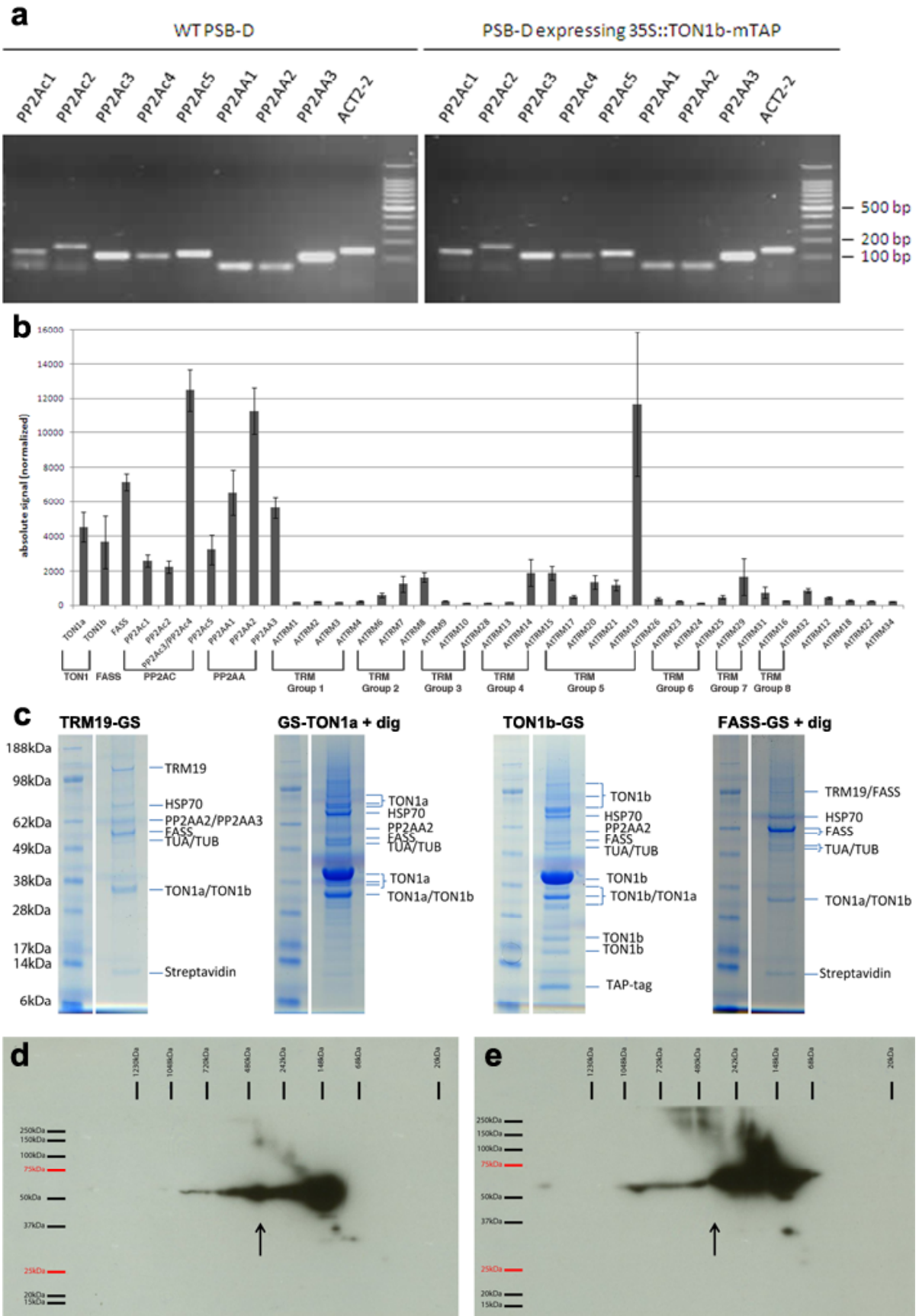


## Supplementary Figure S1



## Supplementary Figure S1: Expression and TAP analysis of TTP members in *Arabidopsis* cell culture

(a) RT-PCR analysis of *Arabidopsis* PP2AA and PP2AC subunits in control cells and in Pro35S:TON1b-mTAP expressing cells.

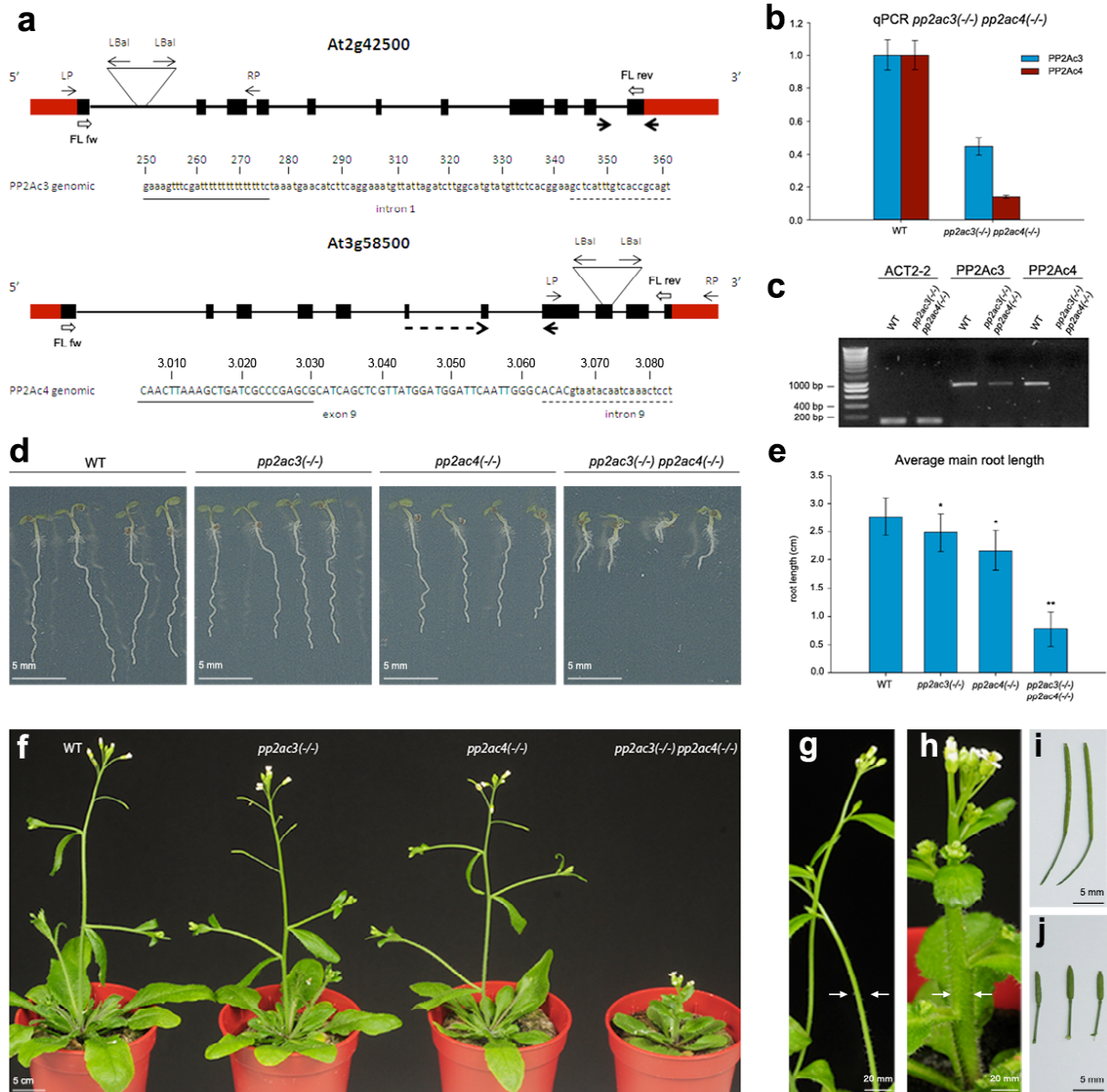
(b) Gene expression analysis of the TTP complex genes in *Arabidopsis* cell culture using the Genevestigator tool<sup>61</sup>. No Affymetrix probes are available for the *TRM5*, *TRM11*, *TRM27*, *TRM30* and *TRM33* genes.

(c) Representative Coomassie stained gels for all bait proteins used in the TAP experiments (TON1a, TON1b, FASS and TRM19) followed by MALDI TOF/TOF analysis. Proteins in the TAP eluates were separated on a 4-12% gradient NuPAGE gel (Invitrogen) and visualized with colloidal Coomassie Brilliant Blue staining. The most abundant proteins detected in each visible band on gel are indicated. In brief, the identified proteins were traced back to their gel slice number via MALDI spot numbers, making it possible to indicate the most abundant identified protein(s) in each visible band on SDS-PAGE. Some of the visible bands on SDS-PAGE correspond to non-specific co-purified proteins like HSP70, TUA/TUB, streptavidin etc... Such background proteins are not presented in the final table of interactors, as known background proteins (obtained from cultures expressing TAP-tagged mock proteins GUS, RFP and GFP) are systematically subtracted from the interactor lists as indicated in the Methods section. Identified proteins which do not give a clear visible band on SDS-PAGE or which are not the most abundant protein in a visible band on the SDS-PAGE gels are not indicated. Left lanes of each panel correspond to molecular weight standards.

(d-e) 2D Blue Native/SDS-PAGE on total protein extracts of PSB-D *Arabidopsis* cell suspension cells expressing Pro35S::TON1b-GS (d) or Pro35S::GS-TON1a (e) followed by Western blot detection using the Peroxidase Anti-Peroxidase (PAP) antibody that binds the GS-tag. The observed smear indicates incorporation of the GS-fused proteins in high molecular weight range complexes with an accumulation at an estimated size of 300-350 kDa (arrows). The strong signal at around 100kDa likely reflects the migration of uncomplexed TON1 proteins as dimers.

<sup>61</sup>Zimmermann, P., Hirsch-Hoffmann, M., Hennig, L., and Gruissem, W. GENEVESTIGATOR. *Arabidopsis* microarray database and analysis toolbox. *Plant Physiol* **136**, 2621-2632 (2004)

## Supplementary Figure S2

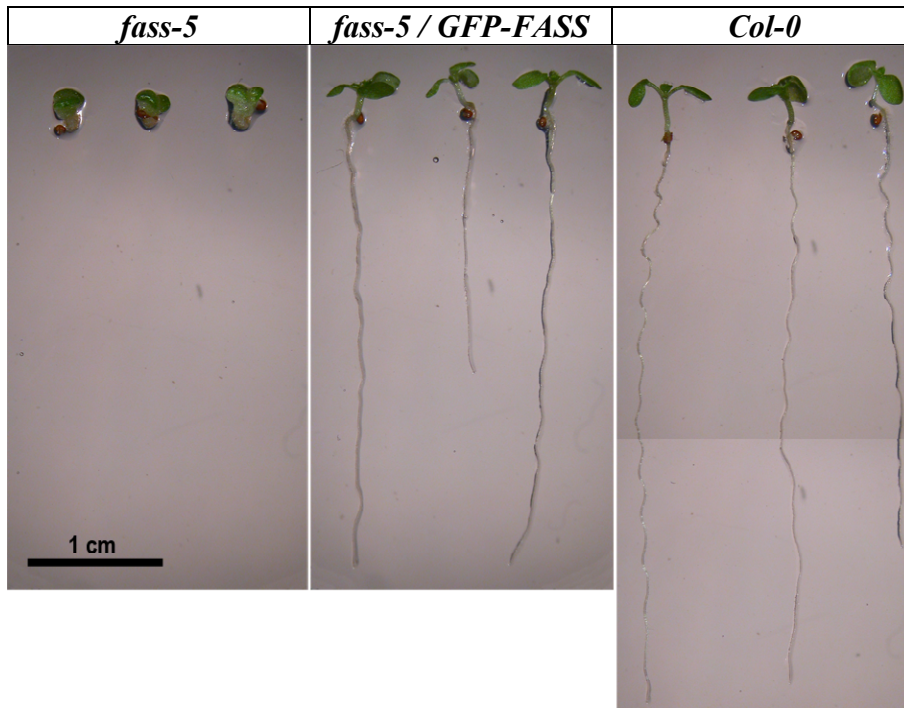


### Supplementary Figure S2 : Analysis of *pp2ac3-c4* double mutant

(a) Gene models of *PP2AC3* and *PP2AC4* with indication of the respective T-DNA insertions and primers used for genotyping (LP, RP, LBAI), qPCR (bold arrows) and RT-PCR (FL fw and FL rev). Sequencing both the LBAI-LP and LBAI-RP PCR products allowed us to determine the exact insertion positions (underlined and dashed sequences) and to confirm the presence of tandem repeat T-DNA insertions in both lines. (b) Quantitative PCR (qPCR) analysis between wild type (Col-0) and the *pp2ac3-c4* double mutant showing significant reduction in *PP2AC3* and *PP2AC4* messenger levels. Expression of *PP2AC3* appears less reduced than *PP2AC4*. Both actin and *eif4A* were used as normalization genes. (c) RT-PCR analysis showing residual expression of full-length *PP2AC3* mRNA, while no full-length *PP2AC4* mRNA could be detected in the *pp2ac3-c4* double mutant. Actin (*Act2-2*) primers were used for normalization. (d) Representative images of wild type (Col-0), *pp2ac3*, *pp2ac4* and *pp2ac3-c4* double mutant seedlings showing severely affected main root growth of the *pp2ac3-c4* double mutant. (e) Average main root length of wild type (n=96), *pp2ac3* (n=38),

*pp2a-c4* (n=35) and the *pp2ac3-ac4* double mutant (n=52). Single *pp2ac* mutants already show a statistical reduction in average main root length compared to wild type (T-test,  $P < 10^{-4}$  and  $P < 10^{-10}$  for *pp2ac3* and *pp2ac4* respectively) and the defects in main root growth are strongly enhanced (T-test,  $P < 10^{-60}$ ) in the *pp2ac3-c4* double mutant. **(f-j)** The *pp2ac3* and *pp2ac4* single mutants develop to wild type looking mature plants while the *pp2ac3-c4* double mutant are dwarfed plant with small, thick and round leaves, thick flowering stems **(g, h)** and shorter siliques **(i, j)**.

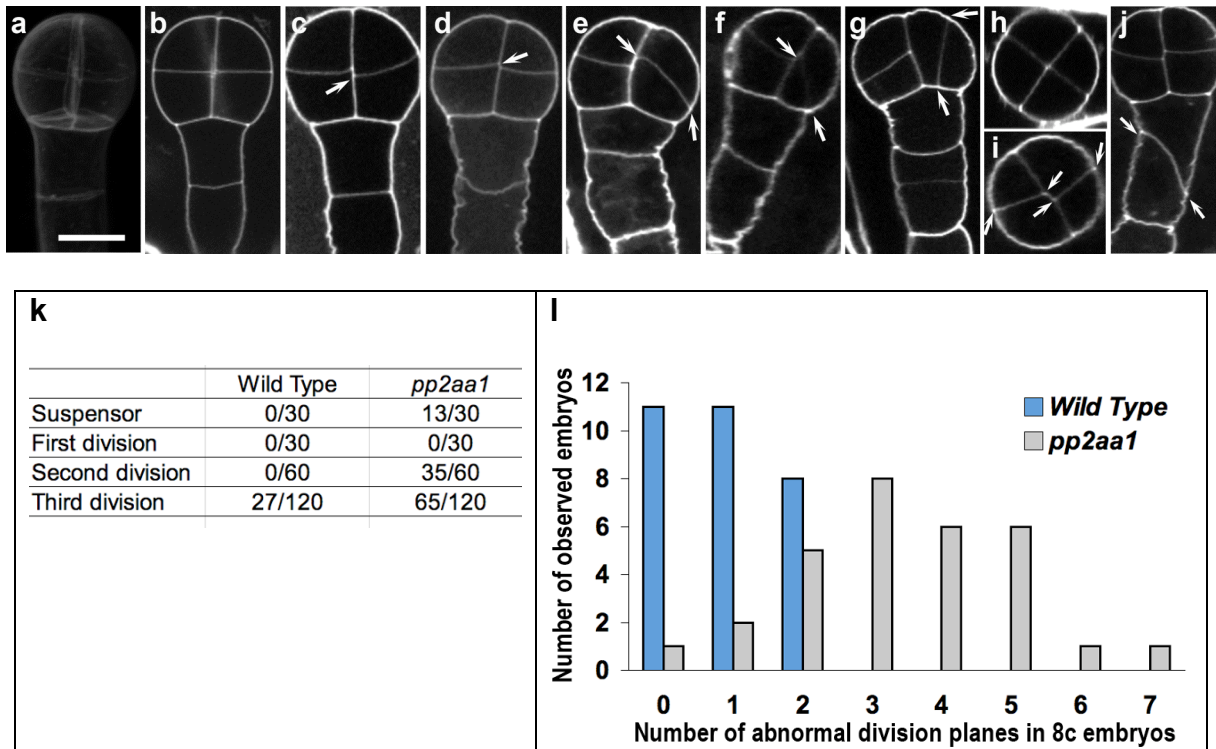
### Supplementary Figure S3



#### Supplementary Figure S3: The GFP-FASS fusion complements the *fass* mutant phenotype

Wild-type *Arabidopsis* and heterozygous *fass-5* plants were infiltrated with *Agrobacterium* carrying the Pro35S:GFP-FASS construct. At least 20 primary transformants were obtained for each genotype. To determine whether the GFP-FASS construct complements the *fass* phenotype, the progenies were analyzed for segregation of the *fass* phenotype and of the hygromycin resistance carried by the T-DNA (GFP-FASS). No plant presenting both hygromycin resistance and the *fass* mutant phenotype was recovered indicating that the GFP-FASS construct complements the *fass* phenotype. Root growth of *fass-5* complemented plants was almost fully restored (compare one-week old seedlings in the middle panel to *fass-5* (left panel) or Col-0 (right panel)).

## Supplementary Figure S4

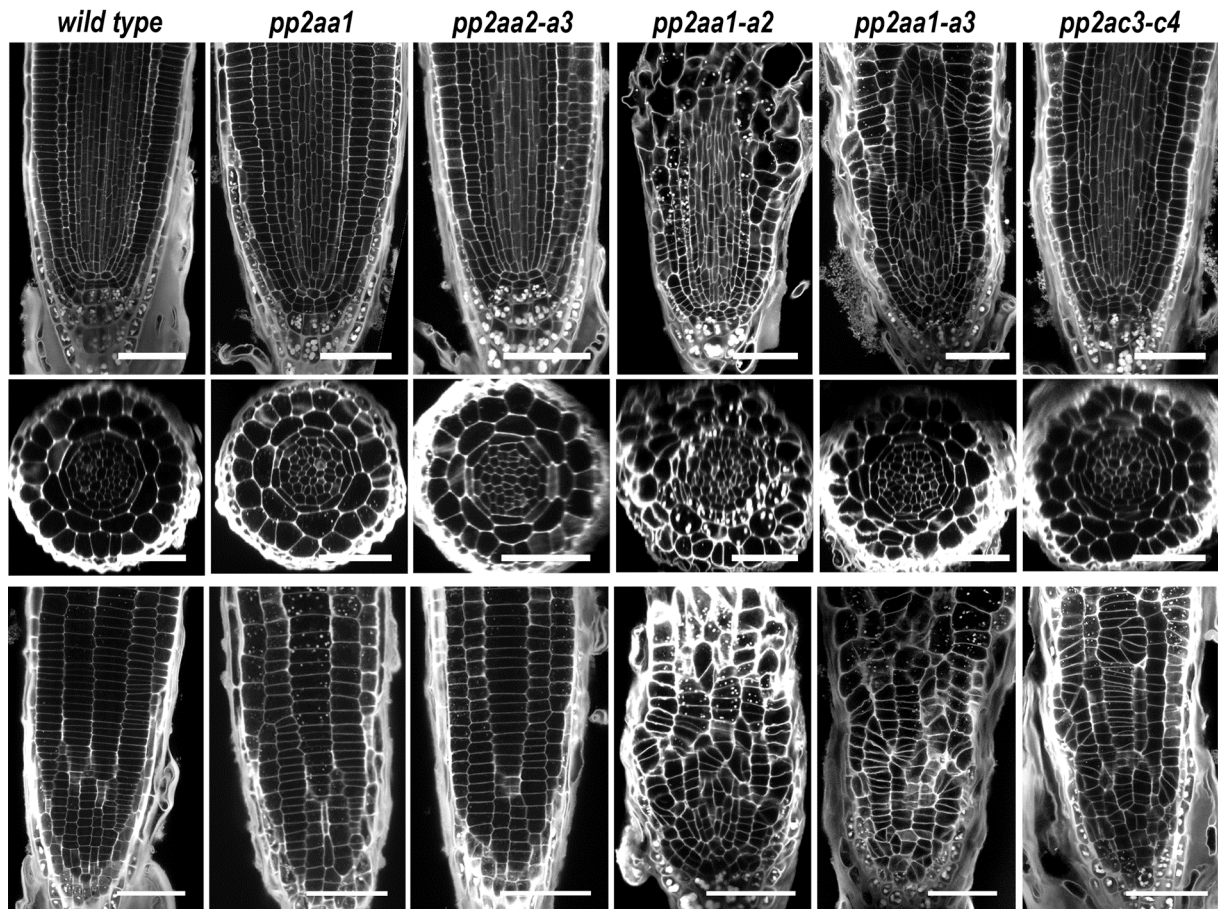


### Supplementary Figure S4: Cell division plane defects in 8-celled *pp2aa1* mutant embryos

(a-j) Propidium Iodide (PI) staining of 8-celled wild type (a-c and h) and *pp2aa1* (d-g, i, j) embryos. In a is shown a 3D reconstruction, in b-g and j, longitudinal views, and in h-i transverse sections of 8-celled embryos. 8-celled embryos result from three rounds of division (seven divisions in total, ie 1+2+4) in the embryo proper. The first division of the apical cell is longitudinal. The two-celled embryo then divides by a second longitudinal division perpendicular to the plane of the previous division. The four resulting cells then divide by transverse divisions, giving rise to the eight-celled embryo. PI imaging and 3D reconstruction of 8-celled embryos allowed to estimate division plane defects during these three rounds of division : Transverse sections enabled to follow the first and second round of division, whereas longitudinal sections revealed abnormal division plane during the third one. The table in k reports cell division plane defects observed in 30 embryos of *pp2aa1* compared to wild type. Abnormal division planes are visible in the second round of divisions in *pp2aa1* embryos where 35 planes were tilted (as shown in i) compared to wild type embryos (h). Defects are even more pronounced during the third round of division. At this stage, slightly tilted planes are present occasionally in wild-type embryos (27/120; as shown in c). Among 65 abnormal division planes observed in *pp2aa1* embryos, 35 appeared slightly tilted as in the wild type (c-d), while, in contrast to wild type, 15 were strongly misplaced as in e and 16, as in f or g. Aberrant division planes are also frequent in the suspensor of *pp2aa1* embryos (in 13/30 embryos as shown in j) but never observed in wild type. (l) Number of embryos displaying 0 to 7 abnormal division planes in wild type compared to *pp2aa1* embryos. Most *pp2aa1* embryos have more than 2 abnormal division planes. Wild type embryos reported here as having abnormal planes only have slightly tilted division planes.

Images in a to j are at the same scale (bar = 10  $\mu$ m in a).

## Supplementary Figure S5

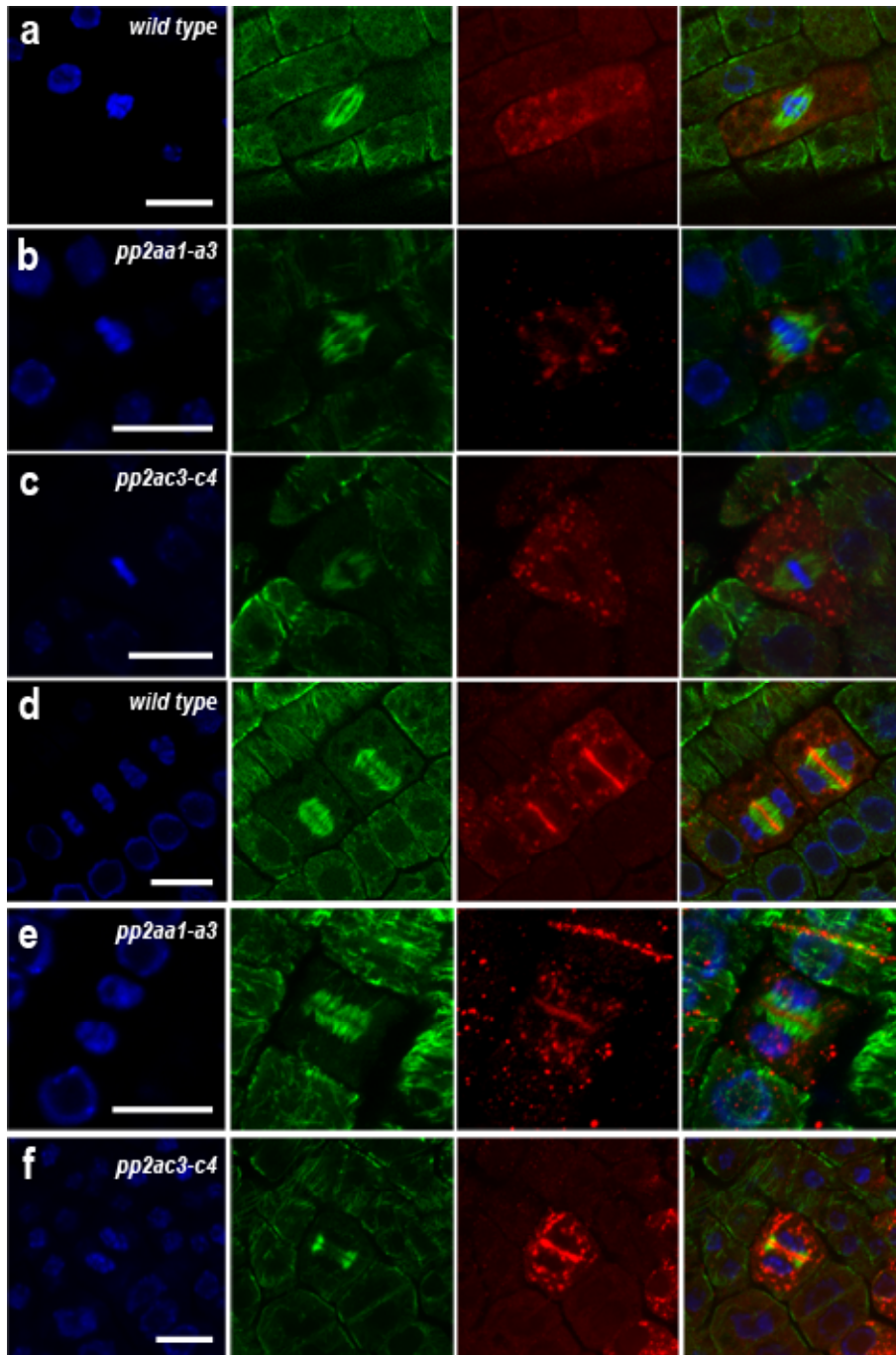


### Supplementary Figure S5 : Cell division plane defects in roots of *pp2aaa* and *pp2ac* mutants

Longitudinal (upper panels), transverse (mid panels), and tangential views of root tips of wild-type (Col-0), *pp2aa1*, *pp2aa2-a3*, *pp2aa1-a2*, *pp2aa1-a3*, *pp2ac3-c4* three-day old seedlings stained with propidium iodide.

Bars = 50  $\mu$ m in all images.

## Supplementary Figure S6

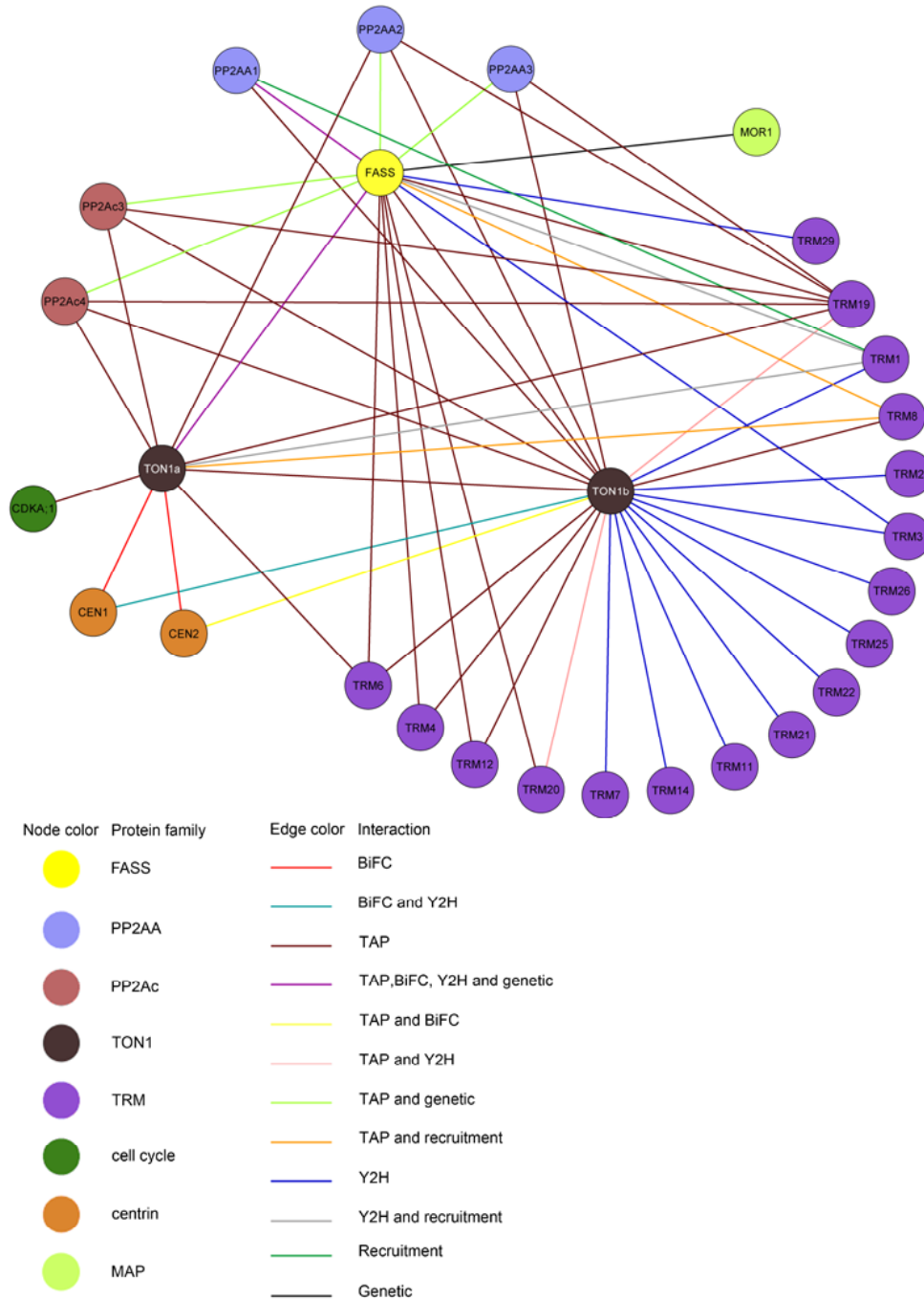


### Supplementary Figure S6 : *pp2aa* and *pp2ac* mutations don't affect spindle and phragmoplast

(a-c) Spindles and (d-f) phragmoplasts. Whole mount immuno-localization of  $\alpha$ -tubulin, and KNOLLE in root tip cells of wild type (a and d), *pp2aa1-a3* (b and e), *pp2ac3-c4* (c and f) double mutants. Here are shown DAPI staining (in blue), anti- $\alpha$ -tubulin (in green), anti-KNOLLE (red), and overlay images. Bars = 10 $\mu$ m in a-f.



## Supplementary Figure S7



### Supplementary Figure S7: Interactome map overview of all TTP protein interactions

This protein-protein interaction network includes all known interactions concerning the TTP proteins. The presented data is a combination of novel data described in this paper and previously published interactions<sup>10,11,12,15,25,43</sup>. The nodes represent individual proteins and the node colors are indicative of the different protein families. The lines connecting the nodes indicate known interactions and are color-coded dependent on the techniques used to identify the interactions.

# Supplementary Table S1 : MALDI MS/MS analysis of TAP eluates

Bait	TAP Protocol	Prev				Database version	PMF data					MSMS data										Variable Modification
		Prey Locus	Prey Name	# Found/ # exp	Protein Score		Expect	RMS error (ppm)	Sequence coverage %	Unique Peptides	Total Ion Score	Peptide Number	Start	End	Observed	Mr(Exp)	Mr(Calc)	Delta	Miss	Ions Score	Expect	
TON1a	-dig	AT3G55005	TON1b	4/4	TAIR8	156	8,20E-12	4	42	8	106	1	237	243	863.44	862.43	862.43	0.00	0	32	2,00E-02	R.NLTSSWR.N
												2	34	45	1394.64	1393.64	1393.64	0.00	0	52	1,00E-04	R.ASVFEAIEEDR.V
												3	126	142	1958.05	1957.04	1957.05	-0.01	0	22	7,40E-02	R.NEDSRPLLDVLEGLK.F
TON1a	-dig	AT5G18580	FASS	4/4	TAIR8	207	6,50E-17	6	25	11	146	1	244	252	1272.56	1271.55	1271.55	0.00	0	15	2,80E-01	K.FFFFCDPHR.R
												2	369	379	1311.65	1310.64	1310.65	-0.01	0	50	2,50E-04	K.DTPEGLTYLFR.C
												3	18	29	1384.77	1383.76	1383.77	-0.01	1	12	9,70E-01	R.KIPPASSMLWVR.N
												4	387	399	1477.77	1476.76	1476.77	-0.01	0	25	7,20E-02	R.GFLTADHSLFR.D
												5	321	334	1656.81	1655.8	1655.8	0.00	0	44	7,30E-04	K.EYADGTLTEIFIER.V
TON1a	-dig	AT3G25800	PP2AA2	1/4	TAIR8	58	5,10E-02	12	32	12	/											
TON1a	-dig	AT2G42500/AT3	PP2A3/PP2A4	1/4	TAIR8	68	4,60E-03	7	28	6	34	1	260	272	1589.78	1588.78	1588.77	0.01	0	34	9,40E-03	K.VVTIFSAFNQYCR.C
TON1a	-dig	AT2G45240	MAP1A	1/4	TAIR8	47	6,50E-01	4	14	4	31	1	92	102	1310.69	1309.68	1309.68	0.00	0	31	1,70E-02	K.LPHFDWTGTLK.Q
TON1a	+dig	AT3G55005	TON1b	4/4	TAIR8	441	2,60E-40	5	64	14	321	1	237	243	863.44	862.43	862.43	0.00	0	41	2,40E-03	R.NLTSSWR.N
												2	222	233	1330.7	1329.69	1329.69	0.00	0	40	2,50E-03	R.ASTALENLQLDR.K
												3	114	125	1441.66	1440.66	1440.66	0.00	0	83	7,40E-08	R.DFSINNGYELNR.N
												4	187	202	1499.7	1498.7	1498.7	0.00	0	66	3,70E-06	R.SVSASQASGAATSGYR.K
												5	126	142	1958.06	1957.05	1957.05	0.00	0	92	7,60E-09	R.NEDSRPLLDVLEGLK.F
TON1a	+dig	AT5G18580	FASS	4/4	TAIR8	390	3,30E-35	4	35	16	274	1	335	341	901.45	900.44	900.45	-0.01	0	24	5,30E-02	R.VFDEHVR.R
												2	457	464	1002.46	1001.45	1001.45	0.00	0	30	9,80E-03	R.GFWAHDNR.E
												3	19	29	1256.69	1255.68	1255.67	0.01	0	19	2,50E-01	K.IPPASSMLWVR.N
												4	244	252	1272.57	1271.57	1271.55	0.02	0	28	1,50E-02	K.FFFFCDPHR.R
												5	369	379	1311.66	1310.65	1310.65	0.00	0	25	7,10E-02	K.DTPEGLTYLFR.C
												6	18	29	1384.77	1383.77	1383.77	0.00	1	0	1,40E+01	R.KIPPASSMLWVR.N
												7	387	399	1477.78	1476.78	1476.77	0.01	0	59	2,50E-05	R.GFLTADHSLFR.D
												8	321	334	1656.82	1655.81	1655.8	0.01	0	52	1,30E-04	K.EYADGTLTEIFIER.V
												9	101	119	2204.08	2203.07	2203.08	-0.01	0	36	3,90E-03	K.QSDLLNADDLAAMWVCLR.E
TON1a	+dig	AT3G25800	PP2AA2	2/4	TAIR8	111	2,60E-07	9	33	13	31	1	279	287	1047.58	1046.57	1046.58	-0.01	0	1	2,30E+01	R.TELVPAYVR.L
TON1a	+dig	AT2G42500/AT3	PP2A3/PP2A4	2/4	TAIR8	120	3,30E-08	4	33	8	67	2	133	142	1121.57	1120.57	1120.57	0.00	0	30	2,80E-02	R.LAAGEWFTAR.V
TON1a	+dig	AT3G06810	IBR3	1/4	TAIR8	84	1,40E-04	6	28	17	/	1	287	298	1456.72	1455.72	1455.73	-0.01	0	66	4,20E-06	R.NHTFIQEPAPR.R
TON1b	-dig	AT3G55000	TON1a	5/5	TAIR8	622	2,10E-58	10	51	15	487	1	241	247	863.43	862.43	862.43	0.00	0	26	8,30E-02	R.NLTSSWR.N
												2	67	75	952.53	951.52	951.52	0.00	1	35	6,00E-03	R.KLHASPSGR.L
												3	207	213	994.48	993.47	993.47	0.00	1	38	3,20E-03	R.KDEFNWR.Q
												4	175	185	1144.57	1143.56	1143.55	0.01	0	33	1,00E-02	R.SSASDSLPPQR.R
												5	226	237	1300.69	1299.68	1299.68	0.00	0	58	3,40E-05	R.ASAALENLQLDR.K
												6	34	45	1394.68	1393.67	1393.64	0.03	0	61	1,80E-05	R.ASVFEAIEEDR.V
												7	114	125	1426.65	1425.65	1425.63	0.02	0	85	5,20E-08	R.DFNSNNGFELNR.N
												8	126	142	1786.97	1785.96	1785.95	0.01	0	87	2,60E-08	R.NGDSGLLDVLEGLK.F
												9	110	125	1911.92	1910.91	1910.89	0.02	1	23	1,00E-01	K.SELRDFNSNNGFELNR.N
TON1b	-dig	AT5G18580	FASS	5/5	TAIR8	568	5,20E-53	4	45	23	353	1	335	341	901.45	900.44	900.45	-0.01	0	33	8,50E-03	R.VFDEHVR.R
												2	457	464	1002.46	1001.45	1001.45	0.00	0	30	9,20E-03	R.GFWAHDNR.E
												3	156	164	1120.52	1119.51	1119.51	0.00	0	31	1,10E-02	R.FFFSPSNFMK.F
												4	244	252	1272.57	1271.56	1271.55	0.01	0	21	7,30E-02	K.FFFFCDPHR.R
												5	369	379	1311.66	1310.65	1310.65	0.00	0	40	2,10E-03	K.DTPEGLTYLFR.C
												6	66	77	1327.66	1326.65	1326.65	0.00	0	18	3,20E-01	K.SQEAGTIPSPFYK.K
												7	444	456	1346.71	1345.7	1345.7	0.00	0	38	4,00E-03	K.QGGTVASMLIDVR.G
												8	18	29	1384.77	1383.77	1383.77	0.00	1	30	1,40E-02	R.KIPPASSMLWVR.N
												9	18	29	1400.77	1399.76	1399.76	0.00	1	(2)	1,30E+01	R.KIPPASSMLWVR.N
												10	387	399	1477.78	1476.77	1476.77	0.00	0	50	2,30E-04	R.GFLTADHSLFR.D
												11	173	184	1514.84	1513.84	1513.84	0.00	0	3	7,20E+00	R.IAILPPYLYVMR.T
												12	227	238	1529.64	1528.63	1528.63	0.00	0	28	6,20E-03	R.DMPFANQMYCR.I
												13	321	334	1656.81	1655.8	1655.8	0.00	0	31	1,90E-02	K.EYADGTLTEIFIER.V
TON1b	-dig	AT3G25800	PP2AA2	4/5	TAIR8	225	1,00E-18	6	36	15	129	1	279	287	1047.58	1046.57	1046.58	-0.01	0	27	5,20E-02	R.TELVPAYVR.L
												2	133	142	1121.57	1120.57	1120.57	0.00	0	38	4,20E-03	R.LAAGEWFTAR.V
												3	537	552	1787.02	1786.01	1786.02	-0.01	0	23	2,20E-02	K.VLQSLIPVDQSVVEK.T
												4	310	325	1858.06	1857.06	1857.05	0.01	0	30	6,70E-03	R.IILNPEIAIHLPCVK.E
TON1b	-dig	AT2G42500/AT3	PP2A3/PP2A4	3/5	TAIR8	218	5,20E-18	5	45	11	131	5	259	278	2340.11	2339.1	2339.09	0.01	0	11	1,40E+00	R.YMVANQLYELCEAVGPEPTR.T
												1	287	298	1456.72	1455.71	1455.73	-0.02	0	34	7,20E-03	R.NHTFIQEPAPR.R
												2	260	272	1589.78	1588.77	1588.77	0.00	0	42	1,30E-03	K.VVTIFSAFNQYCR.C
												3	219	239	2341.09	2340.08	2340.08	0.00	0	35	3,80E-03	R.GAGYTFQDISEQFNHTNKL.L
												4	54	74	2412.18	2411.17	2411.17	0.00	0	20	1,60E-01	K.SPVTICGDIHQFDLAEI.LFR.I
TON1b	-dig	AT1G13320	PP2AA3	1/5	TAIR8	74	1,40E-03	10	13	6	50	1	133	142	1137.58	1136.57	1136.56	0.01	0	34	1,20E-02	R.LSAGEWFTAR.V
												2	259	278	2340.11	2339.1	2339.09	0.01	0	16	3,70E-01	R.YMVANQLYELCEAVGPEPTR.T
												3	259	278	2356.1	2355.1	2355.09					





Protein identification details were obtained with the 4800 MALDI TOF/TOF<sup>TM</sup> Proteomics analyzer (AB SCIEX) and the GPS explorer v3.6 (AB SCIEX) software package combined with search engine Mascot version 2.1 or 2.2 (Matrix Science)

Column headers for Protein and Peptide data are as follows :

**Protein score:** The score calculated by the Mascot search engine for each protein. This score is based on the probability that peptide mass matches are non-random events. If the Protein Score is equal to or greater than the Mascot® Significance Level calculated for the database search, the protein match is considered to be statistically non-random at the 95% confidence interval. Protein score =  $-10 \cdot \log(P)$ , where P is the probability that the observed match is a random event. **Expect:** Protein score expectation value. **RMS error (ppm):** RMS error of the set of matched mass values, in ppm. **Sequence coverage %:** Percentage of protein sequence covered by assigned peptide matches. **Unique peptides:** The number of peptides with unique sequences matching the selected protein. **Total Ion Score:** A score calculated by weighting Ion Scores for all individual peptides matched to a given protein. **Peptide Number:** Peptide index number within the list of peptides associated with a given protein. **Start:** The starting position of the peptide in the protein. **End:** The ending position of the peptide in the protein. **Observed:** The observed monoisotopic mass of the peptide in the spectrum (m/z). **Mr (Exp):** The experimental mass of the peptide calculated from the observed m/z value. **Mr (Calc):** The theoretical mass of the peptide based on its sequence. **Delta (Da):** The difference between the theoretical (Mr (Calc)) and experimental (Mr (Exp)) masses, in daltons. **Miss:** Number of missed Trypsin cleavage sites. **Ions score:** The Ions Score is calculated by the Mascot search engine for each peptide matched from MS/MS peak lists. This score is based on the probability that ion fragmentation matches are non-random events. If the Ion Score is equal to or greater than the Mascot® Significance Level calculated for the database search, the peptide match is considered to be statistically non-random at the 95% confidence interval. Ions score =  $-10 \cdot \log(P)$ , where P is the probability that the observed match is a random event. **Best Ions score:** The highest individual Ion Score for a given protein identification. **Expect:** Ions score expectation value. **Peptide:** The amino acid sequence of the selected peptide. **Variable Modification:** Variable modification type on the peptide.

**Supplementary Table S2: VELOS orbitrap analysis of TAP TON1b eluates**

Accession	Description	Score	Coverage	# Unique Peptides	# Peptides	# AAs					
<b>AT3G55005</b>	<b>TON1b</b>	<b>237,09</b>	<b>64,20</b>	<b>17</b>	<b>18</b>	<b>257</b>					
	Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	$\Delta M$ [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
	High	NEDSRPLLDVLEGLK		1	4,69	3	1958,06035	0,47	653,35830	28,73	0
	High	NEDSRPLLDVLeGFLK	E13(Methyl)	1	4,48	3	1972,07370	-0,69	658,02942	29,64	0
	High	NEDSRPLLDVLEGLKFENmTQVMGGSSR	M21(Oxidation)	1	4,48	4	3398,66884	-0,95	850,42267	29,10	1
	High	ASVFEAIEEEDRVIEENEGPLPALLGScNDR	C28(Carbamidomethyl)	1	4,43	4	3443,63776	-0,23	861,66490	26,02	1
	High	SEIRDFSIINNGYeLNR	E13(Methyl)	1	4,13	3	1940,94687	0,37	647,65381	21,77	1
	High	RSVSASQASGAATSGYR		1	4,08	2	1655,80777	-1,14	828,40752	17,14	1
	High	FENmTQVMGGSSR	M4(Oxidation); M8(Oxidation)	1	3,79	2	1475,62496	-0,32	738,31612	17,30	0
	High	VIeNNeGLPALLGScNDR	E3(Methyl); E6(Methyl); C16(Carbamidomethyl)	1	3,74	2	2096,04546	0,61	1048,52637	23,56	0
	High	LLSALlCeYLDWAQLNHTLK	C7(Carbamidomethyl)	1	3,68	2	2401,25728	-0,51	1201,13228	27,13	0
	High	YDTEdMPEEVmR	M6(Oxidation); M11(Oxidation)	1	3,65	2	1546,60247	-0,78	773,80487	18,49	0
	High	ASVFEAIEEEDRVIEENeGLPALLGScNDR	E18(Methyl); C28(Carbamidomethyl)	1	3,60	3	3457,65342	-0,22	1153,22266	27,17	1
	High	ASTALENLQLDRK		1	3,58	2	1458,78997	-0,81	729,89862	20,22	1
	High	FeNMTQVMGGSSR	E2(Methyl)	1	3,55	2	1457,65068	-0,38	729,32898	21,93	0
	High	SEIRDFSIINNGYeLNR		1	3,55	3	1926,93133	0,43	642,98196	21,38	1
	High	VIENNEGLPALLGScNDR	C16(Carbamidomethyl)	1	3,54	3	2068,01265	-0,11	690,00907	22,47	0
	High	ReSETESSSLDTR	E2(Methyl)	1	3,53	2	1623,78048	-1,00	812,39388	19,08	1
	High	NedSRPLLDVLEGLK	E2(Methyl); D3(Methyl)	1	3,48	2	1986,08820	-1,27	993,54774	29,20	0
	High	ESETESSSLDTR		1	3,46	2	1453,66570	0,25	727,33649	20,03	0
	High	ASTALENLQLDR		1	3,39	2	1330,69585	-0,25	665,85156	21,68	0
	High	SVSASQASGAATSGYR		1	3,39	2	1499,70744	-0,73	750,35736	18,20	0
	High	eSETESSSLDTR	E1(Methyl)	1	3,38	2	1467,68340	1,65	734,34534	20,09	0
	High	FENmTQVMGGSSR	M4(Oxidation)	1	3,37	2	1459,62924	-0,87	730,31826	18,99	0
	High	YDTEdMPEeVMR	M6(Oxidation); E9(Methyl)	1	3,23	2	1544,62241	-1,30	772,81485	20,28	0
	High	ASTALENLQLDR	E6(Methyl)	1	3,20	2	1344,71044	-1,03	672,85886	21,40	0
	High	DFSINNGYeLNR		1	3,20	2	1441,67035	-0,23	721,33881	22,28	0
	High	YDTEdMPEEVmR		1	3,14	2	1514,61221	-1,09	757,80974	21,49	0
	High	SSASDSLPHQR		1	3,14	2	1184,56520	-0,24	592,78624	16,35	0
	High	eSeTESSSLDTR	E1(Methyl); E3(Methyl)	1	3,13	2	1481,69475	-1,27	741,35101	21,14	0
	High	YDTEdMPEeVMR	M6(Oxidation); E8(Methyl); E9(Methyl)	1	3,11	2	1558,63833	-1,12	779,82280	20,89	0
	High	YDTEdMPEEVmR	M6(Oxidation)	1	2,96	2	1530,60719	-1,03	765,80723	19,74	0
	High	ASVFEAIEEEdR	D11(Methyl)	1	2,96	2	1408,65850	-0,45	704,83289	24,03	0
	High	ASVFEAIEEEDR		1	2,92	3	1394,64457	0,77	465,55304	23,22	0
	High	ASVFEAIEeEDRVIEENEGPLPALLGScNDR	E9(Methyl); E10(Methyl); C28(Carbamidomethyl)	1	2,91	4	3471,66665	-0,92	868,67212	26,83	1
	High	RSSASDSLPHQR		1	2,88	2	1340,66502	-1,18	670,83615	15,96	1
	High	YDTEdMPEeVmR	M6(Oxidation); E8(Methyl); E9(Methyl); M11(Oxidation)	1	2,87	2	1574,63445	-0,34	787,82086	19,71	0
	High	YDTEdMPEEVmR	E4(Methyl); M6(Oxidation); M11(Oxidation)	1	2,85	2	1560,61769	-1,05	780,81248	18,77	0
	High	RESeTESSSLDTR	E4(Methyl); E6(Methyl)	1	2,83	3	1637,79674	-0,62	546,60376	19,58	1
	High	FeNmTQVMGGSSR	E2(Methyl); M4(Oxidation)	1	2,82	2	1473,64531	-0,57	737,32629	19,37	0
	High	VIENNeGLPALLGScNDR	E6(Methyl); C16(Carbamidomethyl)	1	2,81	3	2082,02652	-0,97	694,68036	22,75	0
	High	DFSINNGYeLNR	E9(Methyl)	1	2,80	2	1455,68541	-0,63	728,34634	22,51	0
	High	RESETESSSLDTR		1	2,80	2	1609,76513	-0,82	805,38620	19,23	1
	High	YDTEdMPEEVmR	E4(Methyl)	1	2,52	2	1528,62847	-0,68	764,81787	22,07	0
	High	ASVFEAIEeEDR	E8(Methyl); E9(Methyl); E10(Methyl)	1	2,39	2	1436,68880	-1,14	718,84804	25,31	0
<b>AT3G55000</b>	<b>TON1a</b>	<b>129,58</b>	<b>60,00</b>	<b>13</b>	<b>14</b>	<b>260</b>					
	Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	$\Delta M$ [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
	High	ASVFEAIEEEDRVIEENEGPPALLGScNDR	C28(Carbamidomethyl)	1	5,27	3	3427,60361	-1,06	1143,20605	25,47	1
	High	LLSALlCeYLDWAQLNHTLlVYQPESNLPK	C7(Carbamidomethyl)	1	4,89	3	3541,84348	0,96	1181,28601	28,02	0
	High	NGDSGPLLDVLeGFLK	E13(Methyl)	1	4,82	2	1800,97258	-0,95	900,98993	31,08	0
	High	NGDSGPLLDVLEGLK		1	4,53	3	1786,95831	-0,18	596,32429	30,35	0
	High	SELRDFNSNNGFELNR		1	4,47	3	1911,89481	0,18	637,96979	21,01	1
	High	VIeNNeGPPALLGScNDR	E3(Methyl); C16(Carbamidomethyl)	1	4,09	2	2065,99785	0,30	1033,50256	21,50	0
	High	DSETESSSSLESR		1	3,99	2	1413,59636	-0,91	707,30182	17,68	0

High	ASVFEAIEEEDRVIEENEGPPALLGScNDR	E15(Methyl); C28(Carbamidomethyl)	1	3,95	3	3441,62063	-0,66	1147,87839	25,03	1
High	RDESETSSSSLESr		1	3,33	3	1569,69815	-0,39	523,90424	16,67	1
High	VIENNEGPPALLGScNDR	C16(Carbamidomethyl)	1	3,32	2	2051,97952	-1,01	1026,49340	21,15	0
High	SSASDSLPPQR		1	3,31	2	1144,55866	-0,60	572,78297	17,35	0
High	RDESETSSSSLeSR	E12(Methyl)	1	3,31	3	1583,71410	-0,20	528,57622	16,82	1
High	FESmTQmGSSSR	M4(Oxidation); M8(Oxidation)	1	3,30	2	1436,57681	-0,92	718,79204	16,35	0
High	DFNSNNGFELNR		1	3,28	2	1426,63337	-0,89	713,82032	21,74	0
High	RSSASDSLPPQR		1	3,24	2	1300,65837	-1,61	650,83282	16,58	1
High	dSETSSSSLESr	D1(Methyl)	1	3,09	2	1427,61283	-0,32	714,31005	17,43	0
High	RPVSASQASDR		1	2,98	2	1173,59650	-0,53	587,30189	15,45	0
High	ASVFEAIEEEdR	D11(Methyl)	1	2,96	2	1408,65850	-0,45	704,83289	24,03	0
High	ASVFEAIEEEDR		1	2,92	3	1394,64457	0,77	465,55304	23,22	0
High	FESMTQmGSSSR	M8(Oxidation)	1	2,87	2	1420,58199	-0,87	710,79463	17,67	0
High	QGNQDTHeeVTR	E9(Methyl)	1	2,53	2	1427,64940	-1,13	714,32834	16,28	0
High	ASVFEAIEeeDR	E8(Methyl); E9(Methyl); E10(Methyl)	1	2,39	2	1436,68880	-1,14	718,84804	25,31	0
<b>AT5G18580</b>	<b>FASS</b>	<b>127,05</b>	<b>67,50</b>	<b>22</b>	<b>22</b>	<b>480</b>				

Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	ΔM [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
High	EmDFDSFLDFVLALENK	M2(Oxidation)	1	4,90	2	2048,94927	-1,45	1024,97827	30,00	0
High	QELKEYADGTLTEIFIER		1	4,89	3	2155,09128	-0,26	719,03528	24,03	1
High	MNYEDFCHIASVcTEQIGPK	C7(Carbamidomethyl); C13(Carbamidomethyl)	1	4,82	3	2399,04288	-1,53	800,35248	22,80	0
High	YIGSGAGLGSEALMELETK		1	4,44	2	1925,95305	0,25	963,48016	24,17	0
High	KKPEEGSISQR		1	4,27	3	1258,67521	0,12	420,22992	15,40	1
High	mNYEDFCHIASVcTEQIGPK	(Oxidation); C7(Carbamidomethyl); C13(Carbamidomethyl)	1	3,99	3	2415,03824	-1,34	805,68427	21,90	0
High	IcDmFLALDKDmSGSLcK	domethyl); M4(Oxidation); M12(Oxidation); C17(Carba	1	3,93	2	2135,94780	-0,15	1068,47754	21,84	1
High	QSDLLNADDLAAMWvcLR	C17(Carbamidomethyl)	1	3,93	3	2204,08261	-0,61	735,36572	28,42	0
High	YIGSGAGLGSEALMeLETK	M14(Oxidation); E15(Methyl)	1	3,86	2	1955,96282	-0,16	978,48505	24,45	0
High	YIGSGAGLGSEALmELETK	M14(Oxidation)	1	3,78	2	1941,94621	-0,65	971,47675	23,73	0
High	IDmSELDESDGFLHSDemESYIGGLIPNLAQLR	M3(Oxidation)	1	3,70	3	3825,73899	-2,14	1275,91785	27,76	0
High	IcDMFLALDKDmSGSLcK	Carbamidomethyl); M12(Oxidation); C17(Carbamidome	1	3,64	2	2119,94902	-1,97	1060,47815	23,11	1
High	KIPPASmLWVR	M8(Oxidation)	1	3,51	3	1400,77150	-0,32	467,59535	20,34	1
High	ILLSNLQELMELHQESEEEVTDTEQAENWFLSTSAQR	C6(Carbamidomethyl); M11(Oxidation)	1	3,44	4	4524,07680	-1,13	1131,77466	26,34	0
High	WIEGGNYELcIEDVR	C10(Carbamidomethyl)	1	3,43	2	1852,85307	-0,23	926,93018	23,67	0
High	QGGTVASmLIDVR	M8(Oxidation)	1	3,37	2	1362,70378	-0,64	681,85553	21,77	0
High	IDmSELDESDGFLHSDemESYIGGLIPNLAQLR	M3(Oxidation); M19(Oxidation)	1	3,35	3	3841,74643	1,13	1281,25366	26,33	0
High	IAILPFYLYVmR	M11(Oxidation)	1	3,34	2	1514,84295	-0,71	757,92511	26,65	0
High	DmPPAFNQMyCr	M2(Oxidation); C11(Carbamidomethyl)	1	3,21	2	1545,62773	-0,43	773,31750	21,30	0
High	IAILPFYLYVMR		1	3,16	2	1498,84734	-1,18	749,92731	27,98	0
High	DmPPAFNQmYcR	M2(Oxidation); M9(Oxidation); C11(Carbamidomethyl)	1	3,13	2	1561,62419	0,57	781,31573	19,81	0
High	GFWAHDNRENLLQEEEEPEEESQ		1	3,13	3	2912,25840	-0,78	971,42432	22,16	1
High	mNYeDFCHIASVcTEQIGPK	tion); E4(Methyl); C7(Carbamidomethyl); C13(Carbami	1	3,03	3	2429,05442	-1,11	810,35632	22,33	0
High	SQEAGTIPSFYK		1	2,99	2	1327,65325	0,26	664,33026	21,27	0
High	IcDMFLALDK	C2(Carbamidomethyl)	1	2,92	2	1225,59514	-0,39	613,30121	24,30	0
High	YIGSGAGLGSEALMeLeTK	M14(Oxidation); E15(Methyl); E17(Methyl)	1	2,89	2	1969,97820	-0,30	985,49274	25,07	0
High	EYADGTLTEIFIER		1	2,89	3	1656,81076	-0,51	552,94177	24,89	0
High	ITLGDLLGcK	C9(Carbamidomethyl)	1	2,69	2	1089,59714	-0,16	545,30221	24,73	0
High	ILLEIFK		1	2,63	2	875,55997	-0,17	438,28362	24,56	0
High	QGGTVASMLIDVR		1	2,61	2	1346,70891	-0,62	673,85809	23,55	0
High	IcDmFLALDK	C2(Carbamidomethyl); M4(Oxidation)	1	2,56	2	1241,58916	-1,10	621,29822	22,61	0
High	ENCVIDDATGAeK	C3(Carbamidomethyl); E12(Methyl)	1	2,46	2	1435,63603	-0,69	718,32166	20,31	0
High	KPEEGSISQR		1	2,42	2	1130,57930	-0,69	565,79329	15,93	0
<b>AT3G25800</b>	<b>PP2AA2</b>	<b>98,25</b>	<b>44,63</b>	<b>10</b>	<b>16</b>	<b>587</b>				

Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	ΔM [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
High	LAIIIEYIPLLASQLGVGFDDK		1	5,62	3	2422,32859	0,68	808,11438	31,48	0
High	YmVANQLYLcEAVGPEPTR	M2(Oxidation); C11(Carbamidomethyl)	1	4,92	2	2356,09331	-0,67	1178,55029	23,66	0
High	TDVMsmFEDLTQDDQDSVR	M6(Oxidation)	1	4,57	2	2247,93486	-1,46	1124,47107	24,43	0
High	FFANQALQSIDNVmMSS	M15(Oxidation)	1	4,40	2	1918,87187	2,30	959,93958	27,11	0
High	FFANQALQSIDNVmMSS	M14(Oxidation); M15(Oxidation)	1	4,10	2	1934,86165	-0,37	967,93446	24,43	0
High	SLYTQLcQDDmPmVR	C7(Carbamidomethyl); M11(Oxidation); M13(Oxidation)	1	4,07	2	1888,82247	-0,74	944,91487	21,55	0
High	SALASVImGmAPVLGK	M8(Oxidation); M10(Oxidation)	1	4,00	3	1576,84348	-0,22	526,28601	22,87	0

High	LAIIEYIPLLASQLGVGFDDK	D21(Methyl)	1	3,98	3	2436,33982	-1,14	812,78479	31,43	0
High	TIRPGLVELSEDPDVR		1	3,82	2	2010,05022	-0,05	1005,52875	21,89	0
High	ILNPEIAIQHILPcVK	C14(Carbamidomethyl)	1	3,78	2	1858,06035	-0,88	929,53381	24,89	0
High	VLQSLIPIVDQSVVEK		1	3,76	3	1767,02597	-0,19	589,68018	24,53	0
High	SALASVImGMAPVLGK	M8(Oxidation)	1	3,65	2	1560,84636	-1,63	780,92682	25,63	0
High	TIRPGLVELSeDPDVR	E11(Methyl)	1	3,44	3	2024,06595	-0,01	675,36017	22,51	0
High	LLEPQDcVQHILPVIvNFSQDK	C7(Carbamidomethyl)	1	3,32	3	2592,34824	-0,35	864,78760	25,31	0
High	AVSLAPVmgSEITcSK	M9(Oxidation); C15(Carbamidomethyl)	1	3,31	2	1778,90068	-1,18	889,95398	22,42	0
High	ESLDVDFHLSLVK		1	3,29	2	1501,78813	-1,08	751,39771	25,17	0
High	LLRDNEAEVR		1	2,93	2	1214,64739	-1,20	607,82733	17,65	1
High	VSAcGVFHIAyPSAPDmLK	C4(Carbamidomethyl); M17(Oxidation)	1	2,62	2	2079,00127	-1,26	1040,00427	22,52	0
High	FAATVeSAHLK	E6(Methyl)	1	2,61	2	1187,64128	-0,56	594,32428	19,47	0
High	ELSSDSSQHVR		1	2,59	2	1244,58501	-1,29	622,79614	16,21	0
<b>AT3G06810.1</b>	<b>IBR3</b>	<b>86,92</b>	<b>36,65</b>	<b>19</b>	<b>19</b>	<b>824</b>				

Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	ΔM [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
High	VSQFGHGQSNPTFLIEVGSGLSKR		1	4,98	3	2632,35056	1,01	878,12170	21,69	1
High	SGFAmTEPQVASSDATNIEcSIR	M5(Oxidation); C20(Carbamidomethyl)	1	4,74	2	2487,11235	-0,16	1244,05981	21,82	0
High	ALASLHSADVDAIGLEK		1	4,56	2	1709,90544	-0,87	855,45636	21,30	0
High	IADGPDEVHLGTIGK		1	4,36	2	1521,78942	-0,93	761,39835	20,50	0
High	NTGVQANELVESALGYIAR		1	4,26	2	2005,03667	0,84	1003,02197	26,06	0
High	NIPAEDSTGATSGLVHGDFR		1	4,14	2	2043,97295	-0,08	1022,49011	20,79	0
High	FAADNVSGFPTNPSQFK		1	4,09	2	1826,86897	-1,04	913,93812	22,32	0
High	YGNKEQISEWLIPLLEGR		1	3,92	3	2145,13547	0,70	715,71667	26,20	1
High	EGLWNLFPVDSAAR		1	3,90	2	1673,86363	-0,62	837,43545	27,07	0
High	ELAATENKHNLGK		1	3,76	2	1511,77995	-0,90	756,39362	16,56	1
High	SFDQLFGEGLTNLEYGlcEImGR	C19(Carbamidomethyl); M22(Oxidation)	1	3,76	3	2828,28916	-0,55	943,43457	28,54	0
High	FIAQHGSFVSDLAk		1	3,76	2	1519,78874	-1,10	760,39801	21,01	0
High	VSQFGHGQSNPTFLIEVGSGLSK		1	3,39	3	2476,24308	-1,50	826,08588	22,61	0
High	SWWAPQVfNcGAPDTGNmEVILR	C10(Carbamidomethyl); M18(Oxidation)	1	3,19	3	2577,21970	-1,12	859,74475	24,01	0
High	FmETHIYPmENEFSK	M2(Oxidation); M9(Oxidation)	1	2,99	2	1934,82915	-0,44	967,91821	20,30	0
High	LLVLEAADHLDK		1	2,98	2	1336,74687	-0,22	668,87708	22,20	0
High	WlmgNASAGER	M3(Oxidation)	1	2,90	2	1207,55156	-0,78	604,27942	19,23	0
High	QGDSYVINGTK		1	2,78	2	1181,57939	-0,29	591,29333	18,90	0
High	mFELVDWLR	M1(Oxidation)	1	2,67	2	1224,60784	-0,32	612,80756	26,26	0
High	FMETHIYPmENEFSK	M9(Oxidation)	1	2,63	2	1918,83464	-0,23	959,92096	21,06	0
<b>AT3G53540</b>	<b>TRM19</b>	<b>48,20</b>	<b>19,16</b>	<b>12</b>	<b>12</b>	<b>924</b>				

Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	ΔM [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
High	SHSSYNSPEVSITPSSLK		1	4,62	2	2006,96501	-0,79	1003,98615	19,96	0
High	EFNDALEALDSNKDLLK		1	4,55	2	2048,05571	0,48	1024,53149	24,33	1
High	ELQWLSLEDDIEIIGR		1	4,09	2	1928,99358	-1,51	965,00043	27,33	0
High	VNANLTQAEMAFIR		1	3,84	2	1577,80950	-0,65	789,40839	23,49	0
High	HASYDTIDLpNEELR		1	3,82	2	1772,84490	-0,08	886,92609	21,28	0
High	TAmSSEALDLSTVTSVTDpDISR	M3(Oxidation)	1	3,59	2	2412,14263	-0,88	1206,57495	23,27	0
High	LmGLDVLPSQSSSHK	M2(Oxidation)	1	3,57	2	1614,81304	-1,62	807,91016	20,74	0
High	LTHKFEHEIEISR		1	3,28	4	1638,85903	-0,53	410,47021	18,39	1
High	KLLFDQISR		1	2,92	3	1119,65165	-0,42	373,88873	21,21	1
High	TFASPSSSDFEFR		1	2,87	2	1417,62322	0,10	709,31525	20,06	0
High	FLQHPDSLFTK		1	2,74	2	1332,69646	1,31	666,85187	21,15	0
High	FKDVFVLDLAK		1	2,62	2	1310,69878	-0,28	655,85303	23,37	1
<b>AT1G13320</b>	<b>PP2AA3</b>	<b>45,71</b>	<b>26,92</b>	<b>3</b>	<b>10</b>	<b>587</b>				

Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	ΔM [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
High	YmVANQLYELcEAVGPEPTR	M2(Oxidation); C11(Carbamidomethyl)	1	4,92	2	2356,09331	-0,67	1178,55029	23,66	0
High	SALASVImGmAPVLGK	M8(Oxidation); M10(Oxidation)	1	4,00	3	1576,84348	-0,22	526,28601	22,87	0
High	LAIIEYIPLLASQLGVGFDEK		2	3,98	3	2436,33982	-1,14	812,78479	31,43	0
High	SIYGQLcQDDmPmVr	C7(Carbamidomethyl); M11(Oxidation); M13(Oxidation)	1	3,84	2	1844,79436	-1,79	922,90082	21,03	0
High	ILNPEIAIQHILPcVK	C14(Carbamidomethyl)	1	3,78	2	1858,06035	-0,88	929,53381	24,89	0
High	SALASVImGMAPVLGK	M8(Oxidation)	1	3,65	2	1560,84636	-1,63	780,92682	25,63	0
High	YFANQLQSDINvmSS	M14(Oxidation); M15(Oxidation)	1	3,39	2	1950,85198	-2,71	975,92963	23,81	0
High	AVSLAPVmgSEITcSK	M9(Oxidation); C15(Carbamidomethyl)	1	3,31	2	1778,90068	-1,18	889,95398	22,42	0



	High	ESDLVEHFTPLAK		1	2,65	2	1485,75713	-0,89	743,38220	21,97	0
	High	ELSSDSSQHVR		1	2,59	2	1244,58501	-1,29	622,79614	16,21	0
	High	FAATIESAHLK		2	2,55	2	1187,64128	-0,56	594,32428	19,47	0
<b>AT3G58500</b>	<b>PP2A-4</b>	<b>41,53</b>	<b>45,37</b>	<b>10</b>	<b>10</b>	<b>313</b>					
	Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	ΔM [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
	High	GAGYTFGQDISEQFNHTNNLK		1	4,78	2	2341,08257	-0,80	1171,04492	22,30	0
	High	cPDTNYLFmGDYVDR	C1(Carbamidomethyl)	1	4,51	2	1865,78093	-1,30	933,39410	24,23	0
	High	GYYSVETVTLVGLK		1	4,37	2	1641,90911	-0,45	821,45819	25,96	0
	High	SPVTIcGDIHGQFHDLAELFR	C6(Carbamidomethyl)	1	3,71	3	2412,17466	-0,80	804,72974	24,54	0
	High	cPDTNYLFmGDYVDR	C1(Carbamidomethyl); M9(Oxidation)	1	3,55	2	1881,77556	-1,45	941,39142	22,89	0
	High	AHQLVmDGFNWAHEQK	M6(Oxidation)	1	3,27	3	1926,89096	-0,34	642,96851	20,35	0
	High	cGNmASILEVDDcR	Carbamidomethyl); M4(Oxidation); C13(Carbamidomet	1	3,22	2	1655,68266	0,22	828,34497	20,96	0
	High	AKEILmDESNVQPVK	M6(Oxidation)	1	3,18	2	1716,88078	-1,73	858,94403	18,90	1
	High	NHTFIQFEPAPR		1	2,78	2	1456,73174	-1,03	728,86951	20,98	0
	High	EILmDESNVQPVK	M4(Oxidation)	1	2,56	2	1517,74986	-1,19	759,37857	19,86	0
	High	VVTIFSAFNYcYR	C11(Carbamidomethyl)	1	2,47	2	1589,77725	-0,54	795,39226	23,10	0
<b>AT5G26910</b>	<b>TRM8</b>	<b>38,62</b>	<b>21,83</b>	<b>11</b>	<b>11</b>	<b>852</b>					
	Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	ΔM [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
	High	VLSESEVSdCtSFYDK	C11(Carbamidomethyl)	1	4,33	2	1952,84294	-0,05	976,92511	21,00	0
	High	SDSSccASSVTSDDGQGTR	C5(Carbamidomethyl); C6(Carbamidomethyl)	1	4,28	2	1976,75078	-2,50	988,87903	17,47	0
	High	EYALGmATDVLPAFLFDEmEGR	M6(Oxidation); M19(Oxidation)	1	4,24	2	2447,10674	-1,60	1224,05701	25,32	0
	High	GLSSDSLSTQIGIGQDTPSAVSNIGDLSNALLEQK		1	4,21	3	3711,77817	-1,39	1237,93091	25,68	0
	High	LmGLESLPVNPVQePR	M2(Oxidation); E14(Methyl)	1	3,76	2	1808,95256	-2,78	904,97992	23,58	0
	High	WDAYENLGYVNLr		1	3,69	2	1612,77402	-0,91	806,89065	23,68	0
	High	EmDVISFTFSSPIK	M2(Oxidation)	1	3,30	2	1616,78630	-0,87	808,89679	24,57	0
	High	EGVVSALVEPHLLSAF		1	3,19	2	1667,89958	-0,48	834,45343	26,25	0
	High	FQSETFPPR		1	2,60	2	1108,54094	-1,16	554,77411	21,18	0
	High	NPIYVmEAASR	M6(Oxidation)	1	2,56	2	1266,61394	-0,65	633,81061	19,46	0
	High	ITTLTTPSTSK		1	2,44	2	1236,66716	-0,85	618,83722	19,02	0
<b>AT4G28760</b>	<b>TRM20</b>	<b>35,29</b>	<b>13,53</b>	<b>9</b>	<b>9</b>	<b>924</b>					
	Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	ΔM [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
	High	LLQELVEEAVIDLTR		1	4,99	2	1740,97331	-0,56	870,99030	29,93	0
	High	DSGEDmDANSLAAELSVK	M6(Oxidation)	1	4,96	2	1867,82109	-0,85	934,41418	21,79	0
	High	DASQcSsmSQLAAPSPVTLTGK	C5(Carbamidomethyl); M8(Oxidation)	1	3,87	2	2252,05107	-1,05	1126,52917	21,46	0
	High	INSIVSETITTR		1	3,85	2	1321,69499	-0,63	661,35114	19,64	0
	High	TSSTLGEMLALTETK	M8(Oxidation)	1	3,69	2	1597,79778	-0,77	799,40253	21,60	0
	High	AVSSSQSSPR		1	3,06	2	1005,49584	-0,16	503,25156	15,45	0
	High	VSNLFFFK		1	2,70	2	1001,54588	0,36	501,27658	24,50	0
	High	LASSSQETGWGNR		1	2,61	2	1392,64812	-1,56	696,82770	18,83	0
	High	GYFDEPEDVETK		1	2,52	2	1428,61565	-0,66	714,81146	21,22	0
<b>AT2G42500</b>	<b>PP2A-3</b>	<b>28,75</b>	<b>36,74</b>	<b>1</b>	<b>8</b>	<b>313</b>					
	Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	ΔM [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
	High	GAGYTFGQDISEQFNHTNNLK		1	4,78	2	2341,08	-0,80	1171,04	22,30	0
	High	SPVTIcGDIHGQFHDLAELFR	C6(Carbamidomethyl)	1	3,71	3	2412,17	-0,80	804,73	24,54	0
	High	cGNmASILEVDDcR	Carbamidomethyl); M4(Oxidation); C13(Carbamidomet	1	3,22	2	1655,68	0,22	828,34	20,96	0
	High	AKEILmDESNVQPVK	M6(Oxidation)	1	3,18	2	1716,88	-1,73	858,94	18,90	1
	High	IGGmcPDTNYLFmGDYVDR	M4(Oxidation); C5(Carbamidomethyl); M13(Oxidation)	1	2,92	2	2255,94	-0,86	1128,47	23,08	0
	High	NHTFIQFEPAPR		1	2,78	2	1456,73	-1,03	728,87	20,98	0
	High	EILmDESNVQPVK	M4(Oxidation)	1	2,56	2	1517,75	-1,19	759,38	19,86	0
	High	VVTIFSAFNYcYR	C11(Carbamidomethyl)	1	2,47	2	1589,78	-0,54	795,39	23,10	0
<b>AT3G05750</b>	<b>TRM6</b>	<b>25,87</b>	<b>13,23</b>	<b>7</b>	<b>7</b>	<b>801</b>					
	Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	ΔM [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
	High	SDSScSTSTPTSDDGQGSK	C5(Carbamidomethyl)	1	4,16	2	1903,74394	-1,00	952,37561	16,26	0
	High	LmGLESLPVNPALEPR	M2(Oxidation)	1	3,82	2	1751,93328	-1,63	876,47028	23,88	0
	High	QNQFAETSVSNQR		1	3,31	2	1508,70842	-0,29	754,85785	18,55	0
	High	dmDVISFTFSSPIK	D1(Methyl); M2(Oxidation)	2	3,30	2	1616,78630	-0,87	808,89679	24,57	0
	High	DmDVISFTFSSPIK	M2(Oxidation)	3	3,12	2	1602,77104	-0,64	801,88916	25,04	0
	High	SKKPANGVQEAAGVNSDK		1	2,98	3	1728,88565	-1,12	576,96674	15,81	1

	High	GQSPQISNGTcNNK	C11(Carbamidomethyl)	1	2,63	2	1504,68108	0,09	752,84418	16,73	0
	High	FSSSDSSSLPmK	M12(Oxidation)	1	2,55	2	1375,60356	-0,79	688,30542	18,19	0
<b>AT1G74160</b>	<b>TRM4</b>	<b>23,22</b>	<b>9,37</b>	<b>7</b>	<b>7</b>	<b>1025</b>					
	Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	ΔM [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
	High	SLQQVDDQLSQASNESR		1	4,73	2	1904,89250	-1,05	952,94989	20,71	0
	High	SGNWADFSGEmSGLVLDVER	M11(Oxidation)	1	3,99	2	2184,98296	-1,73	1092,99512	24,41	0
	High	YISEILLASGLLLR		1	3,44	2	1560,93437	-1,06	780,97083	27,60	0
	High	LSEFSFEScSSSK	C9(Carbamidomethyl)	1	3,29	2	1521,63555	-1,22	761,32141	18,69	0
	High	ELcSAIETQQK	C3(Carbamidomethyl)	1	2,85	2	1306,62993	-0,66	653,81860	19,09	0
	High	SPSVIEAAK		1	2,54	2	901,49932	0,38	451,25330	18,43	0
	High	QQQSTNFAVQR		1	2,38	2	1306,64934	-0,41	653,82831	18,56	0
<b>AT1G25490</b>	<b>PP2AA1</b>	<b>17,70</b>	<b>13,10</b>	<b>1</b>	<b>5</b>	<b>588</b>					
	Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	ΔM [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
	High	LLNPelaiQHILPcVK	C14(Carbamidomethyl)	1	3,78	2	1858,06035	-0,88	929,53381	24,89	0
	High	LAIIEYVPLLASQLGIGFFDdK	D21(Methyl)	4	2,94	3	2436,33982	-1,14	812,78479	31,43	0
	High	LLRDNEAEVR		1	2,93	2	1214,64739	-1,20	607,82733	17,65	1
	High	ELSSDSSQHVR		1	2,59	2	1244,58501	-1,29	622,79614	16,21	0
	High	YFANQALNSIDGSTAAQS		1	2,54	2	1857,86443	1,62	929,43585	23,14	0
<b>AT5G57580</b>	<b>Calmodulin-binding protein</b>	<b>11,50</b>	<b>6,96</b>	<b>3</b>	<b>3</b>	<b>647</b>					
	Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	ΔM [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
	High	EGVGTGLGELVFTDNSSWIR		1	4,41	2	2080,03520	0,26	1040,52124	26,17	0
	High	LTAEGINTVEDFLR		1	3,87	2	1577,81584	-0,76	789,41156	24,82	0
	High	NLDGNDDQPER		1	3,22	2	1387,57158	-0,38	694,28943	17,51	0
<b>AT1G63670</b>	<b>TRM12</b>	<b>5,77</b>	<b>3,92</b>	<b>2</b>	<b>2</b>	<b>689</b>					
	Peptide confidence	Sequence	Modifications	Rank	XCorr	Charge	MH+ [Da]	ΔM [ppm]	m/z [Da]	RT [min]	# Missed Cleavages
	High	IGLSFDDPNIDLIEK		1	2,91	2	1688,87175	-1,47	844,93951	24,77	0
	High	TSDVSTSTVGHR		1	2,85	2	1246,60198	-0,23	623,80463	16,47	0

Protein identification details were obtained with the LTQ Orbitrap Velos (Thermo Fisher Scientific) and the Proteome Discoverer software package version 1.3 (Thermo Fisher Scientific) combined with search engine SEQUEST.

Proteins and peptides headers generated by the software package and used in the table are listed below.

**Score:** Displays the protein score, which is the sum of the scores of the individual peptides. For SEQUEST results, the score is the sum of all peptide Xcorr values above the specified score threshold. The score threshold is calculated as follows:  $0.8 + \text{peptide\_charge} \times \text{peptide\_relevance\_factor}$  where peptide\_relevance\_factor is an advanced parameter of the SEQUEST node in the "Protein Scoring Option" category with a default value of 0.4. For each spectrum, only the highest-scoring match is used. For each spectrum and sequence, the Proteome Discoverer application uses only the highest scored peptide. When it performs a search using dynamic modifications, one spectrum might have multiple matches because of permutations of the modification site. **Coverage:** Displays by default the percentage of the protein sequence covered by identified peptides. **# Unique Peptides:** Displays the number of peptide sequences unique to a protein group. **# Peptides:** Displays the number of distinct peptide sequences in the protein group. **# AAs:** Shows by default the sequence length of the protein. **Peptide Confidence:** Indicates a confidence level associated with the peptide sequence: high confidence (>99%), medium confidence (>95%, <99%), low confidence (<95%). **Sequence:** Displays the sequence of amino acids that compose the peptide. **Modifications:** Displays the static and dynamic modifications identified in the peptide. **XCorr (search-dependent):** Scores the number of fragment ions that are common to two different peptides with the same precursor mass and calculates the cross-correlation score for all candidate peptides queried from the database (SEQUEST searches only). **Rank:** Displays the ordering of peptides by rank. **# Missed Cleavages:** Displays the number of cleavage sites in a peptide sequence that a cleavage reagent (enzyme) did not cleave. This number excludes cases where an amino acid such as proline inhibits the cleaving enzyme. For example, if proline resides next to lysine or arginine, trypsin does not cleave the lysine or arginine. The # of Missed Cleavages value is 0 for a complete digest of all peptides. **Charge:** Displays the charge state of the peptide. **m/z [Da]:** Displays the mass-to-charge ratio of the precursor ion, in daltons. **MH+ [Da]:** Displays the protonated monoisotopic mass of the peptides, in daltons.  **$\Delta M$  [ppm]:** Displays the difference between the theoretical mass of the peptide and the experimental mass of the precursor ion. **RT [min]:** Displays the retention time when the peptide was observed, in minutes.

### Supplementary Table S3: VELOS orbitrap analysis of TAP TON1a and FASS eluates

Bait	Accession	Description	Score	Coverage %	Unique Peptide #	Peptides #	AAs					
TON1A	AT3G55000	TON1a	13660	67.7	19	21	260					
TON1A	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages
TON1A	High	NGDSGPLLLDVLGFLK		1	139.44	5.60E-13	2	1785,951	1785,951	1E-04	893,983	0
TON1A	High	VIENNEGPPALLGSCNDR	E3(Methyl); C16(Carbamidomethyl)	1	128.48	8.30E-12	2	2064,988	2064,99	-0.002	1033,501	0
TON1A	High	FESMTQMGSSSR		1	115.98	2.00E-11	2	1403,579	1403,581	-0.002	702,797	0
TON1A	High	RDSETESSSLESR	E4(Methyl)	1	114.44	1.20E-10	2	1582,705	1582,707	-0.002	792,36	1
TON1A	High	ASAALENQLDR		1	103.4	2.30E-09	2	1299,678	1299,678	-2E-04	650,8463	0
TON1A	High	ASVFEAIEEEDR	E9(Methyl)	1	100.06	4.80E-09	2	1407,651	1407,652	-6E-04	704,8329	0
TON1A	High	ASAALENQLDRK		1	95.28	1.20E-08	2	1427,772	1427,773	-0.001	714,8933	1
TON1A	High	DSETESSSLESR		1	94.03	5.50E-09	2	1412,591	1412,59	3E-04	707,3026	0
TON1A	High	ASVFEAIEEEDRVIENNEGPPALLGSCNDR	C28(Carbamidomethyl)	1	92.36	2.00E-08	3	3426,594	3426,6	-0.006	1143,205	1
TON1A	High	SSASDSLPPQR		1	86.84	7.40E-08	2	1143,552	1143,552	-2E-04	572,7832	0
TON1A	High	DFNSNNGFELNR		1	78.36	3.50E-07	2	1425,622	1425,627	-0.005	713,8183	0
TON1A	High	AGLSTSGYR		1	76.86	8.20E-07	2	910,4507	910,4508	-2E-04	456,2326	0
TON1A	High	QGNQDTHIEVTR	E8(Methyl)	1	73.14	1.60E-06	2	1426,641	1426,644	-0.003	714,3278	0
TON1A	High	RPVSAQSADR		1	71.07	4.20E-06	2	1172,586	1172,59	-0.003	587,3005	1
TON1A	High	RSSASDSLPPQR		1	55.55	1.30E-04	2	1299,652	1299,653	-0.001	650,8332	1
TON1A	High	DFNSNNGFELNRNGDSGPLLLDVLGFLK		1	53.23	2.20E-04	3	3194,542	3193,568	0.974	1065,855	1
TON1A	High	KLHASPGR		1	51.22	2.00E-04	2	951,5245	951,525	-5E-04	476,7695	1
TON1A	High	RAGLSTSGYR		1	50.03	4.90E-04	2	1066,551	1066,552	-6E-04	534,283	1
TON1A	High	KDEFNWR		1	48.19	7.20E-04	2	993,4675	993,4668	7E-04	497,741	1
TON1A	High	DGTNEEGR		1	40.85	2.70E-04	2	1005,4	1005,4	4E-04	503,7074	0
TON1A	High	EMMDLK	M2(Oxidation)	1	34.9	5.30E-03	2	781,3349	781,335	-1E-04	391,6747	0
TON1A	Accession	Description	Score	Coverage %	Unique Peptide #	Peptides #	AAs					
TON1A	AT3G55005	TON1b	3596	62.3	11	13	257					
TON1A	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages
TON1A	High	SVSASQASGAATSGYR		1	137.32	7.30E-13	2	1498,695	1498,701	-0.006	750,3547	0
TON1A	High	PLDLLDVLGFLK		1	109.56	1.10E-10	2	1355,806	1355,806	-4E-04	678,9103	0
TON1A	High	FENMTQVMGSSSR		1	105.35	5.60E-10	2	1442,627	1442,628	-0.002	722,3206	0
TON1A	High	VIENNEGPPALLGSCNDR	C16(Carbamidomethyl)	1	105.33	1.80E-09	2	2067,005	2067,006	-2E-04	1034,51	0
TON1A	High	ASVFEAIEEEDR	M9(Methyl)	1	100.06	4.80E-09	2	1407,651	1407,652	-6E-04	704,8329	0
TON1A	High	SSASDSLPHQR		1	90.07	3.50E-08	2	1183,558	1183,558	-2E-04	592,7862	0
TON1A	High	NEDSRPLDLLDVLGFLK		1	88.36	4.30E-08	3	1957,052	1957,052	-5E-04	653,3578	1
TON1A	High	ASTALENQLDR		1	87.7	1.00E-07	2	1329,689	1329,689	1E-04	665,8517	0
TON1A	High	YDTEDEPEEVMR		1	74.45	1.70E-07	2	1513,805	1513,807	-0.002	757,8096	0
TON1A	High	DFNSNNGFELNR		1	70.46	5.20E-06	2	1440,65	1440,663	-0.014	721,3321	0
TON1A	High	LLSALICEVLDWAQLNHTLK	C7(Carbamidomethyl)	1	57.52	7.20E-05	3	2400,252	2400,251	4E-04	801,0911	0
TON1A	High	VYQPECNASAK	C8(Carbamidomethyl)	1	56.03	7.80E-05	2	1194,533	1194,534	-8E-04	598,2738	0
TON1A	High	EMMDLK	M2(Oxidation)	1	34.9	5.30E-03	2	781,3349	781,335	-1E-04	391,6747	0
TON1A	Accession	Description	Score	Coverage %	Unique Peptide #	Peptides #	AAs					
TON1A	AT5G18580	FASS	1323	51.5	20	20	480					
TON1A	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages
TON1A	High	YIGSGAGLGSEALMELETK		1	144.7	2.10E-13	2	1924,943	1924,945	-0.002	963,479	0
TON1A	High	QSDLLLNADDLAAMWVCLR	M14(Oxidation); C17(Carbamidomethyl)	1	131.18	4.50E-12	2	2219,071	2219,072	0	1110,543	0
TON1A	High	EMDFDSFLDFLVALENK	M2(Oxidation)	1	115.01	1.40E-10	2	2047,946	2047,945	0.002	1024,981	0
TON1A	High	ENLLQEEEPPEESQ		1	108.71	1.80E-10	2	1927,815	1927,817	-0.002	964,9149	0
TON1A	High	EYADGTLTEIFIR		1	105.95	1.30E-09	2	1655,806	1655,804	0.001	828,91	0
TON1A	High	GGGTVASMLDVR		1	103.75	2.90E-09	2	1345,702	1345,702	-9E-04	673,858	0
TON1A	High	SQEAGTIPSFYK		1	94.53	1.60E-08	2	1326,646	1326,646	-1E-04	664,33	0
TON1A	High	GFLTADHSLFR		1	88.06	1.00E-07	2	1476,772	1476,773	-5E-04	739,3933	0
TON1A	High	MNYEDFCHIASVCTEQIGPK	M1(Oxidation); C7(Carbamidomethyl); C13(Carbamidomethyl)	1	84.05	5.10E-08	3	2414,032	2414,034	-0.003	805,6845	0
TON1A	High	DTPEGLTYLFR		1	80.13	5.10E-07	2	1310,651	1310,651	1E-04	656,3326	0
TON1A	High	ITLDGDLGCK	C9(Carbamidomethyl)	1	73.21	3.20E-06	2	1088,59	1088,59	2E-04	545,3024	0
TON1A	High	ICDMFLALDK	C2(Carbamidomethyl)	1	71.4	3.80E-06	2	1224,589	1224,588	6E-04	613,3017	0
TON1A	High	ENCVIDDATGAEK	C3(Carbamidomethyl)	1	70.28	1.70E-06	2	1420,61	1420,614	-0.004	711,3122	0
TON1A	High	DMPPAFNQMYCR	M2(Oxidation); C11(Carbamidomethyl)	1	54.97	1.40E-05	2	1544,62	1544,621	-0.001	773,3171	0
TON1A	High	ILLEIFK		1	53.31	1.40E-05	2	874,5528	874,5528	1E-04	438,2837	0
TON1A	High	FFFFCDPHR	C5(Carbamidomethyl)	1	44.26	7.60E-04	2	1271,554	1271,555	-7E-04	636,7843	0
TON1A	High	CLDLQGR	C1(Carbamidomethyl)	1	43.46	2.50E-03	2	860,4177	860,4174	3E-04	431,2161	0
TON1A	High	VFDEHVR		1	41.56	2.00E-03	2	900,4455	900,4454	2E-04	451,23	0
TON1A	High	GFWAHDNR		1	39.82	1.50E-03	2	1001,446	1001,447	-5E-04	501,7304	0
TON1A	High	KPEEGSISQR		1	37.23	8.80E-03	2	1129,572	1129,573	-9E-04	565,7932	1
TON1A	Accession	Description	Score	Coverage %	Unique Peptide #	Peptides #	AAs					
TON1A	AT3G25800	PP2AA2	475	29	8	11	587					
TON1A	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages
TON1A	High	TDVMSMFEDLTQDDQDSVR		1	139.42	1.10E-13	2	2230,937	2230,936	0.001	1116,476	0
TON1A	High	ESLDVDFHSLVK		1	85.81	1.50E-07	2	1500,782	1500,782	-5E-04	751,3982	0
TON1A	High	SALASVIMGMAPVLGK		1	77.81	6.90E-07	2	1543,846	1543,847	-3E-04	772,9304	0
TON1A	High	FFANQALQSIDNMMSS	M15(Oxidation)	1	71.83	1.80E-06	2	1917,86	1917,86	-1E-04	959,9373	0
TON1A	High	LAAGEWFTAR		1	67.55	1.20E-05	2	1120,56	1120,567	-0.006	561,2874	0
TON1A	High	TIRPGLVLESDPDVDVR		1	58.51	6.40E-05	2	2009,042	2009,043	-0.001	1005,528	1
TON1A	High	ELSSDSSQHYR		1	56.62	6.90E-05	2	1243,579	1243,579	-3E-04	622,7967	0
TON1A	High	AVSLLAPVMGSEITCSK	C15(Carbamidomethyl)	1	56.53	1.50E-04	2	1761,901	1761,901	1E-04	881,9576	0
TON1A	High	VLQSLIPVDQSVVEK		1	56.33	2.00E-05	2	1766,018	1766,019	-0.001	884,0162	0
TON1A	High	LLEPQDCVQHILPVI/NFVSQDK	C7(Carbamidomethyl)	1	52.79	2.10E-04	3	2591,34	2591,342	-0.002	864,7874	0
TON1A	High	FAATVESHLK		1	52.37	3.50E-04	2	1172,617	1172,619	-0.002	587,316	0

Accession	Description	Score	Coverage	% Unique	Peptide #	Peptides	# AAs				
TON1A	AT5G57580	313	14,8	7	8	647					
TON1A	Peptide confidence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	Δ [Da]	m/z [Da]	# Missed Cleavages
TON1A	High	LTAEAGINTVEDFLR	1	105,31	2,20E-09	2	1576,809	1576,81	-3E-04	789,412	0
TON1A	High	EGVGTGLGVFTDNNSSWIR	1	91,18	4,80E-08	2	2079,028	2079,027	4E-04	1040,521	0
TON1A	High	LYIYYAEDSR	1	78,34	7,10E-07	2	1291,61	1291,608	0,002	646,8123	0
TON1A	High	MWDALVEHAK	1	68,57	5,60E-06	2	1198,58	1198,581	0	600,2975	0
TON1A	High	LCSSELPILR	1	59,75	6,50E-05	2	1186,038	1186,038	-3E-04	594,3261	0
TON1A	High	AVVYGPASAK	1	51,35	5,10E-04	2	1090,565	1090,566	-8E-04	546,2898	0
TON1A	High	LSLPLFTGGK	1	44,04	8,10E-04	2	1031,601	1031,602	-1E-04	516,808	0
TON1A	High	HYPPALNDVVWR	1	38,77	7,20E-03	2	1481,704	1481,705	-0,001	741,8592	0
TON1A	AT3G53540	307	11,8	9	9	924					
TON1A	Peptide confidence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	Δ [Da]	m/z [Da]	# Missed Cleavages
TON1A	High	SHSSYNSPEVSIPTLSLK	1	141,16	4,30E-13	2	2005,959	2005,959	2E-04	1003,987	0
TON1A	High	ELQWLSLEDDIEIGR	1	105,19	1,50E-09	2	1927,988	1927,989	-0,001	965,0011	0
TON1A	High	LLFDQISR	1	57,35	7,70E-05	2	990,5495	990,5498	-3E-04	496,282	0
TON1A	High	EGDQPSPVSVLEASFDVSSGSECFESVSADLR	1	54,53	5,40E-05	3	3617,568	3616,564	1,005	1206,863	0
TON1A	High	MQLQLLK	1	54,28	2,10E-04	2	872,5149	872,5153	-4E-04	437,2647	0
TON1A	High	KLFDQISR	1	47,98	3,90E-04	2	1118,644	1118,645	-3E-04	560,3295	1
TON1A	High	LPCTTHIGR	1	45,34	6,50E-04	2	1041,502	1041,503	-6E-04	521,7583	0
TON1A	High	IQETLR	1	39,35	7,20E-03	2	758,4288	758,4286	2E-04	380,2217	0
TON1A	High	FEHEISR	1	36,28	9,20E-03	2	1158,567	1158,567	-4E-04	580,2905	0
TON1A	AT2G42500/AT3G5850	203	31	8	8	313					
TON1A	Peptide confidence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	Δ [Da]	m/z [Da]	# Missed Cleavages
TON1A	High	QITQVYGFYDECLR	1	86,57	1,00E-07	2	1790,828	1790,83	-0,002	896,4212	0
TON1A	High	NHTFIQFEPAPR	1	64,52	1,80E-05	2	1455,726	1455,726	-4E-04	728,87	0
TON1A	High	VVTIFSAPNYCYR	1	63,31	2,60E-05	2	1588,77	1588,771	-0,001	795,3921	0
TON1A	High	EILMDESINVQPVK	1	47,99	9,40E-04	2	1516,744	1516,744	-8E-04	759,379	0
TON1A	High	YGNANVWK	1	45,9	1,40E-03	2	950,4602	950,461	-8E-04	476,2374	0
TON1A	High	RGEPDVTR	1	41,29	2,00E-03	2	928,4738	928,4726	0,001	465,2442	1
TON1A	High	GAGYTFGQDISEQFNHTNNLK	1	40,63	3,50E-03	3	2340,075	2340,077	-0,002	781,0324	0
TON1A	High	PLSEQQVR	1	40,4	2,60E-03	2	955,5088	955,5087	1E-04	478,7617	0
TON1A	AT3G05750	179	6,4	4	4	801					
TON1A	Peptide confidence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	Δ [Da]	m/z [Da]	# Missed Cleavages
TON1A	High	EFSLGMATDILPLSLFDETEGK	1	101,6	4,10E-09	2	2412,175	2412,177	-0,002	1207,095	0
TON1A	High	QNQFAETSVSNQR	1	83,53	2,10E-07	2	1507,7	1507,702	-0,002	754,8571	0
TON1A	High	QDIFLER	1	66,95	1,60E-05	2	919,4759	919,4763	-4E-04	460,7452	0
TON1A	High	EILADQVLK	1	36,03	6,30E-03	2	1027,591	1027,591	-5E-04	514,8027	0
TON1A	AT5G26910	86	4,6	2	2	853					
TON1A	Peptide confidence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	Δ [Da]	m/z [Da]	# Missed Cleavages
TON1A	High	ETYEEGIDIEGEIVSTLVDLVDLVSGTQR	1	78,91	5,00E-07	3	3407,643	3407,647	-0,003	1136,888	0
TON1A	High	TLDFVFNK	1	37,73	5,20E-03	2	982,5128	982,5124	4E-04	492,2637	0
TON1A	AT2G45240	50	5,3	2	2	398					
TON1A	Peptide confidence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	Δ [Da]	m/z [Da]	# Missed Cleavages
TON1A	High	HATMSGLSVVR	1	41,17	4,30E-03	2	1172,595	1172,597	-0,002	587,3049	0
TON1A	High	AIAIVKPGVR	1	35,45	4,00E-04	2	1022,66	1022,66	-2E-04	512,3372	1
Bait	AT5G18580	3906	81	32	32	480					
FASS	Peptide confidence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	Δ [Da]	m/z [Da]	# Missed Cleavages
FASS	High	YIGSGAGLGSSEALMELETK	1	157,28	1,20E-14	2	1924,945	1924,945	2E-04	963,48	0
FASS	High	QSDLLLNADDLAAMWVCLR	1	146,63	1,40E-13	2	2203,08	2203,077	0,004	1102,547	0
FASS	High	MNYEDFCHIASVCTEQIGPK	1	130,67	1,10E-12	2	2414,032	2414,034	-0,002	1208,023	0
FASS	High	EMDFDSFLDFVLALENK	1	123	2,60E-11	2	2031,951	2031,95	7E-04	1016,983	0
FASS	High	SQEAAGTIPSFYK	1	114,05	1,90E-10	2	1326,645	1326,646	-0,001	664,3296	0
FASS	High	QGGTVAKMLDVR	1	107,68	1,20E-09	2	1345,702	1345,702	-2E-04	673,8584	0
FASS	High	ICDMFLALDKDMSSGLCK	1	107,63	4,80E-10	2	2118,946	2118,946	0	1060,48	1
FASS	High	GFLTADHSLFR	1	100,06	6,00E-09	2	1476,772	1476,773	-2E-04	739,3934	0
FASS	High	EYADGTLTEIFER	1	99,64	5,70E-09	2	1655,804	1655,804	-8E-04	828,909	0
FASS	High	ENCVIDDATGAEK	1	89,89	2,50E-08	2	1434,629	1434,63	-0,001	718,3215	0
FASS	High	WIEGGNYELCIEDVR	1	87,5	7,90E-08	2	1865,861	1865,862	-8E-04	933,9378	0
FASS	High	DTPEGLTYLFR	1	84,2	2,00E-07	2	1324,669	1324,666	0,003	663,342	0
FASS	High	ITLGDLLGCK	1	79,08	8,30E-07	2	1088,59	1088,59	-2E-04	545,3022	0
FASS	High	WIEGGNYELCIEDVRDEIWDVMK	1	75,28	1,10E-06	3	2868,31	2868,31	4E-04	957,1107	1
FASS	High	.SNCLQELMELHQSEEEVTDTEQAENWFSLTSA	1	72,75	1,20E-06	3	4523,093	4523,074	0,019	1508,705	0
FASS	High	ICDMFLALDK	1	71,65	3,60E-06	2	1224,588	1224,588	-2E-04	613,3013	0
FASS	High	IPPASSMLWVR	1	69,56	5,80E-06	2	1271,67	1271,67	1E-04	636,8421	0
FASS	High	KPEEGSISQSR	1	67,63	8,40E-06	2	1129,573	1129,573	0	565,7936	1
FASS	High	IDMSELDESDGFLHSDDEMIESYIGGLIPNLAQLR	1	67,21	4,60E-06	3	3824,738	3824,74	-0,002	1275,92	0
FASS	High	IAILPFYLMVR	1	59,31	4,00E-05	2	1513,836	1513,837	-5E-04	757,9254	0
FASS	High	DEIWDVMKPSDPLK	1	55,9	1,40E-04	2	1671,818	1671,818	-2E-04	836,9161	1
FASS	High	FFFCDPHR	1	52,39	1,20E-04	2	1271,554	1271,555	-4E-04	636,7844	0
FASS	High	VFDEHVR	1	52,06	1,80E-04	2	900,4445	900,4454	-8E-04	451,2295	0
FASS	High	ILLEIFK	1	50,34	2,90E-05	2	874,5529	874,5528	2E-04	438,2837	0
FASS	High	DMPAPFNQMYCR	1	48,04	1,10E-04	2	1528,626	1528,626	-3E-04	765,3202	0
FASS	High	FFSPSNFMK	1	44,83	1,10E-03	2	1103,511	1103,511	1E-04	552,7628	0

FASS	High	GFWAHDNR		1	44.77	5.00E-04	2	1001.447	1001.447	1E-04	501.7307	0
FASS	High	TVLSLQAR		1	44.65	2.70E-03	2	874.4873	874.4872	0	438.2509	0
FASS	High	KIPPASSMLWVR	M8(Oxidation)	1	44.64	1.50E-03	2	1399.764	1399.765	-0.001	700.889	1
FASS	High	CLDLQGR	C1(Carbamidomethyl)	1	43.5	2.50E-03	2	860.4179	860.4174	4E-04	431.2162	0
FASS	High	DMSGSLCK	C7(Carbamidomethyl)	1	40.97	9.60E-04	2	896.3733	896.3732	1E-04	449.1939	0
FASS	High	FEKDEAGR		1	40.95	3.80E-03	2	950.4455	950.4457	-2E-04	476.2301	1
FASS	Accession	Description	Score	Coverage	%f	Unique	Peptide #	Peptides	#	AAs		
FASS	AT3G55000	TON1A	1540	60.8	15	16	260					
FASS	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages
FASS	High	NGDSGPLLLDLVLEGLFK		1	146.46	1.20E-13	2	1785.951	1785.951	-1E-04	893.9828	0
FASS	High	VIENNEGPPALLGSCNDR	C16(Carbamidomethyl)	1	112.98	2.90E-10	2	2050.971	2050.974	-0.003	1026.493	0
FASS	High	DSETSSSSLESR	E11(Methyl)	1	97.46	3.00E-09	2	1426.605	1426.606	-9E-04	714.3098	0
FASS	High	ASAALENLQDR		1	90.73	4.00E-08	2	1299.678	1299.678	1E-04	650.8464	0
FASS	High	FESMTQGMGSSSR	M4(Oxidation)	1	87.95	1.00E-08	2	1419.575	1419.576	-7E-04	710.7949	0
FASS	High	ASVFEAIEEEDR		1	86.91	9.00E-08	2	1393.636	1393.636	0	697.8253	0
FASS	High	RDSETSSSSLESR		1	75.52	7.30E-07	2	1568.691	1568.691	-2E-04	785.3529	1
FASS	High	DFNSNNGFELNRNGDSGPLLLDLVLEGLFK		1	74.46	1.70E-06	3	3194.544	3193.568	0.976	1065.855	1
FASS	High	RPVSAQASDR		1	73.76	2.00E-06	2	1172.589	1172.59	-4E-04	587.302	1
FASS	High	DFNSNNGFELNR		1	72.79	1.30E-06	2	1425.625	1425.627	-0.002	713.8198	0
FASS	High	SSASDSLPPDR		1	64.13	1.40E-05	2	1143.553	1143.552	0.001	572.7838	0
FASS	High	ASLSTSGYR		1	49.7	4.30E-04	2	910.4509	910.4508	1E-04	456.2327	0
FASS	High	ASVFEAIEEEDRVIENNEGPPALLGSCNDR	C28(Carbamidomethyl)	1	48.3	5.10E-04	3	3426.597	3426.6	-0.003	1143.206	1
FASS	High	QGNQDTHVEVTR	E8(Methyl)	1	46.2	8.30E-04	2	1426.644	1426.644	-2E-04	714.329	0
FASS	High	KDEFNWR		1	42.62	2.60E-03	2	993.4677	993.4668	9E-04	497.7411	1
FASS	High	KLHASPSGR		1	35.75	7.20E-03	2	951.5244	951.525	-6E-04	476.7895	1
FASS	Accession	Description	Score	Coverage	%f	Unique	Peptide #	Peptides	#	AAs		
FASS	AT3G55005	TON1b	1096	55.6	11	12	257					
FASS	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages
FASS	High	SVSASQASGAATSGYR		1	128.43	5.70E-12	2	1498.701	1498.701	-7E-04	750.3575	0
FASS	High	PLLLDLVLEGLFK		1	115.16	2.90E-11	2	1355.806	1355.806	-3E-04	678.9103	0
FASS	High	FENMTQVMGSSSR	M4(Oxidation)	1	104	6.60E-10	2	1458.623	1458.623	-4E-04	730.3187	0
FASS	High	VIENNEGLPPALLGSCNDR	C16(Carbamidomethyl)	1	94.86	2.00E-08	2	2067.005	2067.006	-8E-04	1034.51	0
FASS	High	ASVFEAIEEEDR		1	86.91	9.00E-08	2	1393.636	1393.636	0	697.8253	0
FASS	High	NEDSRPLLDVLEGLFK		1	84.31	1.20E-07	2	1957.052	1957.052	3E-04	979.5335	1
FASS	High	SSASDSLPHDR		1	79.87	4.50E-07	2	1193.559	1193.559	-5E-04	592.7861	0
FASS	High	ASTALENLQDR		1	74.85	2.00E-06	2	1329.688	1329.689	-8E-04	665.8513	0
FASS	High	YDTEDEPVEVMR	M6(Oxidation); E8(Methyl); M11(Oxidation)	1	69.58	3.20E-07	2	1559.611	1559.612	-6E-04	780.813	0
FASS	High	DFSINNGYELNR		1	67.6	7.00E-06	2	1440.66	1440.663	-0.004	721.3371	0
FASS	High	VYQPECNSAK	C6(Carbamidomethyl)	1	53.83	1.30E-04	2	1194.533	1194.534	-8E-04	598.2738	0
FASS	High	QLHASPSGR		1	44.3	1.80E-03	2	951.4883	951.4886	-3E-04	476.7514	0
FASS	Accession	Description	Score	Coverage	%f	Unique	Peptide #	Peptides	#	AAs		
FASS	AT3G25800	PP2AA2	1076	43.6	11	17	587					
FASS	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages
FASS	High	YMVANQLYELCEAVGPEPTR	M2(Oxidation); C11(Carbamidomethyl)	1	138.34	6.20E-13	2	2355.087	2355.088	-4E-04	1178.551	0
FASS	High	TDVMSMFEDLTQDDQDSVR	M4(Oxidation); M6(Oxidation)	1	132.03	2.90E-13	2	2262.925	2262.926	-8E-04	1132.47	0
FASS	High	FFANQALQSIDNVMMS	M14(Oxidation); M15(Oxidation)	1	109.61	2.40E-10	2	1933.855	1933.855	-3E-04	967.9346	0
FASS	High	ESLDVDFHLSLTK		1	94.85	1.90E-08	2	1500.781	1500.782	-0.002	751.3975	0
FASS	High	ILNPEIAQHILPCVK	C14(Carbamidomethyl)	1	89.37	1.40E-08	2	1857.054	1857.055	-9E-04	929.5341	0
FASS	High	AVSLLAPVMGSEITCSK	C15(Carbamidomethyl)	1	87.06	1.40E-07	2	1761.901	1761.901	3E-04	881.9576	0
FASS	High	LQQVNVQIGIDLLSQSLPAIVELAEADR		1	85.5	3.40E-08	3	3060.67	3060.671	-4E-04	1021.231	0
FASS	High	SLYTQLQDDMPMVR	C7(Carbamidomethyl); M11(Oxidation); M13(Oxidation)	1	77.52	3.40E-07	2	1887.815	1887.817	-0.001	944.9149	0
FASS	High	SALASVIMGMAPVLGK	M8(Oxidation)	1	76.94	9.30E-07	2	1559.842	1559.842	3E-04	780.9282	0
FASS	High	PGLVELSEDPDQVDR		1	76.52	1.40E-06	2	1638.81	1638.81	3E-04	820.4125	0
FASS	High	VLQSLPIVDQSVVEK		1	70.37	8.10E-07	2	1766.018	1766.019	-7E-04	884.0164	0
FASS	High	LAAGEWFTAR		1	70.01	6.30E-06	2	1120.567	1120.567	1E-04	561.2906	0
FASS	High	ELSSDSSQVVR		1	61.6	2.20E-05	2	1243.578	1243.579	-0.001	622.7964	0
FASS	High	LLEPQDCVQHILPVIVNFSQDK	C7(Carbamidomethyl)	1	55.14	1.30E-04	3	2591.34	2591.342	-0.002	864.7873	0
FASS	High	FAATVESAHLK		1	54.88	1.90E-04	2	1172.618	1172.619	-0.001	587.3162	0
FASS	High	NDDIQLR		1	52.64	1.80E-04	2	872.4355	872.4352	3E-04	437.225	0
FASS	High	TIRPGLVELSEDPDQVDR		1	51.26	3.30E-04	2	2009.042	2009.043	-7E-04	1005.528	1
FASS	Accession	Description	Score	Coverage	%f	Unique	Peptide #	Peptides	#	AAs		
FASS	AT3G53540	TRM19	876	25.8	17	17	924					
FASS	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages
FASS	High	SHSSYNSPEVSIPTLSLK		1	154.42	2.00E-14	2	2005.958	2005.959	-0.002	1003.986	0
FASS	High	TAMSSSEALDSTVTSVDPDISR	M3(Oxidation)	1	148.1	8.20E-14	2	2411.136	2411.137	-0.001	1206.575	0
FASS	High	SFSGDANSDEEDSSASDDIK		1	133.68	1.20E-13	2	2133.847	2133.846	0.001	1067.931	0
FASS	High	HASYDITDLPNEELR		1	112.37	2.70E-10	2	1771.837	1771.838	-6E-04	886.9258	0
FASS	High	PASFNGLSFEDGISK		1	106.28	1.30E-09	2	1567.748	1567.752	-0.004	784.8813	0
FASS	High	ELQWLSEDDIEIGR		1	104.2	1.90E-09	2	1927.989	1927.989	-4E-04	965.0016	0
FASS	High	EFNDALEALDSNK		1	97.02	8.10E-09	2	1464.673	1464.673	-3E-04	733.3437	0
FASS	High	TFASPSSSDSEFR		1	74.08	6.40E-07	2	1416.615	1416.616	-8E-04	709.3147	0
FASS	High	SGTLAEMLATSDR		1	64.42	2.10E-05	2	1350.645	1350.645	-1E-04	676.3297	0
FASS	High	LLFDQISR		1	63.55	1.80E-05	2	990.5501	990.5498	3E-04	496.2823	0
FASS	High	AMSFETSQFR		1	51.24	1.20E-04	2	1131.501	1131.502	-0.001	566.7576	0
FASS	High	MQLQLK		1	50.94	4.50E-04	2	872.5149	872.5153	-4E-04	437.2647	0
FASS	High	LPCTTTHGR		1	46.37	6.40E-04	2	1041.502	1041.503	-3E-04	521.7584	0
FASS	High	FEHEIEISR		1	41.34	3.30E-03	2	1158.567	1158.567	0	580.2907	0
FASS	High	FENNIQWPELPEPVGISSR		1	40.7	4.90E-03	2	2211.092	2211.096	-0.004	1106.553	0
FASS	High	TIMNQESAGGYTIVLPK	M3(Oxidation); E6(Methyl)	1	40.33	5.40E-03	2	1850.943	1850.945	-0.002	926.4789	0
FASS	High	FKDVFVLDLAK		1	37.39	7.90E-03	2	1309.685	1309.692	-0.007	655.8497	1
FASS	Accession	Description	Score	Coverage	%f	Unique	Peptide #	Peptides	#	AAs		
FASS	AT1G13320	PP2AA3	665	32.2	7	12	587					
FASS	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages

FASS	High	YMVANQLYELCEAVGPEPTR	M2(Oxidation); C11(Carbamidomethyl)	1	138,34	6,20E-13	2	2355,087	2355,088	-4E-04	1178,551	0
FASS	High	ILNPELAHQHILPCVK	C14(Carbamidomethyl)	1	89,37	1,40E-08	2	1857,054	1857,055	-9E-04	929,5341	0
FASS	High	AVSLLAPVMGSEITCSK	C15(Carbamidomethyl)	1	87,06	1,40E-07	2	1761,901	1761,901	3E-04	881,9576	0
FASS	High	LDQVNVIGIDLLSQSLPAIVELAEDR		1	85,5	3,40E-08	3	3060,67	3060,671	-4E-04	1021,231	0
FASS	High	SALASVIMGMAPVLGK	M8(Oxidation)	1	76,94	9,30E-07	2	1559,842	1559,842	3E-04	780,9282	0
FASS	High	LLEPQDCVAHILPVVNFSDQK	C7(Carbamidomethyl)	1	67,41	7,60E-06	3	2534,319	2534,32	-0,001	845,7803	0
FASS	High	LSAGEWFTAR		1	63,26	2,80E-05	2	1136,561	1136,561	-3E-04	569,2879	0
FASS	High	ELSSDSSSHYVR		1	61,5	2,20E-05	2	1243,578	1243,579	-0,001	622,7964	0
FASS	High	ESDLYVEHFTPLAK		1	58,29	1,10E-04	2	1484,751	1484,751	4E-04	743,3827	0
FASS	High	FAATIESAHLK		1	53,26	2,90E-04	2	1186,634	1186,635	-3E-04	594,3244	0
FASS	High	SIYGQLQDDMPMVR	C7(Carbamidomethyl); M11(Oxidation); M13(Oxidation)	1	48,29	2,30E-04	2	1843,788	1843,79	-0,002	922,9013	0
FASS	High	LLPAVITASK		1	29,37	5,30E-03	2	1011,633	1011,633	-3E-04	506,8235	0
FASS												
FASS	Accession	Description	Score	Coverage	%f	Unique Peptide	#	Peptides	#	AAs		
FASS	AT3G58500	PP2A4	477	40,9	2	10	313					
FASS	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages
FASS	High	GAGYTFQDDISEQFNHTNNLK		1	118,69	5,40E-11	2	2340,075	2340,077	-0,002	1171,045	0
FASS	High	EILMDESNVQPK	M4(Oxidation)	1	98,09	9,30E-09	2	1516,743	1516,744	-0,002	759,3786	0
FASS	High	CPDNTYLFMGDYVDR	C1(Carbamidomethyl); M9(Oxidation)	1	97,65	1,40E-09	2	1880,77	1880,771	-0,001	941,392	0
FASS	High	QITQVYGFYDECLR	C12(Carbamidomethyl)	1	89,96	4,60E-08	2	1790,83	1790,83	0	896,4222	0
FASS	High	VVTFISAPNYCYR	C11(Carbamidomethyl)	1	66,82	1,20E-05	2	1588,77	1588,771	-0,001	795,3921	0
FASS	High	YGNANWIK		1	66,68	1,10E-05	2	950,4605	950,461	-5E-04	476,2375	0
FASS	High	NHTFIQIEFPAPR		1	64,58	1,80E-05	2	1455,726	1455,726	-4E-04	728,37	0
FASS	High	PLSEQQVR		1	52,18	1,80E-04	2	955,5086	955,5087	-1E-04	478,7616	0
FASS	High	RGEPDVTR		1	50,65	2,60E-04	2	928,4722	928,4726	-5E-04	465,2434	1
FASS	High	AHQLVMDGFNWAHEQK	M6(Oxidation)	1	42,95	2,00E-03	3	1925,883	1925,884	-9E-04	642,9684	0
FASS												
FASS	Accession	Description	Score	Coverage	%f	Unique Peptide	#	Peptides	#	AAs		
FASS	AT5G26910	TRM8	366	22,3	13	13	853					
FASS	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages
FASS	High	ETVEEGIDIEGIEVSTLVDLNDLVSGTOR		1	105,63	1,20E-09	3	3408,656	3407,647	1,009	1137,226	0
FASS	High	SDSSCCASSVTSDDGGQTR	C5(Carbamidomethyl); C6(Carbamidomethyl)	1	83,18	7,70E-09	2	1975,749	1975,748	9E-04	988,8819	0
FASS	High	EYALGMATDVLPAFLFDEMEGR	M6(Oxidation); M19(Oxidation)	1	77,96	5,30E-07	2	2446,108	2446,103	0,004	1224,061	0
FASS	High	DNQPSMTSVLNQK	M6(Oxidation)	1	76,26	1,20E-06	2	1476,888	1476,888	4E-04	739,3514	0
FASS	High	EGVVSALVEPHLSAF		1	63,41	1,60E-05	2	1666,894	1666,893	8E-04	834,4542	0
FASS	High	QLGLSTASAEK		1	58,88	2,50E-04	2	1103,582	1103,582	-8E-04	552,798	0
FASS	High	DWLAEELNLR		1	47,62	7,20E-04	2	1144,551	1144,551	1E-04	573,283	0
FASS	High	VSLVEIDEIK		1	46,62	1,30E-03	2	1200,66	1200,66	-6E-04	601,3371	0
FASS	High	GGFLFEQR		1	41,95	2,90E-03	2	952,4771	952,4767	5E-04	477,2458	0
FASS	High	CEQFMFGSCR	Carbamidomethyl); M4(Oxidation); M6(Oxidation); C9(Carbamidomethyl)	1	41,54	7,00E-05	2	1336,466	1336,467	-4E-04	669,2405	0
FASS	High	ITSLTTPSTSK		1	40,77	3,90E-03	2	1235,663	1235,661	0,002	618,8387	0
FASS	High	VPVESGSISK		1	40,69	5,40E-03	2	1001,539	1001,539	-6E-04	501,7766	0
FASS	High	QNSNDFNLK		1	38,21	7,60E-03	2	1179,551	1179,552	-7E-04	590,7829	0
FASS												
FASS	Accession	Description	Score	Coverage	%f	Unique Peptide	#	Peptides	#	AAs		
FASS	AT2G45240	MAP1A	270	26,4	8	8	398					
FASS	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages
FASS	High	VEPNSDLQHVVEIK		1	83,72	2,70E-07	2	1605,834	1605,836	-0,002	803,9243	0
FASS	High	LPSSPDVYFPWLTK		1	79,16	7,90E-07	2	1501,781	1501,782	-4E-04	751,8979	0
FASS	High	WVPAIEKPKWIDGTPK		1	76,65	9,50E-07	2	1964,024	1964,026	-0,001	983,0194	1
FASS	High	SAQFEHTLVTGVEVLTAR		1	74,98	1,20E-06	3	2300,204	2300,201	0,003	767,7419	0
FASS	High	VIHPGVTTDEIDR		1	54,41	2,40E-04	2	1450,741	1450,742	-9E-04	726,3776	0
FASS	High	EVLDAAR		1	52,36	2,00E-04	2	843,4437	843,445	-0,001	422,7292	0
FASS	High	EIGEIVNR		1	44,47	1,20E-03	2	928,4976	928,4978	-1E-04	465,2561	0
FASS	High	AIIVKPGVR		1	36,16	3,40E-04	2	1022,66	1022,66	-3E-04	512,3372	1
FASS												
FASS	Accession	Description	Score	Coverage	%f	Unique Peptide	#	Peptides	#	AAs		
FASS	AT3G05750	TRM6	190	7,9	6	6	801					
FASS	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages
FASS	High	QNQFAETSVSNQR		1	64,48	1,80E-05	2	1507,702	1507,702	3E-04	754,8582	0
FASS	High	FSSSDSSSLPMK	M12(Oxidation)	1	58,48	2,40E-05	2	1374,596	1374,597	-9E-04	688,3055	0
FASS	High	QDIFLER		1	53,34	3,70E-04	2	919,4763	919,4763	0	460,7455	0
FASS	High	SPGFVQSR		1	46,03	2,00E-03	2	876,4449	876,4454	-4E-04	439,2298	0
FASS	High	CEQMFMGTCK	Carbamidomethyl); M4(Oxidation); M6(Oxidation); C9(Carbamidomethyl)	1	40,14	9,70E-05	2	1322,476	1322,476	-3E-04	662,2453	0
FASS												
FASS	Accession	Description	Score	Coverage	%f	Unique Peptide	#	Peptides	#	AAs		
FASS	AT4G28760	TRM20	186	6	4	4	924					
FASS	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages
FASS	High	LLQELVEEAVIDLTR		1	112,2	1,20E-10	2	1739,968	1739,967	0,001	870,9912	0
FASS	High	INSIVSETTTTR		1	90,69	5,40E-08	2	1320,687	1320,689	-0,001	661,351	0
FASS	High	VSNLFFFK		1	52,73	2,70E-04	2	1000,538	1000,538	-1E-04	501,2763	0
FASS	High	GTEEHTVQPTR		1	38,82	6,30E-03	2	1253,599	1253,6	-0,001	627,8068	0
FASS												
FASS	Accession	Description	Score	Coverage	%f	Unique Peptide	#	Peptides	#	AAs		
FASS	AT1G63670	TRM12	131	7,3	4	4	689					
FASS	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages
FASS	High	IGLSFDPNIDLEIK		1	72,17	3,80E-06	2	1697,865	1687,867	-0,002	844,9397	0
FASS	High	TSDVSTVYVYVHR		1	65,85	6,50E-06	2	1245,595	1245,595	-1E-04	623,8047	0
FASS	High	VLNLESSLR		1	63,7	2,30E-04	2	1029,584	1029,582	0,002	515,7993	0
FASS	High	QEQQPSPVSVLER		1	43,53	2,70E-03	2	1624,803	1624,806	-0,003	813,4088	0
FASS												
FASS	Accession	Description	Score	Coverage	%f	Unique Peptide	#	Peptides	#	AAs		
FASS	AT1G74160	TRM4	85	2,3	2	2	1025					
FASS	Peptide confidence	Sequence	Modifications	Rank	Peptide score	Expect	Charge	Mr(Exp)	Mr(Calc)	ΔM [Da]	m/z [Da]	# Missed Cleavages
FASS	High	SENFLEEEDDFLK		1	81,76	2,80E-07	2	1726,795	1726,794	0,002	864,4049	0
FASS	High	SILAEVDVTR		1	39,09	6,50E-03	2	1115,62	1115,619	0,002	558,8174	0

Protein identification details were obtained with the LTQ Orbitrap Velos (Thermo Fisher Scientific) and Mascot Distiller software (version 2.3.2.0, Matrix Science) combined with the Mascot search engine (Matrix Science) using the Mascot Daemon interface (version 2.4, Matrix Science).

Proteins and peptides headers used in the table are listed below.

**Score:** Displays the protein score. **Coverage %:** Displays by default the percentage of the protein sequence covered by identified peptides. **# Unique Peptides:** Displays the number of peptide sequences unique to a protein group. **# Peptides:** Displays the number of distinct peptide sequences in the protein group. **# AAs:** Shows by default the sequence length of the protein. **Peptide Confidence:** Indicates a confidence level associated with the peptide sequence: high confidence (>99%), medium confidence (>95%, <99%), low confidence (<95%). **Sequence:** Displays the sequence of amino acids that compose the peptide. **Modifications:** Displays the static and dynamic modifications identified in the peptide. **Rank:** Displays the ordering of peptides by rank. **Peptide score:** Displays the peptide ions score. **Expect:** Expectation value for the peptide match. (The number of times we would expect to obtain an equal or higher score, purely by chance. The lower this value, the more significant the result.) **Charge:** Displays the charge state of the peptide. **Mr(Exp):** experimental relative molecular mass, in daltons. **Mr(Calc):** calculated relative molecular mass, in daltons.  **$\Delta M$  [Da]:** difference (error) between the experimental and calculated masses, in daltons. **m/z [Da]:** Displays the mass-to-charge ratio of the precursor ion, in daltons. **# Missed Cleavages:** Displays the number of cleavage sites in a peptide sequence that a cleavage reagent (enzyme) did not cleave. This number excludes cases where an amino acid such as proline inhibits the cleaving enzyme. For example, if proline resides next to lysine or arginine, trypsin does not cleave the lysine or arginine. The # of Missed Cleavages value is 0 for a complete digest of all peptides.



**Supplementary Table S4: List of primers used in this study**

Primer name	Construct	Sequence
TON1a_fwd	Entry vectors of TON1a, TON1b, FASS and TRM19 to generate TAP constructs	published Van Leene et. al. 2007
TON1a_nostop_rev		published Van Leene et. al. 2007
TON1a_stop_rev		AGAAAGCTGGGTCTCAATCTTCCCTTCTTCTTC
TON1b_fwd		AAAAAGCAGGCTCCACCATTGACATTATACAAGAGAG
TON1b_stop_rev		AGAAAGCTGGGTCTCAATCTTCCCTTCTTCTTCTTC
TON1b_nostop_rev		AGAAAGCTGGGTCTTCTTCCCTTCTTCTTCTTCAC
FASS_fwd		AAAAAGCAGGCTCCACCATTGATAGCGGATCTAGCGATGG
FASS_stop_rev		AGAAAGCTGGGTCTCACTGAGACTTCTCTCAGG
FASS_nostop_rev		AGAAAGCTGGGTCTGAGACTTCTCTCAGGTGG
AT3G53540_fwd		AAAAAGCAGGCTCCACCATTGACAGATTTGCACTCAGC
AT3G53540_stop_rev		AGAAAGCTGGGTCTAGAAAGATTGCACCTACCACAAGC
AT3G53540_nostop_rev		AGAAAGCTGGGTGCGAAGATTGCACCTACCACAAGC
attB1-F		GGGGACAAGTTTGTACAAAAAAGCAGGCTCCACC
attB2-R		GGGGACCACTTTGTACAAGAAAGCTGGGTCT
FASSGW (+1)		Entry vectors of FASS and truncated versions of FASS
FASSGW (-480)	GGGGACCACTTTGTACAAGAAAGCTGGGTACTGAGACTTCTCTCAG	
FASSGW (-185)	GGGGACCACTTTGTACAAGAAAGCTGGGTCCGTGCGCATCACATAAG	
FASSGW (+186)	GGGGACAAGTTTGTACAAAAAAGCAGGCTACAAAATGGTGTCCCTTACACAAGC	
TRM1 (+1)	Entry vectors of TRM1 and truncated versions of TRM1	GGGGACAAGTTTGTACAAAAAAGCAGGCTACAAAATGTCTGCAAAACTTCT
TRM1 (+587)		GGGGACAAGTTTGTACAAAAAAGCAGGCTACAAAATGAGCTTGGGATCGAATGTT
TRM1 (-698)		GGGGACCACTTTGTACAAGAAAGCTGGGTATTATCCCGGTTACTCTGAGG
TRM1 (-827)		GGGGACCACTTTGTACAAGAAAGCTGGGTATTTTAAAGTGGATTTGC
TRM1 (-905)		GGGGACCACTTTGTACAAGAAAGCTGGGTAGCAGAAAGCAAACTTCAT
TON1a (+1)	Entry vectors of TON1a and truncated versions of TON1a	GGGGACAAGTTTGTACAAAAAAGCAGGCTTGTAGGACATTATACAAG
TON1a (-107)		GGGGACCACTTTGTACAAGAAAGCTGGGTACGAAATCTTTTGGCAA
TON1a (+108)		GGGGACAAGTTTGTACAAAAAAGCAGGCTACAAAATGTGGAAGTCTGAGTTACG
TON1a (-260)		GGGGACCACTTTGTACAAGAAAGCTGGGTAACTTCTTCCCTTCTTC
RCN1GW FWD		PP2AA1 entry vector
RCN1GW REV	PP2AA1 entry vector	GGGGACCACTTTGTACAAGAAAGCTGGGTAGGATTTGCTGCTGTGG
TRM8 (+1)	TRM8 entry vector	GGGGACAAGTTTGTACAAAAAAGCAGGCTACAAAATGGAGTTGTTGAGAGG
TRM8 (-854)	TRM8 entry vector	GGGGACCACTTTGTACAAGAAAGCTGGGTAGAATGCCGAGAGAAGAT
R1	Genotyping of <i>pp2aa1</i> mutant	AGCACATCCTTCTTGTGTGAAGG
R2		AACTTGCTTATGATGTTAAGGCGC
T (T-DNA)	Genotyping of <i>pp2aa2</i> mutant	TGTCGCCGCTCATCGGGGGGGT
oA2RAV		GCCTGCGGTGTCTCTTCTTGACC
oA2new3'	Genotyping of <i>pp2aa2</i> mutant	CACATTAGTAGCAAGACAATGGACAAAAACCCG
oA3leader	Genotyping of <i>pp2aa3</i> mutant	GATCGCTCGGAACCTTGGAAAGCAGC
oA3DDEV		GCCAAAAGCACCTCATCGTCATCGTC
SALK_069250 RP	Genotyping of <i>pp2ac3</i> mutant	TGTTTCTGATCTGTTTCCG
SALK_069250 LP		TAATTGGTATCAGGGCACTGC
SALK_035009 RP	Genotyping of <i>pp2ac4</i> mutant	GTGGATTATCACCATCCATCG
SALK_035009 LP		GCTTGAAGAACAGCATTTCG
LB-Salk2	Genotyping of <i>pp2aa2</i> , <i>pp2aa3</i> , <i>pp2ac3</i> and <i>pp2ac4</i> mutants	GCTTCTTCCCTTCTTCTCTC
ACT2_2_FWD	RT-PCR primers for actin, eif4A, PP2Ac1-c5 and PP2AA1-A3	TTGACTACGAGCAGGAGATGG
ACT2_2_REV		ACAACAGAGGGCTGGAAACAAG
eif4A-FWD		ACGGAGACATGGACCAAGAAC
eif4A-REV		GCTGAGTTGGGAGATCGAAG
at3g58500_LP1		TCTGCGAAAGTATGGCAATG
at3g58500_RP1		CCACCGTGAAGGCAGAATA
at2g42500_LP2		GAACCACACCTTCATCCAGTTT
at2g42500_RP2		TGAGATGAGTCTGTGCCAGTGA
at3g25800_fwd		CGATGCTGTGCTAACAATC
at3g25800_rev		ATTGCCATTGAGGACCA
at1g10430_fwd		CCAGAGAAGCTAGGAGGATTTGATC
at1g10430_rev		TTGAAAAGCCACCACCGATCC
at1g59830_fwd		TATGAGAGAAAACAAAACCAAGACG
at1g59830_rev		TCTCGTCTGCTCTCTTTAAC
at1g69960_fwd		ATGGACCAGAAATTCCTTCAGTTTG
at1g69960_rev		TAGTGTGTGGAATTGCAGCTTTAG
at1g13320_fwd		GACCGGAGCCAACCTAGGAC
at1g13320_rev		AAAACCTTGGTAACTTTTCCAGCA
at1g25490-fwd		ATCCTTGTACCAGGCAAT
at1g25490-rev		TGCAAGTCGGACTCTCCAG

### **Supplementary Methods for the LC-MS/MS VELOS1 analysis (related to Supplementary Table S2)**

Tuning of equipment and sample analyses were performed using a standard operation procedure. Quality performance of the LC MS analyses was assessed using control standards. Nanoflow LC-MS/MS analyses were performed on a dual channel NanoLC Ultra 2D system (Eksigent, Dublin, California, USA) and connected to an LTQ Orbitrap Velos mass spectrometer (Thermo, Bremen, Germany) using a chip based nano electrospray source (Triversa Advion Biosystems, Ithaca, NY, USA) operated at 1.8 kV. Mobile phase A (CH<sub>3</sub>CN : CHOOH : H<sub>2</sub>O in a 2:0.1:97.9 (v/v) ratio) and mobile phase B (CH<sub>3</sub>CN : CHOOH : H<sub>2</sub>O in a 90:0.1:9.9 (v/v) ratio) were reconstituted from ULC/MS-grade CH<sub>3</sub>CN, CHOOH and H<sub>2</sub>O (Biosolve-chemicals, Valkenswaard, the Netherlands). All standards and samples were dissolved in mobile phase A. A volume of 10 µL of each sample or control standard was loaded on a C<sub>18</sub> precolumn (PepMap 100, 5 µm particles, 20 mm x 200 µm ID, Dionex, Sunnyvale, California, USA) using a flow rate of 6 µL/min (100% mobile phase A). Five minutes after injection the trapping column was switched in line with an analytical C<sub>18</sub> column (Acclaim Pepmap 100, 3 µm particles, 150 mm x 75 µm ID, Dionex). Effective elution of peptides was accomplished within a 45 minutes gradient flow (95% A to 50% A) at a flow rate of 350 nl/min.

MS spectra were recorded in the orbitrap FT analyzer with a resolution of 60000 (at m/z 400) and an AGC target setting of 500000. The maximum injection time was set to 500 msec and lock mass was enabled (polysiloxane ion at m/z 445.12024). Collision induced dissociation MS/MS spectra were acquired using the ion trap in data dependent mode selecting up to the 20 most abundant multiply charged precursor ions from the MS spectrum. The maximum injection time was set to 50 msec and an AGC setting of 7500. Fragmentation was accomplished by CID wideband activation at normalized collision energy of 35 eV and a 30msec activation time. After MS/MS the precursor m/z's were excluded for 30 seconds. Peak lists were generated and submitted for protein identification with Proteome Discoverer 1.3.0.339 (Thermo, Bremen, Germany). Spectrum grouping was allowed with a maximum retention time of 1 minute and a precursor mass tolerance of 2 ppm. Peak lists were generated only for MS/MS spectra containing more than 5 peaks, with a total intensity above 50. The relative signal-to-noise limit was set to 5. Peak lists were submitted for protein identification against TAIR10 containing 35386 sequence entries with search engine SEQUEST. Enzyme was set to trypsin, allowing for maximum 1 missed cleavage site. Precursor mass tolerance was set to 3 ppm and fragment mass tolerance at 0.8 Da. Fixed modifications were set to carbamidomethylation of cysteines. Variable modifications were set to methionine oxidation and methylation of aspartic acid and glutamic acid, with a maximum of 4 modifications per peptide. Peptides were validated using a decoy database search. The strict target False Discovery Rate (FDR) was set to 0.01 (or 1%), the relaxed FDR was set to 0.05 (or 5%). Only high confident (>99%) peptides were withheld. Only proteins with at least two matched high confident peptides were retained.

### **Supplementary Methods for the LC-MS/MS VELOS2 analysis (related to Supplementary Table S3)**

The obtained peptide mixtures were introduced into an LC-MS/MS system, the Ultimate 3000 RSLC nano (Dionex, Amsterdam, The Netherlands) in-line connected to an LTQ Orbitrap Velos (Thermo Fisher Scientific, Bremen, Germany). The sample mixture was loaded on a trapping column (made in-house, 100 µm internal diameter (I.D.) x 20 mm (length), 5 µm

C18 Reprosil-HD beads, Dr. Maisch GmbH, Ammerbuch-Entringen, Germany). After back-flushing from the trapping column, the sample was loaded on a reverse-phase column (made in-house, 75  $\mu$ m I.D. x 150 mm, 5  $\mu$ m C18 Reprosil-HD beads, Dr. Maisch). Peptides were loaded with solvent A (0.1% trifluoroacetic acid, 2% acetonitrile), and separated with a linear gradient from 2% solvent A' (0.1% formic acid) to 50% solvent B' (0.1% formic acid and 80% acetonitrile) at a flow rate of 300 nl/min, followed by a wash step reaching 100% solvent B'.

The mass spectrometer was operated in data-dependent mode, automatically switching between MS and MS/MS acquisition for the ten most abundant peaks in a given MS spectrum. In the LTQ-Orbitrap Velos, full scan MS spectra were acquired in the Orbitrap at a target value of 1E6 with a resolution of 60,000. The ten most intense ions were then isolated for fragmentation in the linear ion trap, with a dynamic exclusion of 20 seconds. Peptides were fragmented after filling the ion trap at a target value of 1E4 ion counts.

From the MS/MS data in each LC run, Mascot Generic Files were created using the Mascot Distiller software (version 2.3.2.0, Matrix Science, [www.matrixscience.com/Distiller.html](http://www.matrixscience.com/Distiller.html)). When generating these peak lists, grouping of spectra was allowed with a maximum intermediate retention time of 30 seconds and a maximum intermediate scan count of 5 was used where possible. Grouping was done with 0.005 Da precursor tolerance. A peak list was only generated when the MS/MS spectrum contained more than 10 peaks. There was no de-isotoping and the relative signal-to-noise limit was set to 2. These peak lists were then searched with the Mascot search engine (MatrixScience, [www.matrixscience.com](http://www.matrixscience.com), PMID 10612281) using the Mascot Daemon interface (version 2.4, Matrix Science). Spectra were searched against the TAIR10 database containing 35386 sequence entries. Variable modifications were set to methionine oxidation and methylation of aspartic acid and glutamic acid. Fixed modifications were set to carbamidomethylation of cysteines. Mass tolerance on MS was set to 10 ppm (with Mascot's C13 option set to 1) and the MS/MS tolerance at 0.5 Da. The peptide charge was set to 1+, 2+ and 3+ and the instrument setting was set to ESI-TRAP. Trypsin was set as the protease used, allowing for 1 missed cleavage, and also cleavage was allowed when arginine or lysine is followed by proline. Only high confident peptides, ranked one and with scores above the threshold score, set at 99% confidence, were withheld. Only proteins with at least two matched high confident peptides were retained.