

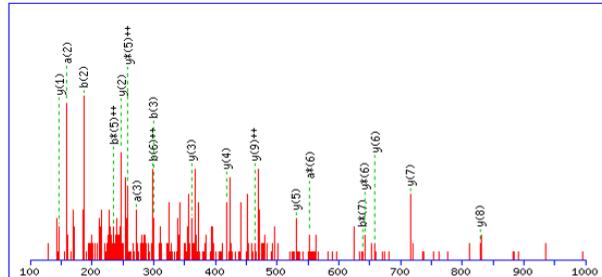
Supplemental Figure S1A

Peak ID: 1740

MS/MS Fragmentation of **SVLGQLGITK**

Found in **gi|24438**, alpha-1-antitrypsin (aa 268-394) [Homo sapiens]

Match to Query 6: 1014.645800 from(508.330176,2+)



Monoisotopic mass of neutral peptide Mr(calc): 1014.6073

Ions Score: 15 Expect: 47

Matches (**Bold**): 19/92 fragment ions using 81 most intense peaks

#	a	a ⁺⁺	a*	a* ⁺⁺	b	b ⁺⁺	b*	b* ⁺⁺	Seq.	y	y ⁺⁺	y*	y* ⁺⁺	#
1	60.0444	30.5258			88.0393	44.5233			S					10
2	159.1128	80.0600			187.1077	94.0575			V	928.5826	464.7949	911.5560	456.2817	9
3	272.1969	136.6021			300.1918	150.5995			L	829.5142	415.2607	812.4876	406.7475	8
4	329.2183	165.1128			357.2132	179.1103			G	716.4301	358.7187	699.4036	350.2054	7
5	457.2769	229.1421	440.2504	220.6288	485.2718	243.1395	468.2453	234.6263	Q	659.4087	330.2080	642.3821	321.6947	6
6	570.3610	285.6841	553.3344	277.1709	598.3559	299.6816	581.3293	291.1683	L	531.3501	266.1787	514.3235	257.6654	5
7	627.3824	314.1949	610.3559	305.6816	655.3774	328.1923	638.3508	319.6790	G	418.2660	209.6366	401.2395	201.1234	4
8	740.4665	370.7369	723.4400	362.2236	768.4614	384.7343	751.4349	376.2211	I	361.2445	181.1259	344.2180	172.6126	3
9	841.5142	421.2607	824.4876	412.7475	869.5091	435.2582	852.4825	426.7449	T	248.1605	124.5839	231.1339	116.0706	2
10									K	147.1128	74.0600	130.0863	65.5468	1

Matched peptides shown in **Bold Red**

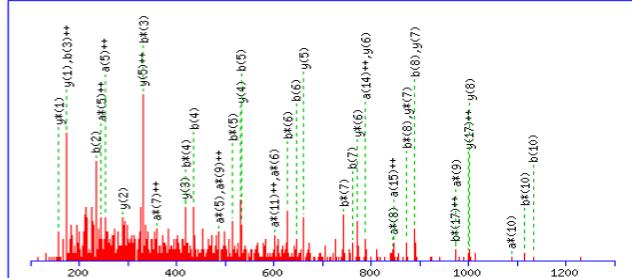
1 THDIITKFLE NEDRRSASLH LPKLSITGTY DLK**SVLGQLG ITKVFSNGAD**
 51 LSGVTEEAPL KLSKAVHKAV LTIDEKGTEA AGAMFLEAIP MSIPPEVKFN
 101 KPFVFLMIEQ NTKSPLFMGK VVNPTQK

Supplemental Figure S1B

Peak ID: 11165

MS/MS Fragmentation of **HPNSPLDEENLTQENQDR**

Found in **gi|443345**, Chain A, Alpha1 Antichymotrypsin
Match to Query 2: 2134.980774 from(712.667534,3+)



Monoisotopic mass of neutral peptide Mr(calc): 2134.9515

Ions Score: 28 Expect: 0.93

Matches (**Bold**): 41/196 fragment ions using 64 most intense peaks

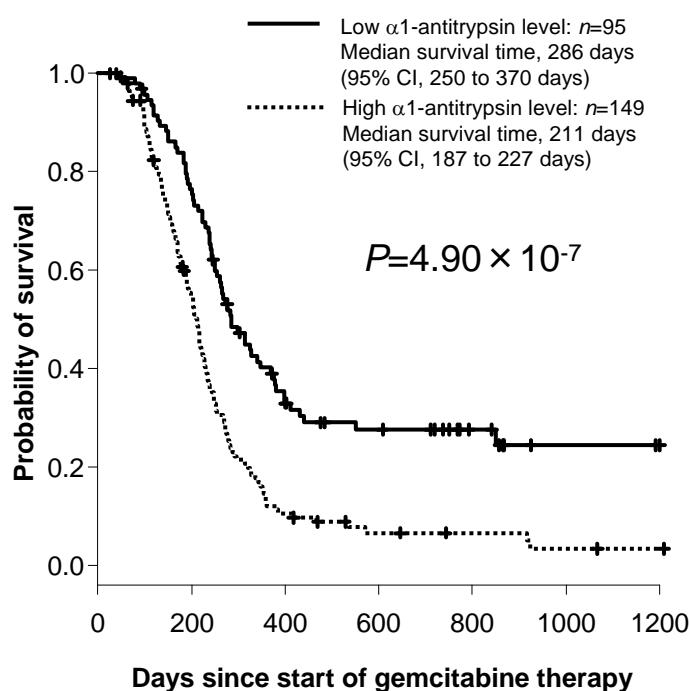
#	a	a ⁺⁺	a*	a ⁺⁺⁺	b	b ⁺⁺	b*	b ⁺⁺⁺	Seq.	y	y ⁺⁺	y*	y ⁺⁺⁺	#
1	110.0713	55.5393			138.0662	69.5367			H					18
2	207.1240	104.0657			235.1190	118.0631			P	1998.8999	999.9536	1981.8734	991.4403	17
3	321.1670	161.0871	304.1404	152.5738	349.1619	175.0846	332.1353	166.5713	N	1901.8472	951.4272	1884.8206	942.9139	16
4	408.1990	204.6031	391.1724	196.0899	436.1939	218.6006	419.1674	210.0873	S	1787.8042	894.4058	1770.7777	885.8925	15
5	505.2518	253.1295	488.2252	244.6162	533.2467	267.1270	516.2201	258.6137	P	1700.7722	850.8897	1683.7456	842.3765	14
6	618.3358	309.6715	601.3093	301.1583	646.3307	323.6690	629.3042	315.1557	L	1603.7194	802.3634	1586.6929	793.8501	13
7	733.3628	367.1850	716.3362	358.6717	761.3577	381.1825	744.3311	372.6692	D	1490.6354	745.8213	1473.6088	737.3080	12
8	862.4054	431.7063	845.3788	423.1930	890.4003	445.7038	873.3737	437.1905	E	1375.6084	688.3079	1358.5819	679.7946	11
9	991.4479	496.2276	974.4214	487.7143	1019.4429	510.2251	1002.4163	501.7118	E	1246.5658	623.7866	1229.5393	615.2733	10
10	1105.4909	553.2491	1088.4643	544.7358	1133.4858	567.2465	1116.4592	558.7333	N	1117.5232	559.2653	1100.4967	550.7520	9
11	1218.5749	609.7911	1201.5484	601.2778	1246.5699	623.7886	1229.5433	615.2753	L	1003.4803	502.2438	986.4538	493.7305	8
12	1319.6226	660.3149	1302.5961	651.8017	1347.6175	674.3124	1330.5910	665.7991	T	890.3963	445.7018	873.3697	437.1885	7
13	1447.6812	724.3442	1430.6546	715.8310	1475.6761	738.3417	1458.6496	729.8284	Q	789.3486	395.1779	772.3220	386.6646	6
14	1576.7238	788.8655	1559.6972	780.3523	1604.7187	802.8630	1587.6922	794.3497	E	661.2900	331.1486	644.2634	322.6354	5
15	1690.7667	845.8870	1673.7402	837.3737	1718.7616	859.8845	1701.7351	851.3712	N	532.2474	266.6273	515.2209	258.1141	4
16	1818.8253	909.9163	1801.7987	901.4030	1846.8202	923.9137	1829.7937	915.4005	Q	418.2045	209.6059	401.1779	201.0926	3
17	1933.8522	967.4298	1916.8257	958.9165	1961.8472	981.4272	1944.8206	972.9139	D	290.1459	145.5766	273.1193	137.0633	2
18									R	175.1190	88.0631	158.0924	79.5498	1

Matched peptides shown in **Bold Red**

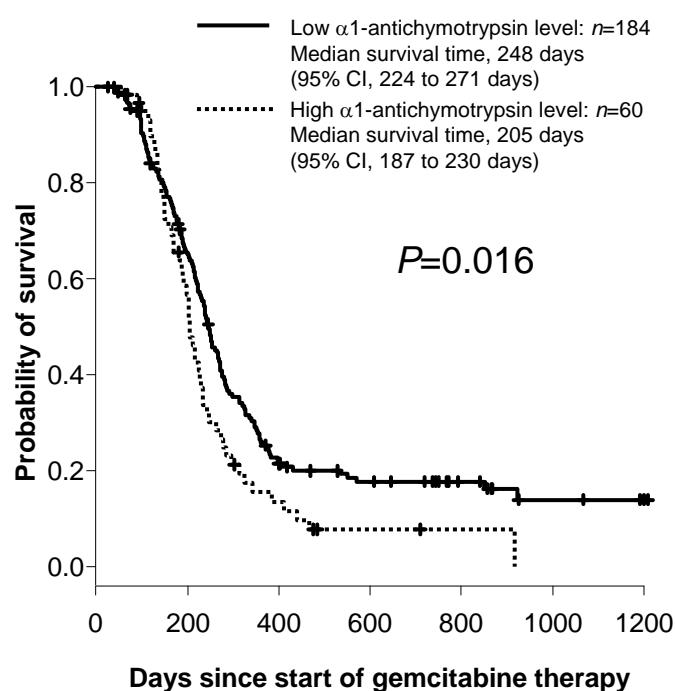
1 **HPNSPLDEEN LTQENQDRGT HVDLGLASAN VDFAFSLYKQ LVLKAPDKNV**
51 **IFSPPLSISTA LAFLSLGAHN TTLTEILKGL KFNLTETSEA EIHQSFQHLL**
101 **RTLNQSSDEL QLSMGNAMFV KEQLSLLDRF TEDAKRLYGS EAFATDFQDS**
151 **AAAKKLINDY VKNGTRGKIT DLIKDLDQSQT MMVLVNYIFF KAKWEMPFDP**
201 **QDTHQSRFYI SKKKWVMVPM MSLHHLTIPY FRDEELSCTV VELKYTGNAS**
251 **ALFILPDQDK MEEVEAMLLP ETLKWRDSL EFREIGELYL PKFSISRDYN**
301 **LNDILLQLGI EEAFTSKADL SGITGARNLA VSQVVHKAVL DVFEEGTEAS**
351 **AATAVKITLL**

Supplemental Figure S2

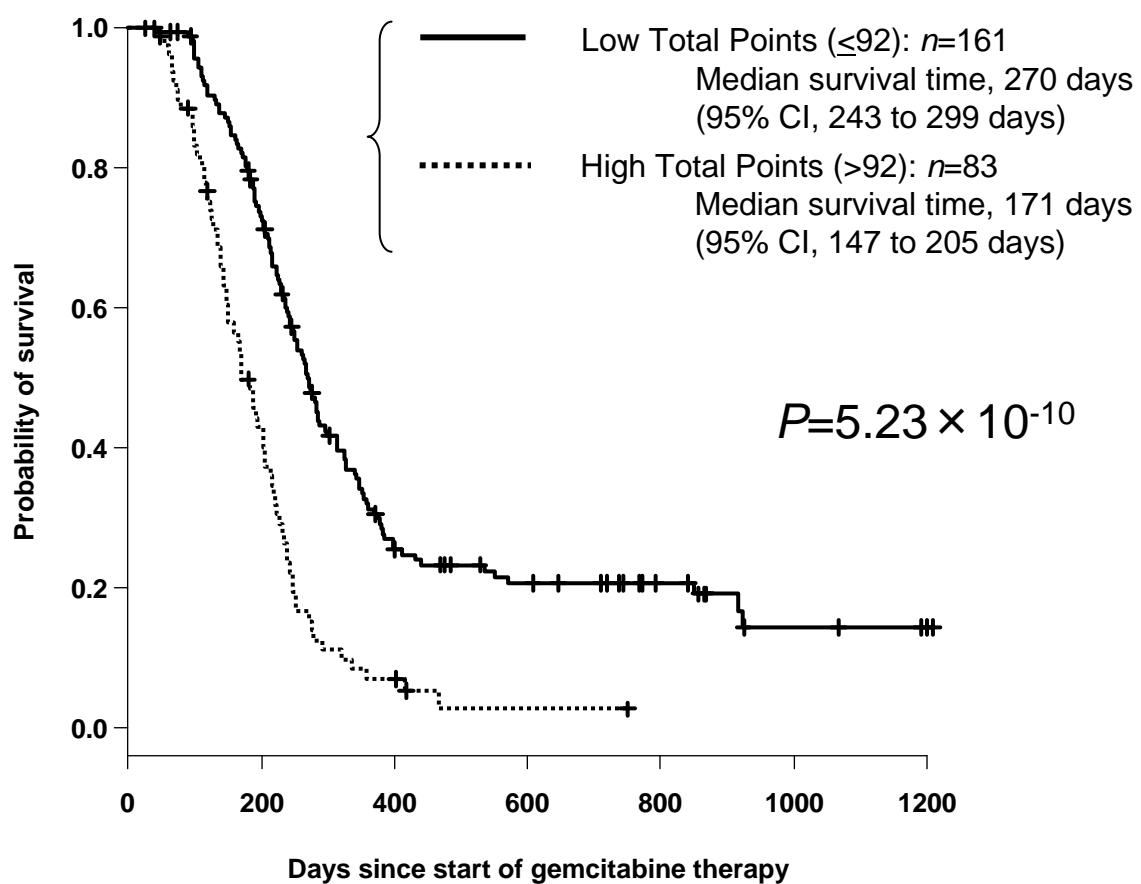
A



B

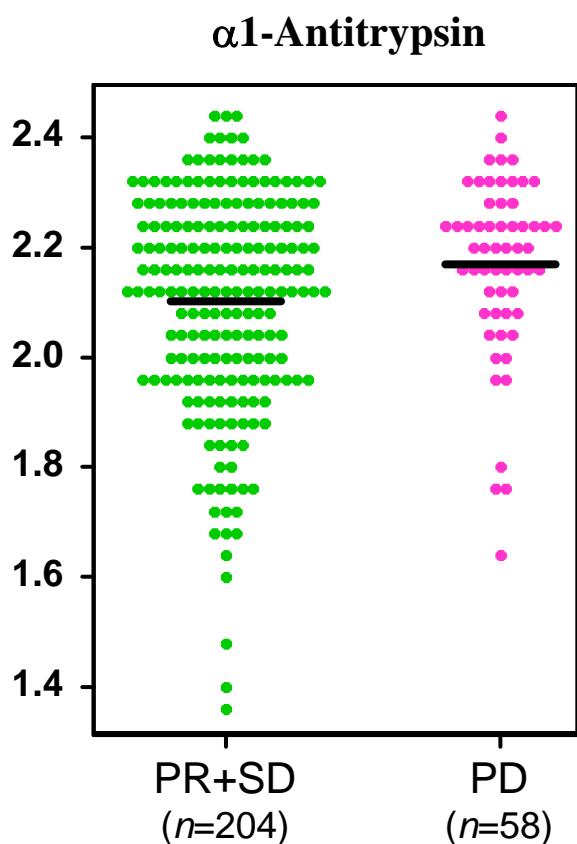


C

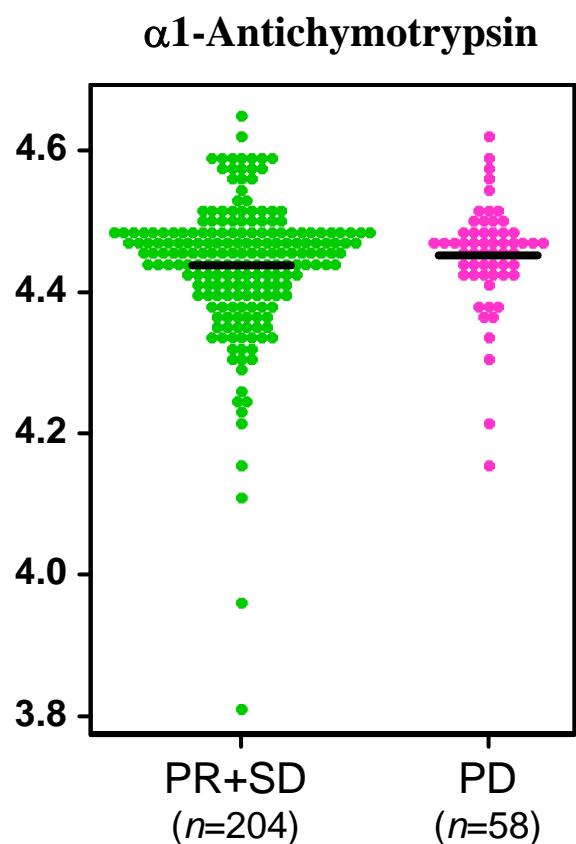


Supplemental Figure S3

A



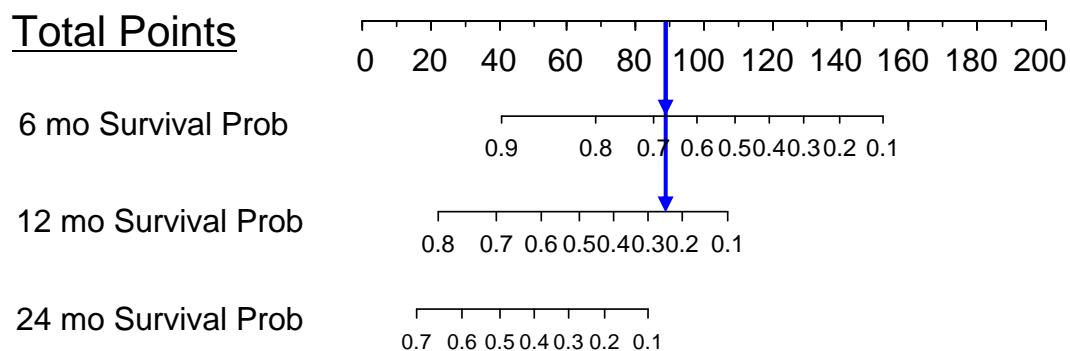
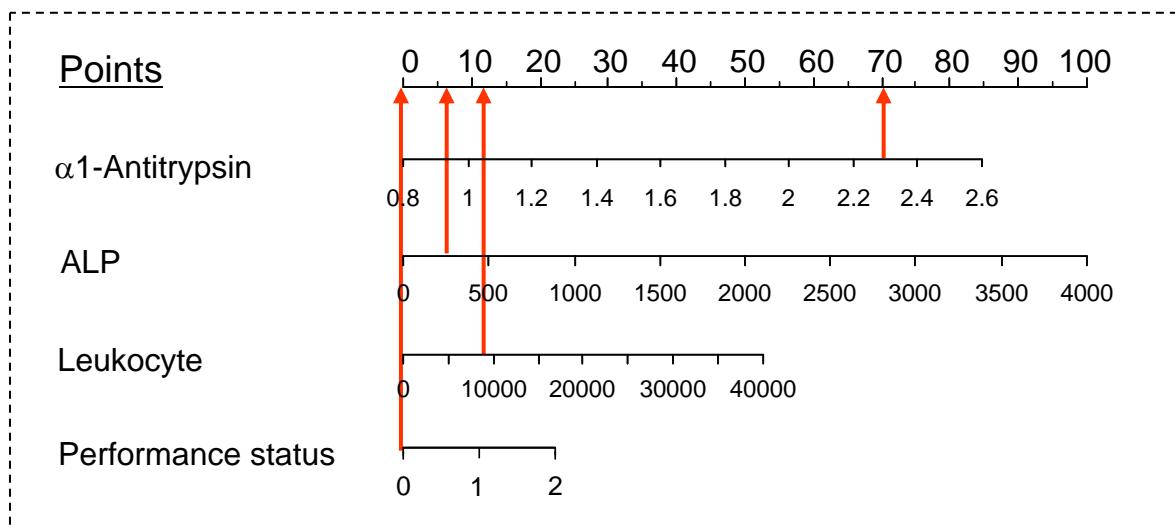
B



$P=0.020$

$P=0.312$

Supplemental Figure S4



LEGENDS FOR SUPPLEMENTAL FIGURES ONLINE

Supplemental Figure S1 online

MS/MS spectra and database search results of two representative MS peaks [ID 1740 (A) and ID 11165 (B)] identified as being derived from α 1-antitrypsin and α 1-antichymotrypsin, respectively. Peptides that matched the amino acid sequences in the database are highlighted in red.

Supplemental Figure S2 online

Kaplan-Meier plots of overall survival in relation to α 1-antitrypsin (A), α 1-antichymotrypsin (B) data, and Total Point score (C) in 244 samples that were not used in the 2DICAL analysis.

Supplemental Figure S3 online

Plasma/serum levels of α 1-antitrypsin and α 1-antichymotrypsin according to tumor response (RECIST guideline). Horizontal lines represent the average levels. Forty-two patients whose tumor response was not evaluable were excluded. PR, partial response; SD, stable disease; PD, progressive disease.

Supplemental Figure S4 online

First, draw a vertical line (*red*) upward to the Points row (*top*) to obtain the number of points for each variable (*top*). Then, total the points and drop a vertical line (*blue*) from the Total Points row to obtain the estimated probability (Prob) of survival 6, 12, and 24 months (mo) after gemcitabine treatment (*bottom*). This particular patient was estimated to have no chance of surviving more than 24 months using the current gemcitabine treatment.