



Full wwPDB/EMDatabank EM Map/Model Validation Report ⓘ

Feb 8, 2018 – 12:28 PM EST

PDB ID : 6BF7
EMDB ID: : EMD-7091
Title : Cryo-EM structure of human insulin degrading enzyme in complex with FAB H11-E heavy chain, FAB H11-E light chain
Authors : Liang, W.G.; Zhang, Z.; Bailey, L.J.; Kossiakoff, A.A.; Tan, Y.Z.; Wei, H.; Carragher, B.; Potter, S.C.; Tang, W.J.
Deposited on : 2017-10-26
Resolution : 6.50 Å(reported)

This is a Full wwPDB/EMDatabank EM Map/Model Validation Report
for a publicly released PDB/EMDB entry.

We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

MolProbity : 4.02b-467
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030736

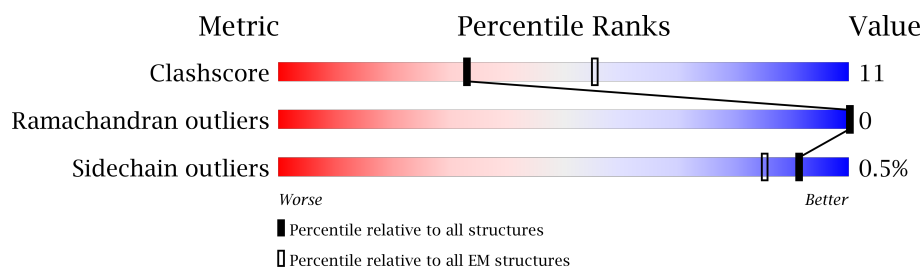
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 6.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	125131	1336
Ramachandran outliers	121729	1120
Sidechain outliers	121581	1026

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	966	69% 28% .
1	B	966	71% 26% .
2	C	218	81% 18%
2	E	218	72% 27%
3	D	211	76% 24%
3	F	211	76% 24%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 21866 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Insulin-degrading enzyme.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	949	Total	C	N	O	S	0	0
			7730	4980	1299	1429	22		
1	B	940	Total	C	N	O	S	0	0
			7662	4939	1285	1417	21		

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	110	LEU	CYS	engineered mutation	UNP P14735
A	171	SER	CYS	engineered mutation	UNP P14735
A	178	ALA	CYS	engineered mutation	UNP P14735
A	257	VAL	CYS	engineered mutation	UNP P14735
A	414	LEU	CYS	engineered mutation	UNP P14735
A	573	ASN	CYS	engineered mutation	UNP P14735
A	590	SER	CYS	engineered mutation	UNP P14735
A	789	SER	CYS	engineered mutation	UNP P14735
A	812	ALA	CYS	engineered mutation	UNP P14735
A	819	ALA	CYS	engineered mutation	UNP P14735
A	904	SER	CYS	engineered mutation	UNP P14735
B	110	LEU	CYS	engineered mutation	UNP P14735
B	171	SER	CYS	engineered mutation	UNP P14735
B	178	ALA	CYS	engineered mutation	UNP P14735
B	257	VAL	CYS	engineered mutation	UNP P14735
B	414	LEU	CYS	engineered mutation	UNP P14735
B	573	ASN	CYS	engineered mutation	UNP P14735
B	590	SER	CYS	engineered mutation	UNP P14735
B	789	SER	CYS	engineered mutation	UNP P14735
B	812	ALA	CYS	engineered mutation	UNP P14735
B	819	ALA	CYS	engineered mutation	UNP P14735
B	904	SER	CYS	engineered mutation	UNP P14735

- Molecule 2 is a protein called Fab H11-E heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	217	Total	C	N	O	S	0	0
			1614	1020	270	319	5		
2	E	217	Total	C	N	O	S	0	0
			1614	1020	270	319	5		

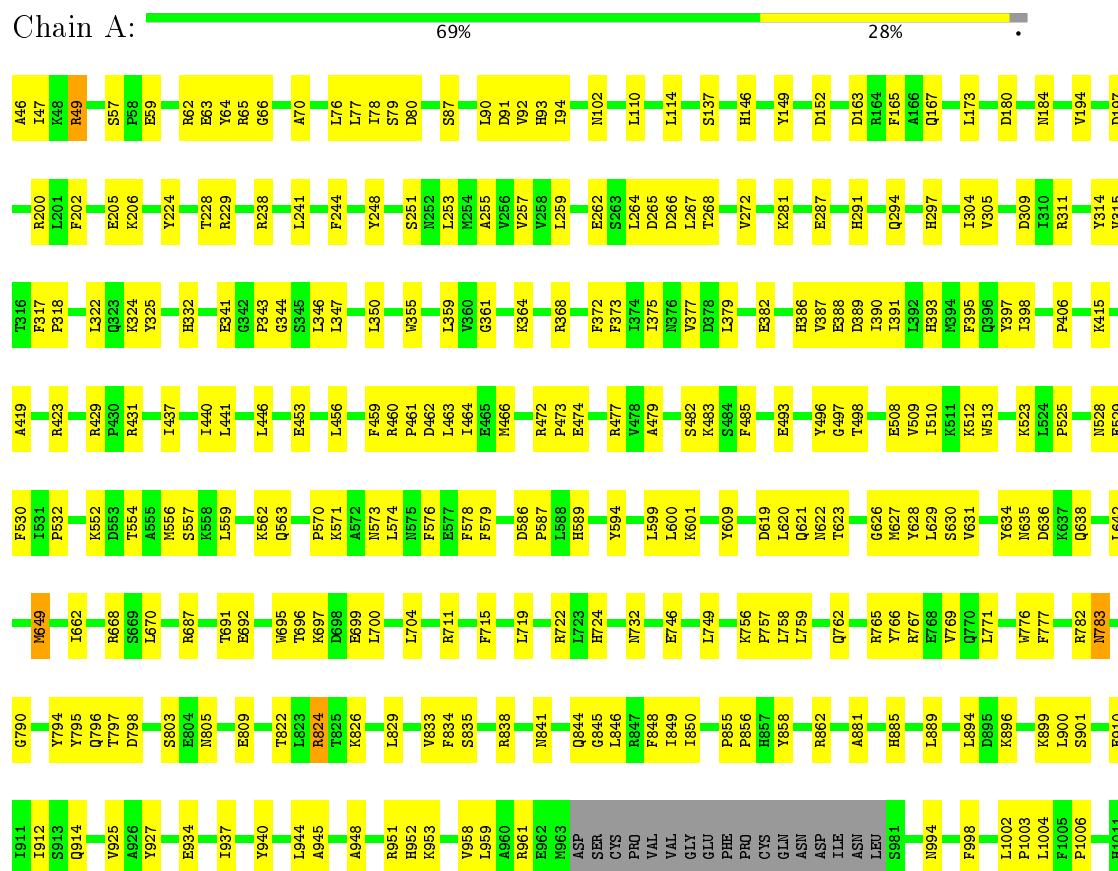
- Molecule 3 is a protein called Fab H11-E light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	211	Total	C	N	O	S	0	0
			1623	1016	272	330	5		
3	F	211	Total	C	N	O	S	0	0
			1623	1016	272	330	5		

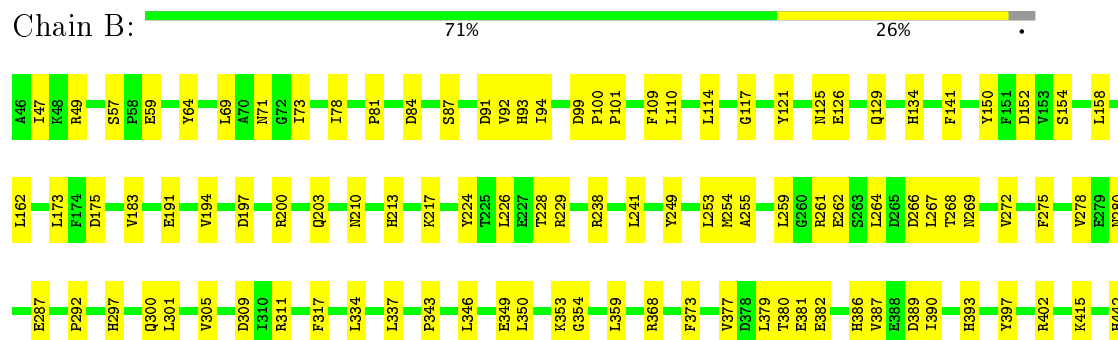
3 Residue-property plots

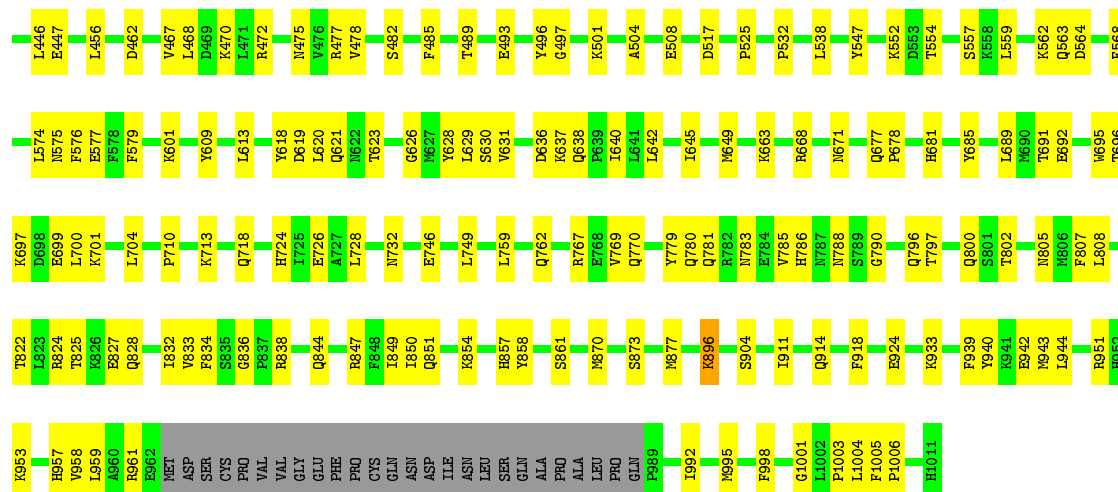
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Insulin-degrading enzyme



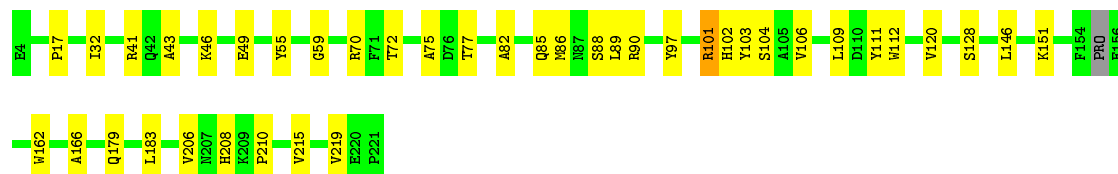
• Molecule 1: Insulin-degrading enzyme





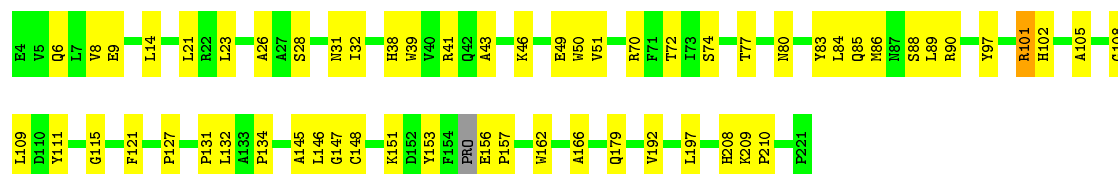
• Molecule 2: Fab H11-E heavy chain

Chain C: 81% 18%



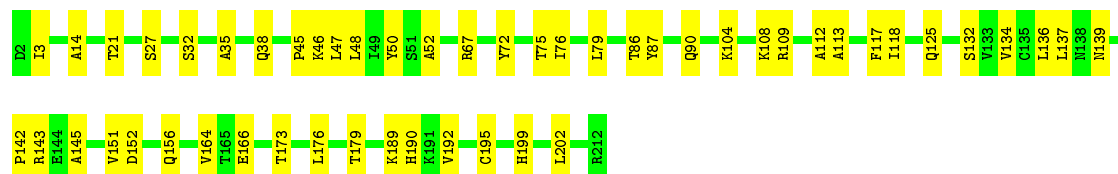
• Molecule 2: Fab H11-E heavy chain

Chain E: 72% 27%



• Molecule 3: Fab H11-E light chain

Chain D: 76% 24%



• Molecule 3: Fab H11-E light chain

Chain F: 76% 24%



Y147	K150	V151	D152	E162	S163	V164	T173	L176	D186	K191	V192	Y193	A194	H199	L202	R212
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4 Experimental information

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of particles used	16944	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	7.9, 6.8	Depositor
Minimum defocus (nm)	940	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	46598	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k), GATAN K2 SUMMIT (4k x 4k)	Depositor

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 2$	RMSZ	$\# Z > 2$
1	A	0.47	0/7924	0.74	0/10721
1	B	0.45	0/7854	0.73	0/10624
2	C	0.39	0/1653	0.66	0/2254
2	E	0.39	0/1653	0.64	0/2254
3	D	0.37	0/1658	0.64	0/2252
3	F	0.33	0/1658	0.61	0/2252
All	All	0.43	0/22400	0.71	0/30357

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	7730	0	7639	186	0
1	B	7662	0	7568	157	0
2	C	1614	0	1566	25	0
2	E	1614	0	1566	40	0
3	D	1623	0	1580	30	0
3	F	1623	0	1580	32	0
All	All	21866	0	21499	456	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (456) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:41:ARG:HB3	2:C:49:GLU:HB3	1.68	0.75
1:A:711:ARG:HH22	1:A:715:PHE:HB2	1.52	0.75
1:B:621:GLN:HB3	1:B:628:TYR:HB3	1.68	0.75
1:A:62:ARG:HG2	1:A:80:ASP:HB2	1.72	0.70
1:A:759:LEU:H	1:A:762:GLN:HE21	1.39	0.70
1:B:834:PHE:HB3	1:B:849:ILE:HB	1.71	0.70
1:A:317:PHE:HB2	1:A:373:PHE:HB3	1.72	0.70
1:A:314:TYR:HB2	1:A:479:ALA:HB3	1.73	0.69
2:E:41:ARG:HB3	2:E:49:GLU:HB3	1.74	0.69
1:A:346:LEU:HD11	1:A:393:HIS:HB3	1.75	0.69
1:B:759:LEU:H	1:B:762:GLN:HE21	1.41	0.68
1:A:309:ASP:O	1:A:668:ARG:NH1	2.26	0.68
3:F:117:PHE:HB2	3:F:136:LEU:HB3	1.76	0.68
1:B:317:PHE:HB2	1:B:373:PHE:HB3	1.74	0.67
1:A:576:PHE:HB2	1:A:629:LEU:HB3	1.76	0.67
2:C:32:ILE:HD11	2:C:77:THR:HA	1.76	0.67
1:A:389:ASP:O	1:A:393:HIS:ND1	2.27	0.66
1:A:881:ALA:O	1:A:885:HIS:HB2	1.95	0.66
1:A:224:TYR:O	1:A:228:THR:HB	1.96	0.65
1:A:782:ARG:HA	1:A:959:LEU:HB2	1.77	0.65
1:A:508:GLU:OE1	2:C:102:HIS:ND1	2.30	0.65
1:A:635:ASN:ND2	1:A:732:ASN:O	2.30	0.65
1:B:134:HIS:HB3	1:B:154:SER:HB2	1.78	0.65
2:E:156:GLU:HB2	2:E:210:PRO:HD3	1.78	0.65
2:E:70:ARG:HH22	2:E:90:ARG:HG2	1.63	0.64
1:B:575:ASN:HB2	1:B:728:LEU:HB3	1.79	0.64
2:E:102:HIS:HA	2:E:109:LEU:HA	1.79	0.64
2:E:131:PRO:HA	2:E:148:CYS:HA	1.80	0.63
1:B:287:GLU:HB3	1:B:368:ARG:HD3	1.79	0.63
1:B:538:LEU:H	1:B:732:ASN:HD21	1.46	0.63
1:B:769:VAL:O	1:B:796:GLN:NE2	2.32	0.63
1:B:346:LEU:HD11	1:B:393:HIS:HB3	1.80	0.63
2:E:8:VAL:HB	2:E:26:ALA:HB3	1.82	0.62
1:B:579:PHE:HB3	1:B:724:HIS:HB3	1.80	0.62
1:B:415:LYS:NZ	1:B:456:LEU:O	2.32	0.62
1:B:359:LEU:HB2	1:B:377:VAL:HG12	1.82	0.61
1:B:389:ASP:O	1:B:393:HIS:ND1	2.33	0.61
3:D:152:ASP:HA	3:D:192:VAL:HB	1.82	0.61
1:A:805:ASN:HA	1:A:844:GLN:HE22	1.66	0.61
1:B:759:LEU:HB2	1:B:762:GLN:HG2	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:601:LYS:HD2	1:A:620:LEU:HB2	1.82	0.60
2:C:146:LEU:HB2	2:C:219:VAL:HG11	1.84	0.60
1:A:359:LEU:HB2	1:A:377:VAL:HG22	1.83	0.60
1:A:579:PHE:N	1:A:724:HIS:O	2.34	0.59
1:B:779:TYR:HB3	1:B:992:ILE:HB	1.83	0.59
1:A:769:VAL:O	1:A:796:GLN:NE2	2.33	0.59
1:A:287:GLU:HB3	1:A:368:ARG:HD3	1.85	0.59
3:D:137:LEU:HB2	3:D:176:LEU:HB3	1.83	0.59
1:A:341:GLU:HG2	1:A:347:LEU:HD12	1.85	0.59
3:F:150:LYS:HB3	3:F:194:ALA:HB3	1.84	0.59
1:A:1004:LEU:HD12	1:B:1004:LEU:HD12	1.84	0.59
1:A:759:LEU:HB2	1:A:762:GLN:HG2	1.84	0.59
1:A:766:TYR:O	1:A:841:ASN:ND2	2.35	0.59
1:B:337:LEU:O	1:B:397:TYR:OH	2.19	0.59
1:B:264:LEU:O	1:B:268:THR:OG1	2.15	0.58
1:B:574:LEU:HB2	1:B:631:VAL:HB	1.85	0.58
1:B:508:GLU:OE1	2:E:102:HIS:ND1	2.35	0.58
1:B:93:HIS:HB3	1:B:253:LEU:HB3	1.84	0.58
1:A:570:PRO:HA	1:A:732:ASN:HD22	1.68	0.58
2:E:74:SER:HB2	2:E:83:TYR:HB2	1.85	0.58
1:B:305:VAL:HG22	1:B:485:PHE:HB2	1.86	0.58
1:B:73:ILE:HG12	1:B:254:MET:HB2	1.86	0.57
1:A:1002:LEU:O	1:B:767:ARG:NH1	2.36	0.57
1:A:322:LEU:HA	1:A:325:TYR:HD2	1.70	0.57
1:A:795:TYR:HB2	1:A:846:LEU:HB3	1.84	0.57
1:A:163:ASP:O	1:A:167:GLN:NE2	2.36	0.57
1:B:309:ASP:O	1:B:668:ARG:NH1	2.37	0.57
1:B:834:PHE:N	1:B:849:ILE:O	2.37	0.57
3:D:117:PHE:HB2	3:D:136:LEU:HB3	1.86	0.57
2:E:32:ILE:HD11	2:E:77:THR:HA	1.87	0.57
1:A:173:LEU:O	1:A:238:ARG:NH2	2.38	0.57
2:C:43:ALA:HB3	2:C:46:LYS:HB2	1.86	0.57
2:C:72:THR:HB	2:C:85:GLN:HB3	1.87	0.57
1:A:1003:PRO:HG3	1:B:1006:PRO:HG3	1.86	0.57
1:A:332:HIS:HE1	1:A:453:GLU:HG3	1.69	0.57
1:A:344:GLY:HA3	1:A:523:LYS:HB2	1.86	0.57
1:B:110:LEU:HD22	1:B:249:TYR:HE2	1.69	0.57
1:B:940:TYR:HA	1:B:944:LEU:HD12	1.85	0.57
3:F:113:ALA:HB1	3:F:202:LEU:HG	1.86	0.57
1:A:91:ASP:HB3	1:A:255:ALA:HB3	1.87	0.56
1:A:355:TRP:NE1	1:A:382:GLU:OE1	2.37	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:101:ARG:HD2	2:E:111:TYR:HD2	1.70	0.56
1:A:315:VAL:HB	1:A:375:ILE:HB	1.86	0.56
2:C:102:HIS:HA	2:C:109:LEU:HA	1.86	0.56
1:A:493:GLU:HB3	1:A:497:GLY:H	1.71	0.56
1:A:587:PRO:HG3	1:A:695:TRP:HB3	1.88	0.56
1:B:576:PHE:HB2	1:B:629:LEU:HB3	1.86	0.56
2:E:31:ASN:OD1	2:E:80:ASN:ND2	2.38	0.56
1:A:388:GLU:HA	1:A:391:ILE:HD12	1.87	0.56
1:B:601:LYS:HD2	1:B:620:LEU:HB2	1.87	0.56
1:B:642:LEU:HA	1:B:645:ILE:HD12	1.88	0.56
2:E:101:ARG:HB3	2:E:111:TYR:HB2	1.87	0.56
1:B:92:VAL:HG12	1:B:94:ILE:H	1.70	0.55
1:B:822:THR:O	1:B:827:GLU:N	2.32	0.55
1:B:788:ASN:HB2	1:B:851:GLN:HE21	1.71	0.55
1:B:911:ILE:O	1:B:914:GLN:NE2	2.39	0.55
1:A:771:LEU:HB2	1:A:952:HIS:HD2	1.72	0.55
1:B:47:ILE:HG23	1:B:272:VAL:HG22	1.88	0.55
2:E:43:ALA:HB3	2:E:46:LYS:HB2	1.89	0.55
1:B:343:PRO:HG2	1:B:525:PRO:HA	1.89	0.55
1:A:574:LEU:HB2	1:A:631:VAL:HB	1.89	0.55
1:A:571:LYS:HA	1:A:634:TYR:HA	1.89	0.55
1:B:877:MET:O	1:B:933:LYS:NZ	2.34	0.54
1:B:575:ASN:ND2	1:B:904:SER:OG	2.39	0.54
1:B:942:GLU:HG2	1:B:943:MET:HG3	1.88	0.54
2:C:70:ARG:HH12	2:C:90:ARG:HE	1.53	0.54
1:A:600:LEU:HD11	1:A:649:MET:HA	1.89	0.54
1:B:262:GLU:HB2	1:B:267:LEU:HD23	1.89	0.54
2:C:55:TYR:O	2:C:59:GLY:N	2.41	0.54
1:A:194:VAL:HA	1:A:496:TYR:HE1	1.73	0.54
3:F:118:ILE:HA	3:F:135:CYS:HA	1.90	0.54
2:E:72:THR:N	2:E:85:GLN:O	2.40	0.53
1:A:200:ARG:HH21	1:A:498:THR:HB	1.73	0.53
1:B:311:ARG:NH1	1:B:381:GLU:OE2	2.42	0.53
1:A:509:VAL:HG12	1:A:513:TRP:CE2	2.44	0.53
1:B:269:ASN:HA	1:B:272:VAL:HG12	1.90	0.53
1:A:532:PRO:HB3	1:A:636:ASP:HB2	1.89	0.53
1:B:87:SER:HA	1:B:152:ASP:HA	1.90	0.53
1:B:697:LYS:HA	1:B:700:LEU:HD12	1.89	0.53
1:B:800:GLN:HA	1:B:844:GLN:HE21	1.74	0.53
1:A:765:ARG:NH1	1:A:912:ILE:O	2.40	0.53
3:F:121:PRO:HB3	3:F:132:SER:H	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:621:GLN:O	1:A:628:TYR:N	2.38	0.53
1:A:697:LYS:HA	1:A:700:LEU:HD12	1.91	0.53
1:B:71:ASN:HA	1:B:280:ASN:HB2	1.90	0.53
2:E:132:LEU:N	2:E:147:GLY:O	2.38	0.53
1:A:746:GLU:HA	1:A:749:LEU:HD12	1.90	0.52
1:B:57:SER:O	1:B:59:GLU:N	2.43	0.52
1:A:47:ILE:HG22	1:A:272:VAL:HG12	1.91	0.52
3:F:119:PHE:N	3:F:134:VAL:O	2.42	0.52
1:A:638:GLN:O	1:A:642:LEU:CB	2.57	0.52
1:A:934:GLU:HA	1:A:937:ILE:HD12	1.92	0.52
2:E:39:TRP:O	2:E:51:VAL:N	2.39	0.52
1:B:951:ARG:HH21	1:B:953:LYS:HD3	1.74	0.52
1:B:311:ARG:HH21	1:B:380:THR:HA	1.73	0.52
1:A:343:PRO:HG2	1:A:525:PRO:HA	1.91	0.52
1:B:696:THR:N	1:B:699:GLU:OE2	2.33	0.52
1:A:91:ASP:OD2	1:A:146:HIS:ND1	2.43	0.52
1:A:262:GLU:HB2	1:A:267:LEU:HD23	1.91	0.52
3:F:107:ILE:O	3:F:141:TYR:OH	2.21	0.52
3:F:191:LYS:HG2	3:F:212:ARG:HH22	1.75	0.52
1:A:834:PHE:N	1:A:849:ILE:O	2.32	0.52
2:C:70:ARG:NH2	2:C:88:SER:O	2.42	0.52
2:C:72:THR:O	2:C:85:GLN:N	2.41	0.52
1:A:834:PHE:HB3	1:A:849:ILE:HB	1.91	0.51
1:B:858:TYR:O	1:B:861:SER:OG	2.23	0.51
2:E:208:HIS:CE1	2:E:210:PRO:HD2	2.45	0.51
2:E:50:TRP:HB3	3:F:97:ILE:HB	1.91	0.51
3:F:14:ALA:HA	3:F:108:LYS:HG3	1.92	0.51
1:A:315:VAL:O	1:A:375:ILE:N	2.40	0.51
1:A:834:PHE:O	1:A:849:ILE:N	2.30	0.51
1:B:84:ASP:OD2	1:B:896:LYS:NZ	2.44	0.51
1:B:162:LEU:HD11	1:B:275:PHE:HE2	1.75	0.51
1:A:528:ASN:HD21	1:A:530:PHE:HB2	1.75	0.51
1:B:311:ARG:HB3	1:B:379:LEU:HB2	1.93	0.51
2:C:208:HIS:CE1	2:C:210:PRO:HD2	2.45	0.51
1:A:47:ILE:HB	1:A:49:ARG:HH22	1.75	0.51
1:B:382:GLU:O	1:B:386:HIS:ND1	2.22	0.51
1:B:121:TYR:HB3	1:B:126:GLU:HG2	1.93	0.51
1:B:334:LEU:HD11	1:B:467:VAL:HG21	1.93	0.51
1:B:663:LYS:HE3	1:B:704:LEU:HB3	1.93	0.50
1:A:687:ARG:HA	1:A:838:ARG:HH22	1.76	0.50
1:A:696:THR:N	1:A:699:GLU:OE2	2.32	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:9:GLU:OE2	2:E:115:GLY:N	2.45	0.50
3:F:139:ASN:HA	3:F:173:THR:HB	1.93	0.50
1:A:63:GLU:HB2	1:A:79:SER:HB3	1.94	0.50
1:B:780:GLN:HA	1:B:957:HIS:HB2	1.92	0.50
2:E:127:PRO:HB3	2:E:153:TYR:HB3	1.94	0.50
1:B:301:LEU:HA	1:B:478:VAL:HB	1.94	0.50
1:B:877:MET:HB3	1:B:933:LYS:HE3	1.93	0.50
3:F:86:THR:HA	3:F:104:LYS:HA	1.94	0.50
1:B:785:VAL:O	1:B:961:ARG:NH1	2.45	0.49
1:A:724:HIS:HB2	1:A:758:LEU:HD12	1.92	0.49
1:B:349:GLU:O	1:B:353:LYS:N	2.42	0.49
3:F:152:ASP:HA	3:F:192:VAL:HB	1.93	0.49
1:A:762:GLN:HB3	1:B:695:TRP:HH2	1.77	0.49
3:D:38:GLN:HB2	3:D:48:LEU:HD11	1.94	0.49
1:A:66:GLY:HA3	1:A:446:LEU:HB2	1.94	0.49
1:A:76:LEU:HB3	1:A:257:VAL:HG23	1.95	0.49
2:E:74:SER:O	2:E:83:TYR:N	2.40	0.49
1:A:304:ILE:HD11	1:A:479:ALA:HB1	1.94	0.49
1:A:776:TRP:CD1	1:A:953:LYS:HG2	2.47	0.49
1:B:493:GLU:HB3	1:B:497:GLY:H	1.76	0.49
3:D:113:ALA:HB1	3:D:202:LEU:HG	1.95	0.49
3:F:18:ASP:H	3:F:79:LEU:H	1.61	0.49
1:A:255:ALA:HB1	1:A:441:LEU:HD23	1.94	0.49
1:B:802:THR:HG23	1:B:924:GLU:HG2	1.94	0.49
3:D:38:GLN:O	3:D:46:LYS:N	2.46	0.49
3:F:137:LEU:HB2	3:F:176:LEU:HB3	1.95	0.49
1:A:578:PHE:HB2	1:A:627:MET:HB2	1.95	0.49
1:A:251:SER:HB3	1:A:281:LYS:HB2	1.95	0.49
1:A:833:VAL:HA	1:A:850:ILE:HG12	1.95	0.49
1:B:783:ASN:HD21	1:B:786:HIS:HB2	1.77	0.49
3:D:21:THR:HG22	3:D:75:THR:HG23	1.94	0.49
1:A:291:HIS:O	1:A:294:GLN:NE2	2.41	0.48
1:A:317:PHE:N	1:A:373:PHE:O	2.39	0.48
1:A:638:GLN:O	1:A:642:LEU:HB2	2.13	0.48
1:A:824:ARG:HB2	1:A:833:VAL:HG11	1.94	0.48
2:E:70:ARG:NH2	2:E:88:SER:O	2.47	0.48
1:B:117:GLY:HA3	1:B:173:LEU:HB2	1.95	0.48
1:B:350:LEU:O	1:B:354:GLY:N	2.47	0.48
1:A:948:ALA:HB3	1:A:951:ARG:HB2	1.95	0.48
2:C:101:ARG:HD2	2:C:111:TYR:HD2	1.78	0.48
1:A:46:ALA:HB1	1:A:70:ALA:HB3	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:72:THR:O	2:E:85:GLN:N	2.47	0.48
1:A:78:ILE:HB	1:A:259:LEU:HA	1.95	0.48
3:D:76:ILE:HG21	3:D:79:LEU:HD23	1.95	0.48
3:F:142:PRO:O	3:F:199:HIS:NE2	2.36	0.48
1:A:395:PHE:HD1	1:A:398:ILE:HD12	1.78	0.48
3:D:142:PRO:O	3:D:199:HIS:NE2	2.39	0.48
1:A:790:GLY:N	1:A:958:VAL:O	2.41	0.47
1:A:305:VAL:HG22	1:A:485:PHE:HB2	1.97	0.47
1:B:532:PRO:HB3	1:B:636:ASP:HB2	1.97	0.47
3:D:118:ILE:HD12	3:D:195:CYS:HB2	1.96	0.47
2:E:41:ARG:HG2	2:E:97:TYR:CE1	2.49	0.47
1:A:419:ALA:O	1:A:423:ARG:HB2	2.14	0.47
3:D:32:SER:HB2	3:D:52:ALA:HB3	1.96	0.47
1:A:587:PRO:HB3	1:A:700:LEU:HD23	1.95	0.47
1:A:92:VAL:HG12	1:A:94:ILE:H	1.79	0.47
2:E:151:LYS:HE3	2:E:179:GLN:HE22	1.78	0.47
2:E:153:TYR:HB2	2:E:208:HIS:CD2	2.49	0.47
1:A:197:ASP:HA	1:A:200:ARG:HD3	1.97	0.47
1:A:777:PHE:HD2	1:A:998:PHE:HE1	1.62	0.47
1:B:78:ILE:HB	1:B:259:LEU:HA	1.97	0.47
3:D:125:GLN:HE22	3:D:132:SER:HB2	1.80	0.47
3:F:145:ALA:HB2	3:F:199:HIS:HD2	1.78	0.47
1:A:264:LEU:O	1:A:268:THR:OG1	2.25	0.47
1:A:594:TYR:HD1	1:A:622:ASN:HD21	1.62	0.47
1:B:317:PHE:N	1:B:373:PHE:O	2.36	0.47
1:A:899:LYS:HG3	1:A:901:SER:H	1.80	0.47
1:A:858:TYR:OH	1:A:862:ARG:NH1	2.47	0.47
1:B:183:VAL:HG23	1:B:226:LEU:HD22	1.97	0.47
1:B:637:LYS:HG2	1:B:640:ILE:HD12	1.97	0.47
2:C:41:ARG:HG2	2:C:97:TYR:CE1	2.50	0.47
1:B:609:TYR:O	1:B:613:LEU:HB2	2.15	0.46
1:B:781:GLN:O	1:B:959:LEU:N	2.48	0.46
2:E:153:TYR:HE2	2:E:157:PRO:HA	1.80	0.46
1:A:350:LEU:HD23	1:A:350:LEU:HA	1.81	0.46
1:A:460:ARG:HB3	1:A:463:LEU:HD13	1.98	0.46
1:A:894:LEU:HD11	1:A:925:VAL:HG21	1.96	0.46
1:B:746:GLU:HA	1:B:749:LEU:HD12	1.96	0.46
1:A:361:GLY:HA2	1:A:375:ILE:HA	1.97	0.46
1:A:508:GLU:O	1:A:512:LYS:HB2	2.16	0.46
1:A:795:TYR:N	1:A:846:LEU:O	2.46	0.46
1:B:91:ASP:O	1:B:255:ALA:N	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:ASP:O	1:A:184:ASN:ND2	2.48	0.46
1:A:387:VAL:HA	1:A:390:ILE:HD12	1.97	0.46
1:A:429:ARG:HG3	1:A:431:ARG:H	1.80	0.46
1:B:854:LYS:HD2	1:B:858:TYR:HE2	1.81	0.46
2:C:101:ARG:HE	2:C:103:TYR:HA	1.81	0.46
3:D:86:THR:HA	3:D:104:LYS:HA	1.98	0.46
3:D:145:ALA:HB2	3:D:199:HIS:HD2	1.81	0.46
1:A:110:LEU:O	1:A:114:LEU:N	2.47	0.46
1:A:415:LYS:NZ	1:A:456:LEU:O	2.35	0.46
1:A:461:PRO:HA	1:A:464:ILE:HD12	1.97	0.46
1:A:573:ASN:HD22	1:A:900:LEU:HD22	1.81	0.46
1:B:767:ARG:HD3	1:B:1005:PHE:O	2.15	0.46
1:A:382:GLU:O	1:A:386:HIS:ND1	2.35	0.46
2:C:179:GLN:HE21	2:C:183:LEU:HB2	1.81	0.45
3:D:14:ALA:HA	3:D:108:LYS:HB2	1.98	0.45
1:B:191:GLU:HA	1:B:194:VAL:HG23	1.98	0.45
1:A:670:LEU:HD11	1:A:704:LEU:HD12	1.98	0.45
1:B:769:VAL:HA	1:B:1004:LEU:HD23	1.98	0.45
1:B:158:LEU:O	1:B:162:LEU:HB2	2.15	0.45
1:B:203:GLN:HG2	1:B:496:TYR:HE2	1.81	0.45
1:B:790:GLY:N	1:B:958:VAL:O	2.44	0.45
3:F:16:VAL:HG13	3:F:80:GLN:HA	1.98	0.45
1:A:205:GLU:OE1	1:A:364:LYS:NZ	2.49	0.45
1:A:994:ASN:O	1:A:998:PHE:N	2.46	0.45
2:C:206:VAL:HB	2:C:215:VAL:H	1.81	0.45
1:A:711:ARG:NH2	1:A:715:PHE:HB2	2.27	0.45
1:B:601:LYS:HE3	1:B:618:TYR:CE2	2.52	0.45
1:B:623:THR:N	1:B:626:GLY:O	2.36	0.45
1:A:711:ARG:NE	1:B:718:GLN:OE1	2.33	0.45
1:A:552:LYS:HB3	1:A:559:LEU:HB3	1.99	0.45
1:B:564:ASP:OD2	1:B:568:PHE:N	2.50	0.45
3:D:35:ALA:HB3	3:D:90:GLN:HE21	1.82	0.45
1:A:238:ARG:HA	1:A:241:LEU:HD12	1.97	0.44
1:A:437:ILE:HA	1:A:440:ILE:HG13	1.98	0.44
1:B:446:LEU:HD12	1:B:447:GLU:HG3	1.99	0.44
1:B:81:PRO:HA	1:B:261:ARG:HG3	1.99	0.44
1:A:57:SER:O	1:A:59:GLU:N	2.49	0.44
1:B:552:LYS:HB3	1:B:559:LEU:HB3	1.98	0.44
1:B:805:ASN:N	1:B:844:GLN:HE22	2.15	0.44
3:F:114:PRO:HG2	3:F:202:LEU:HD12	1.99	0.44
3:F:117:PHE:HD2	3:F:136:LEU:HD23	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:92:VAL:HG22	1:B:254:MET:HG3	1.99	0.44
1:B:489:THR:HA	1:B:501:LYS:HG2	1.99	0.44
3:D:14:ALA:HA	3:D:108:LYS:HE2	1.99	0.44
1:B:387:VAL:HA	1:B:390:ILE:HD12	1.99	0.44
1:B:836:GLY:N	1:B:847:ARG:O	2.50	0.44
1:A:692:GLU:HB2	1:A:767:ARG:H	1.82	0.44
1:B:210:ASN:HB3	1:B:213:HIS:HB2	1.99	0.44
1:B:554:THR:OG1	1:B:557:SER:N	2.40	0.44
1:B:64:TYR:HE1	1:B:78:ILE:HG23	1.82	0.44
2:C:128:SER:O	2:C:151:LYS:N	2.43	0.44
3:D:143:ARG:NE	3:D:166:GLU:OE2	2.50	0.44
2:E:162:TRP:O	2:E:166:ALA:N	2.45	0.44
1:A:510:ILE:O	1:A:513:TRP:HB2	2.18	0.44
1:B:577:GLU:N	1:B:726:GLU:O	2.35	0.44
2:E:86:MET:HB3	2:E:89:LEU:HD21	2.00	0.44
1:B:710:PRO:HA	1:B:713:LYS:HE2	1.99	0.44
1:A:346:LEU:HD22	1:A:397:TYR:HB2	1.99	0.44
1:B:681:HIS:CE1	1:B:685:TYR:HE2	2.35	0.43
2:C:86:MET:HB3	2:C:89:LEU:HD21	1.99	0.43
1:A:65:ARG:NH2	1:A:265:ASP:OD1	2.35	0.43
1:A:829:LEU:HD21	1:A:862:ARG:HH21	1.83	0.43
1:A:940:TYR:O	1:A:945:ALA:N	2.50	0.43
1:B:125:ASN:O	1:B:129:GLN:N	2.42	0.43
1:B:69:LEU:HB2	1:B:73:ILE:HB	2.00	0.43
1:A:472:ARG:HB2	1:A:474:GLU:HG2	2.00	0.43
1:A:63:GLU:O	1:A:79:SER:N	2.51	0.43
2:C:162:TRP:O	2:C:166:ALA:N	2.48	0.43
2:E:105:ALA:HB3	2:E:108:GLY:HA3	2.00	0.43
1:B:482:SER:HB3	1:B:485:PHE:CZ	2.53	0.43
1:B:619:ASP:HB3	1:B:630:SER:HB3	1.99	0.43
3:D:3:ILE:HG23	3:D:27:SER:HB2	2.01	0.43
2:E:38:HIS:CD2	2:E:102:HIS:HB2	2.54	0.43
1:A:318:PRO:HA	1:A:372:PHE:HD1	1.84	0.43
1:A:317:PHE:O	1:A:373:PHE:N	2.51	0.43
1:A:482:SER:HB3	1:A:485:PHE:CZ	2.53	0.43
1:A:528:ASN:OD1	1:A:529:GLU:N	2.52	0.43
1:A:586:ASP:CG	1:A:589:HIS:HD1	2.22	0.43
1:A:93:HIS:HB3	1:A:253:LEU:HB3	2.01	0.43
1:B:238:ARG:HA	1:B:241:LEU:HD12	2.00	0.43
3:F:63:PHE:HD1	3:F:76:ILE:HG12	1.83	0.43
3:D:67:ARG:HG3	3:D:72:TYR:CE1	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:146:LEU:HD23	2:E:192:VAL:HG21	2.00	0.43
1:A:406:PRO:HB2	1:A:459:PHE:HZ	1.82	0.43
1:A:711:ARG:O	1:A:711:ARG:NH2	2.51	0.43
1:B:770:GLN:HB3	1:B:1003:PRO:HB2	2.00	0.43
1:B:620:LEU:HA	1:B:620:LEU:HD13	1.80	0.43
1:B:689:LEU:HD23	1:B:689:LEU:HA	1.84	0.43
1:B:825:THR:O	1:B:828:GLN:NE2	2.39	0.43
1:B:547:TYR:CD1	1:B:918:PHE:HB3	2.54	0.43
3:D:32:SER:HB3	3:D:67:ARG:HD3	1.98	0.43
1:A:556:MET:HB2	1:A:757:PRO:HG3	2.01	0.43
1:A:783:ASN:O	1:A:961:ARG:HA	2.19	0.43
1:A:794:TYR:OH	1:A:845:GLY:HA3	2.19	0.43
1:B:517:ASP:OD1	1:B:517:ASP:N	2.48	0.43
1:B:575:ASN:OD1	1:B:628:TYR:OH	2.37	0.43
1:A:152:ASP:N	1:A:152:ASP:OD1	2.51	0.43
1:A:483:LYS:HE3	1:A:483:LYS:HB2	1.92	0.43
1:A:803:SER:HA	1:A:927:TYR:CE2	2.54	0.43
1:B:346:LEU:HD22	1:B:397:TYR:HB2	2.00	0.43
2:E:14:LEU:HB3	2:E:121:PHE:HE1	1.83	0.43
1:A:586:ASP:OD2	1:A:589:HIS:ND1	2.44	0.42
1:A:623:THR:N	1:A:626:GLY:O	2.35	0.42
1:B:213:HIS:CG	1:B:292:PRO:HG3	2.54	0.42
1:A:462:ASP:O	1:A:466:MET:HB2	2.19	0.42
1:A:508:GLU:HB3	2:C:104:SER:HA	2.00	0.42
1:B:197:ASP:HA	1:B:200:ARG:HD3	2.01	0.42
3:F:119:PHE:HB2	3:F:134:VAL:HB	2.02	0.42
1:A:364:LYS:HB3	1:A:372:PHE:HB2	2.01	0.42
1:A:722:ARG:HH21	1:A:756:LYS:HD3	1.85	0.42
1:B:71:ASN:HB2	1:B:278:VAL:HG12	2.01	0.42
1:B:575:ASN:HD22	1:B:728:LEU:HD23	1.84	0.42
1:B:99:ASP:O	1:B:217:LYS:NZ	2.48	0.42
1:A:940:TYR:HA	1:A:944:LEU:HB2	2.01	0.42
1:B:470:LYS:O	1:B:475:ASN:ND2	2.41	0.42
1:B:857:HIS:CD2	1:B:857:HIS:H	2.37	0.42
1:B:462:ASP:OD1	1:B:462:ASP:N	2.52	0.42
3:F:65:GLY:HA2	3:F:74:LEU:HA	2.00	0.42
1:A:297:HIS:HB2	1:A:477:ARG:HH12	1.85	0.42
1:A:554:THR:HB	1:A:557:SER:H	1.84	0.42
1:A:621:GLN:HB3	1:A:628:TYR:HB3	2.02	0.42
1:B:402:ARG:HA	1:B:468:LEU:HD11	2.01	0.42
1:B:797:THR:HG21	1:B:808:LEU:HD12	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:ARG:HB3	1:A:379:LEU:HB2	2.02	0.42
1:A:341:GLU:HB2	1:A:609:TYR:CG	2.54	0.42
1:A:715:PHE:CE2	1:A:719:LEU:HB2	2.54	0.42
1:A:762:GLN:HB3	1:B:695:TRP:CH2	2.55	0.42
1:B:175:ASP:N	1:B:175:ASP:OD1	2.47	0.42
3:D:189:LYS:HE3	3:D:190:HIS:CE1	2.55	0.42
1:A:202:PHE:O	1:A:206:LYS:HG2	2.20	0.42
1:A:574:LEU:HB3	1:A:576:PHE:CE2	2.55	0.42
1:A:64:TYR:HE1	1:A:78:ILE:HG23	1.85	0.42
1:A:881:ALA:O	1:A:885:HIS:CB	2.65	0.42
1:A:809:GLU:HB3	1:A:889:LEU:HD11	2.00	0.42
2:E:145:ALA:HB3	3:F:119:PHE:HZ	1.85	0.42
1:A:619:ASP:HB3	1:A:630:SER:HB3	2.01	0.42
1:B:297:HIS:HB2	1:B:477:ARG:HH12	1.84	0.42
2:E:156:GLU:HB2	2:E:209:LYS:HB2	2.02	0.42
1:B:100:PRO:HA	1:B:101:PRO:HD3	1.93	0.41
3:F:164:VAL:HG22	3:F:176:LEU:HG	2.01	0.41
1:A:822:THR:O	1:A:826:LYS:HB3	2.20	0.41
1:A:90:LEU:HD13	1:A:165:PHE:HE2	1.85	0.41
1:A:910:GLU:O	1:A:914:GLN:N	2.53	0.41
1:B:833:VAL:HG13	1:B:850:ILE:HG12	2.00	0.41
2:C:17:PRO:HG3	2:C:120:VAL:HG12	2.02	0.41
3:D:164:VAL:HG22	3:D:176:LEU:HG	2.01	0.41
1:A:562:LYS:NZ	1:A:563:GLN:O	2.45	0.41
1:B:402:ARG:CZ	1:B:472:ARG:HH12	2.33	0.41
1:B:671:ASN:HB3	1:B:701:LYS:NZ	2.35	0.41
3:D:139:ASN:HA	3:D:173:THR:HB	2.03	0.41
2:E:134:PRO:HB2	2:E:197:LEU:HD22	2.02	0.41
3:F:162:GLU:HG2	3:F:176:LEU:HD21	2.03	0.41
1:B:110:LEU:O	1:B:114:LEU:N	2.54	0.41
1:B:832:ILE:HB	1:B:851:GLN:HB3	2.02	0.41
1:B:995:MET:HA	1:B:998:PHE:HB3	2.03	0.41
3:F:38:GLN:HB2	3:F:48:LEU:HD11	2.01	0.41
1:B:224:TYR:HA	1:B:228:THR:HB	2.03	0.41
1:B:368:ARG:HH12	1:B:442:HIS:HE1	1.67	0.41
1:B:691:THR:OG1	1:B:692:GLU:N	2.53	0.41
1:B:870:MET:HA	1:B:873:SER:HB3	2.02	0.41
1:A:359:LEU:HA	1:A:377:VAL:HA	2.01	0.41
1:A:599:LEU:HD12	1:A:662:ILE:HG13	2.03	0.41
2:C:75:ALA:HA	2:C:82:ALA:HA	2.03	0.41
2:C:112:TRP:CZ2	3:D:45:PRO:HB2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:47:LEU:HD21	3:D:50:TYR:HB3	2.02	0.41
1:B:638:GLN:O	1:B:642:LEU:CB	2.68	0.41
1:B:800:GLN:HA	1:B:844:GLN:NE2	2.34	0.41
3:F:3:ILE:HG23	3:F:27:SER:HB2	2.03	0.41
3:F:21:THR:HG22	3:F:75:THR:HG23	2.03	0.41
1:A:638:GLN:O	1:A:642:LEU:HB3	2.20	0.41
1:A:65:ARG:HB3	1:A:77:LEU:HD12	2.03	0.41
1:B:805:ASN:OD1	1:B:844:GLN:NE2	2.54	0.41
2:E:23:LEU:HB2	2:E:84:LEU:HB3	2.02	0.41
1:A:102:ASN:OD1	1:A:229:ARG:NH1	2.53	0.41
1:A:244:PHE:O	1:A:248:TYR:HB2	2.21	0.41
1:A:395:PHE:CZ	1:A:473:PRO:HG3	2.56	0.41
1:A:692:GLU:HB2	1:A:767:ARG:N	2.36	0.41
1:A:835:SER:HA	1:A:848:PHE:HA	2.02	0.41
1:B:266:ASP:OD1	1:B:267:LEU:N	2.52	0.41
3:D:38:GLN:HG3	3:D:87:TYR:HE1	1.86	0.41
2:E:21:LEU:O	2:E:85:GLN:NE2	2.53	0.41
3:F:15:SER:N	3:F:18:ASP:OD2	2.48	0.41
1:A:137:SER:N	1:A:152:ASP:OD1	2.54	0.40
1:A:266:ASP:OD1	1:A:267:LEU:N	2.50	0.40
1:A:324:LYS:HB3	1:A:324:LYS:HE2	1.83	0.40
1:A:691:THR:OG1	1:A:692:GLU:N	2.54	0.40
1:A:771:LEU:HD12	1:A:952:HIS:HB3	2.03	0.40
1:B:300:GLN:HE22	1:B:504:ALA:HB2	1.86	0.40
1:B:807:PHE:HB3	1:B:939:PHE:CE1	2.56	0.40
3:F:186:ASP:O	3:F:193:TYR:OH	2.39	0.40
1:A:797:THR:OG1	1:A:798:ASP:N	2.54	0.40
1:B:109:PHE:HB2	1:B:226:LEU:HD11	2.03	0.40
1:B:677:GLN:HA	1:B:678:PRO:HD3	1.86	0.40
1:A:600:LEU:HD21	1:A:649:MET:HB2	2.03	0.40
1:A:620:LEU:HA	1:A:620:LEU:HD13	1.86	0.40
1:A:855:PRO:HA	1:A:856:PRO:HD3	1.81	0.40
1:B:699:GLU:HG2	1:B:699:GLU:H	1.75	0.40
1:B:833:VAL:HG22	1:B:850:ILE:HG23	2.02	0.40
1:A:512:LYS:NZ	2:C:106:VAL:O	2.42	0.40
1:A:1006:PRO:HB3	1:B:1001:GLY:C	2.42	0.40
1:A:87:SER:O	1:A:259:LEU:N	2.54	0.40
1:A:90:LEU:O	1:A:149:TYR:N	2.44	0.40
1:B:141:PHE:HD1	1:B:150:TYR:HE1	1.70	0.40
1:B:562:LYS:NZ	1:B:563:GLN:O	2.52	0.40
1:B:574:LEU:HB3	1:B:576:PHE:CE2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:109:ARG:HH22	3:D:112:ALA:HB2	1.87	0.40
3:D:134:VAL:HG22	3:D:179:THR:HG22	2.03	0.40
3:D:151:VAL:HG22	3:D:156:GLN:HB2	2.03	0.40
2:E:14:LEU:HB3	2:E:121:PHE:CE1	2.56	0.40
2:E:6:GLN:HB3	2:E:28:SER:HB2	2.04	0.40
3:F:147:VAL:HG21	3:F:176:LEU:HD22	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	945/966 (98%)	913 (97%)	32 (3%)	0	100	100
1	B	936/966 (97%)	902 (96%)	34 (4%)	0	100	100
2	C	213/218 (98%)	208 (98%)	5 (2%)	0	100	100
2	E	213/218 (98%)	204 (96%)	9 (4%)	0	100	100
3	D	209/211 (99%)	205 (98%)	4 (2%)	0	100	100
3	F	209/211 (99%)	206 (99%)	3 (1%)	0	100	100
All	All	2725/2790 (98%)	2638 (97%)	87 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	835/861 (97%)	830 (99%)	5 (1%)	89	94
1	B	827/861 (96%)	821 (99%)	6 (1%)	87	93
2	C	179/181 (99%)	178 (99%)	1 (1%)	89	94
2	E	179/181 (99%)	178 (99%)	1 (1%)	89	94
3	D	187/187 (100%)	187 (100%)	0	100	100
3	F	187/187 (100%)	187 (100%)	0	100	100
All	All	2394/2458 (97%)	2381 (100%)	13 (0%)	91	95

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	ARG
1	A	649	MET
1	A	783	ASN
1	A	824	ARG
1	A	896	LYS
1	B	49	ARG
1	B	229	ARG
1	B	649	MET
1	B	824	ARG
1	B	838	ARG
1	B	896	LYS
2	C	101	ARG
2	E	101	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	93	HIS
1	A	190	HIS
1	A	332	HIS
1	A	681	HIS
1	A	762	GLN
1	A	952	HIS
1	B	294	GLN
1	B	300	GLN
1	B	332	HIS
1	B	681	HIS
1	B	730	HIS
1	B	732	ASN

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Mol	Chain	Res	Type
1	B	743	GLN
1	B	762	GLN
1	B	844	GLN
1	B	851	GLN
1	B	857	HIS
3	D	90	GLN
3	D	125	GLN
2	E	179	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.