprotein A-IV OS=Homo sapiens OX=9606 GN=APOA4 PE=1 SV=3

Database: SwissProt
 Score: 76
 Monoisotopic mass (M_r): 45371
 Calculated pI: 5.28
 Taxonomy: [Homo sapiens](#)

Sequence similarity is available as [an NCBI BLAST search of APOA4_HUMAN against nr.](#)

Search parameters

MS data file: DATA.TXT
 Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
 Fixed modifications: [Carbamidomethyl \(C\)](#)
 Variable modifications: [Oxidation \(M\)](#)

Protein sequence coverage: 2%

Matched peptides shown in **bold red**.

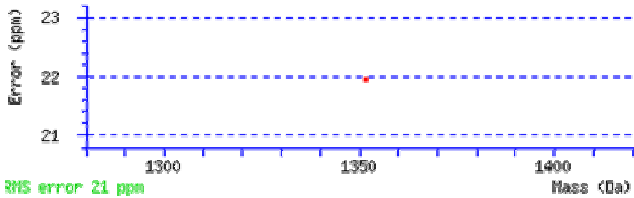
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1 MFLKAVVLTLL ALVAVAGARA EVSADQVATV MWDYFSQLSN NAKEAVEHLQ
51 KSELTQQLNA LFQDKLGEVN TYAGDLQKKL VPFATELHER LAKDSEKLKE
101 EIGKELEELR ARLLPHANEV SQKIGDNLRE LQQRLEPYAD QLRTQVNTQA
151 EQLRRQLTPY AQRMERVLRE NADSLQASLR PHADELKAKI DQNVVEELKGR
201 LTPYADEFKV KIDQTVEELR RSLAPYAQDT QEKLNHQLEG LTFQMKNNAE
251 ELKARISASA EELRQLAPL AEDVRGNLRG NTEGLQKSLA ELGGHLDQQV
301 EEFRRRVEPEY GENFNKALVQ QMEQLRQKLG PHAGDVEGHL SFLEKDLRDK
351 VNSFFSTFKE KESQDKTSL PELEQQQEQQ QEQQQEQQVM LAPLES
    
```

Unformatted sequence string: [396 residues](#) (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Query	Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Score	Expect	Rank	U	Peptide
1	306 - 316	1352.6890	1351.6817	1351.6520	22.0	1	76	7.9e-07	1	U	R.RVEPEYGENFNK.A



Protein View: HPT_HUMAN

Haptoglobin OS=Homo sapiens OX=9606 GN=HP PE=1 SV=1

Database: SwissProt
 Score: 75
 Monoisotopic mass (M_r): 45861
 Calculated pI: 6.13
 Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of HPT_HUMAN against nr.](#)

Search parameters

MS data file: DATA.TXT
 Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
 Fixed modifications: Carbamidomethyl (C)
 Variable modifications: Oxidation (M)

Protein sequence coverage: 3%

Matched peptides shown in **bold red**.

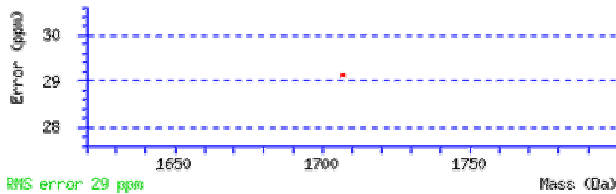
```

1 MSALGAVIAL LLWGQLFAVD SGNDVTDIAD DGCCKPPEIA HGYVEHSVRY
51 QCKNYKLRTE EGDGVYTLND KKQWINKAVG DKLPECEADD GCPKPEIAH
101 GYVEHSVRYQ CKNYYKLRTE GDGVYTLNNE KQWINKAVGD KLPECEAVCG
151 KPNPANPVQ RILGGHLDK GSPFWQAKMV SHHNLTTGAT LINEQWLLTT
201 AKNLFLNHSE NATAKDIAPT LTLYVGKKQL VEIEKVVLP NYSQVDIGLI
251 KLRQKVSUNE RVMPICLPSK DYAEVGRVGY VSGWGRNANF KFTDHLKYVM
301 LPVADQDQCI RHYEGSTVPE KKTTPKSPVGV QPILNEHTFC AGMSKYQEDT
351 CYGDAGSFAFA VHDLEEDTWY ATGILSFDKS CAVAEGVYV KVTSIQDWVQ
401 KTIAEN
    
```

Unformatted sequence string: 406 residues (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Query	Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Score	Expect	Rank	U	Peptide
1	298 - 311	1707.8690	1706.8617	1706.8120	29.1	0	75	9.6e-07	1	U	K.YVMLPVADQDQCIR.H



Protein View: RET4_HUMAN

Retinol-binding protein 4 OS=Homo sapiens OX=9606 GN=RBP4 PE=1 SV=3

Database: SwissProt
 Score: 65
 Monoisotopic mass (M_r): 23337
 Calculated pI: 5.76
 Taxonomy: [Homo sapiens](#)

Sequence similarity is available as [an NCBI BLAST search of RET4_HUMAN against nr.](#)

Search parameters

MS data file: DATA.TXT
 Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
 Fixed modifications: [Carbamidomethyl \(C\)](#)
 Variable modifications: [Oxidation \(M\)](#)

Protein sequence coverage: 10%

Matched peptides shown in **bold red**.

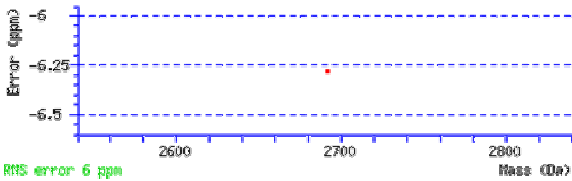
```

1 MKWVWALLLL AALGSGRAER DCRVSSFRVK ENFDKARFSG TWYAMAKKDP
51 EGLFLQDNIV AEFVDEETGQ MSATAKGRVR LLNNDVDCAD MVGTFTDTED
101 PAKFKMKYWG VASFLQKND DHWIVDTDYD TYAVQYSCRL LNLDTGCADS
151 YSFVFSRDPN GLPPEAQKIV RQRQEELCLA RQYRLIVHNG YCDGRSERNL
201 L
    
```

Unformatted sequence string: [201 residues](#) (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Query	Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Score	Expect	Rank	U	Peptide
1	118 - 139	2693.1040	2692.0967	2692.1136	-6.28	0	65	1.3e-06	1	U	K.GNDDHWIVDTDYDTYAVQYSCRL.L



Protein View: HPT_HUMAN

Haptoglobin OS=Homo sapiens OX=9606 GN=HP PE=1 SV=1

Database: SwissProt
 Score: 70
 Monoisotopic mass (M_r): 45861
 Calculated pI: 6.13
 Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of HPT_HUMAN against nr.](#)

Search parameters

MS data file: DATA.TXT
 Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
 Fixed modifications: Carbamidomethyl (C)
 Variable modifications: Oxidation (M)

Protein sequence coverage: 7%

Matched peptides shown in **bold red**.

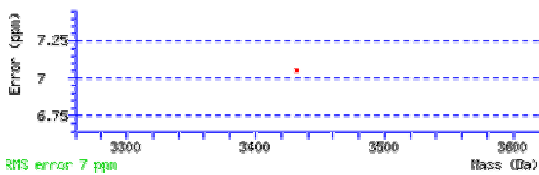
```

1 MSALGAVIAL LLWQLFAVD SGNDVTDIAD DGC PKPPEIA HGYVEHSVRY
51 QCKNYKLRTE EGDGVYTLND KKQWINKAVG DKLPECEADD GCPKPEIAH
101 GYVEHSVRYQ CKNYKLRTE GDGVYTLNNE KQWINKAVGD KLPECEAVCG
151 KPKNPANPVQ RILGGHLDK GSPFWQAKMV SHHNLTTGAT LINEQWLLTT
201 AKNLFNLHSE NATAKDIAPT LTLVVGKKQL VEIEKVVLP NYSQVDIGLI
251 KLRQKQSVNE RVMPICLPSK DYAEVGRVGY VSGWGRNANF KFTDHLKYVM
301 LPVADQDQCI RHYEGSTVPE KKTTPKSPVGV QPILNEHTFC AGMSKYQEDT
351 CYGDAGSAPA VHDLEEDTWY ATGILSFDKS CAVA EYGVVY KVTSIQDWVQ
401 KTIAEN
    
```

Unformatted sequence string: **406 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Query	Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Score	Expect	Rank	U	Peptide
1	78 - 108	3432.6189	3431.6116	3431.5874	7.05	1	70	1.2e-06	1	U	K.AVGDKLPECEADDGCPKPEIAHGYVEHSVRY



Protein View: TTHY_HUMAN

Transthyretin OS=Homo sapiens OX=9606 GN=TTR PE=1 SV=1

Database: SwissProt
 Score: 189
 Monoisotopic mass (M_r): 15991
 Calculated pI: 5.52
 Taxonomy: [Homo sapiens](#)

Sequence similarity is available as [an NCBI BLAST search of TTHY_HUMAN against nr.](#)

Search parameters

MS data file: DATA.TXT
 Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
 Fixed modifications: [Carbamidomethyl \(C\)](#)
 Variable modifications: [Oxidation \(M\)](#)

Protein sequence coverage: 15%

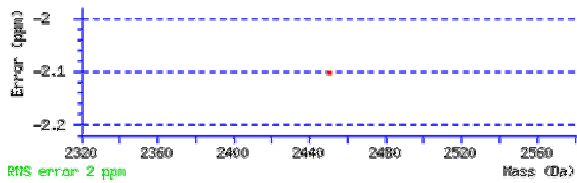
Matched peptides shown in **bold red**.

1 MASHRLLLLC LAGLVFVSEA GPTGTGESKC PLMVKVLDAV RGSPAINVAV
 51 HVFRKAADDT WEPFASGKTS ESGELHGLTT EEEFVEGIYK VEIDTKSYWK
 101 **ALGISPFHEH AEVVFTANDS GPRRYTIAAL** LSPYSYSTTA VVTNPKE

Unformatted sequence string: **147 residues** (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Query	Start - End	Observed	Mr (expt)	Mr (calc)	ppm	M	Score	Expect	Rank	U	Peptide
1	101 - 123	2451.2000	2450.1927	2450.1979	-2.10	0	189	2.5e-18	1	U	K.ALGISPFHEHAEEVVFTANDSGPR.R



Protein View: HPT_HUMAN

Haptoglobin OS=Homo sapiens OX=9606 GN=HP PE=1 SV=1

Database: SwissProt
 Score: 91
 Monoisotopic mass (M_r): 45861
 Calculated pI: 6.13
 Taxonomy: Homo sapiens

Sequence similarity is available as [an NCBI BLAST search of HPT_HUMAN against nr.](#)

Search parameters

MS data file: DATA.TXT
 Enzyme: Trypsin: cuts C-term side of KR unless next residue is P.
 Fixed modifications: Carbamidomethyl (C)
 Variable modifications: Oxidation (M)

Protein sequence coverage: 3%

Matched peptides shown in **bold red**.

```

1 MSALGAVIAL LLWQQLFAVD SGNDVTDIAD DGCCKPPEIA HGYVEHSVRY
51 QCKNYKLRTE EGDGVYTLND KKQWINKAVG DKLPECEADD GCPKPEIAH
101 GYVEHSVRYQ CKNYYKLRTE GDGVYTLNNE KQWINKAVGD KLPECEAVCG
151 KPNPANPVQ RILGGHLDK GSPFWQAKMV SHHNLTTGAT LINEQWLLTT
201 AKNLFNLHSE NATAKDIAPT LTLYVGKKQL VEIEKVVLP NYSQVDIGLI
251 KLKQKVSUNE RVMPICLPSK DYAEVGRVGY VSGWGRNANF KFTDHLKYVM
301 LPVADQDQCI RHYEGSTVPE KKTPKSPVGV QPILNEHTFC AGMSKYQEDT
351 CYGDAGSFAFA VHDLEEDTWY ATGILSFDKS CAVAEGVYV KVTSIQDWVQ
401 KTIAEN
    
```

Unformatted sequence string: 406 residues (for pasting into other applications).

Sort by residue number increasing mass decreasing mass
 Show matched peptides only predicted peptides also

Query	Start - End	Observed	Mr(expt)	Mr(calc)	ppm	M	Score	Expect	Rank	U	Peptide
1	117 - 131	1708.9120	1707.9047	1707.8428	36.3	1	91	3.1e-08	1	U	K.LRTEGDGVYTLNNEK.Q

