



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:09 PM GMT

PDB ID : 3T2N
Title : Human hepsin protease in complex with the Fab fragment of an inhibitory antibody
Authors : Koschubs, T.; Dengl, S.; Duerr, H.; Kaluza, K.; Georges, G.; Hartl, C.; Jennewein, S.; Lanzendoerfer, M.; Auer, J.; Stern, A.; Huang, K.-S.; Kostrewa, D.; Ries, S.; Hansen, S.; Kohnert, U.; Cramer, P.; Mundigl, O.
Deposited on : 2011-07-22
Resolution : 2.55 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

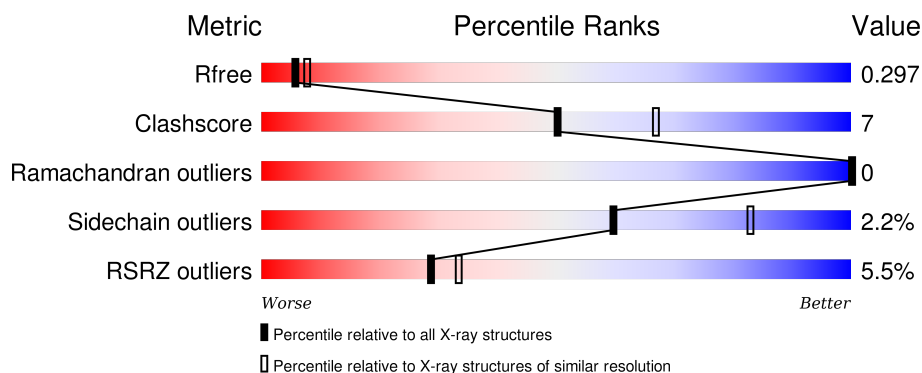
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	4549 (2.58-2.50)
Clashscore	102246	5292 (2.58-2.50)
Ramachandran outliers	100387	5194 (2.58-2.50)
Sidechain outliers	100360	5196 (2.58-2.50)
RSRZ outliers	91569	4561 (2.58-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	372	<div> <div>7%</div> <div> <div></div> <div>75%</div> <div>16%</div> <div>9%</div> </div> </div>
1	B	372	<div> <div>6%</div> <div> <div></div> <div>75%</div> <div>15%</div> <div>10%</div> </div> </div>
2	H	225	<div> <div>4%</div> <div> <div></div> <div>75%</div> <div>15%</div> <div>10%</div> </div> </div>
2	I	225	<div> <div>3%</div> <div> <div></div> <div>79%</div> <div>12%</div> <div>9%</div> </div> </div>
3	L	215	<div> <div>2%</div> <div> <div></div> <div>81%</div> <div>16%</div> <div>3%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
3	M	215	<div><div></div><div>6%</div><div>81%</div><div>16%</div><div>..</div></div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 11749 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 Serine protease hepsin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	339	Total	C	N	O	S	0	0	0
			2605	1638	470	478	19			
1	B	335	Total	C	N	O	S	0	0	0
			2576	1622	464	471	19			

- Molecule 2 is a protein called Antibody, Fab fragment, Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	203	Total	C	N	O	S	0	0	0
			1535	980	250	299	6			
2	I	205	Total	C	N	O	S	0	0	0
			1547	987	252	302	6			

- Molecule 3 is a protein called Antibody, Fab fragment, Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	211	Total	C	N	O	S	0	0	0
			1582	996	267	315	4			
3	M	212	Total	C	N	O	S	0	0	0
			1589	1000	268	317	4			

- Molecule 4 is water.

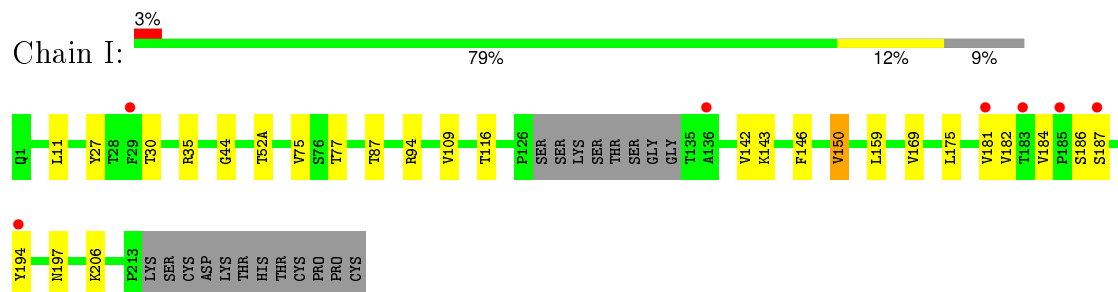
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	64	Total	O	0	0
			64	64		
4	B	74	Total	O	0	0
			74	74		
4	H	46	Total	O	0	0
			46	46		
4	I	36	Total	O	0	0
			36	36		

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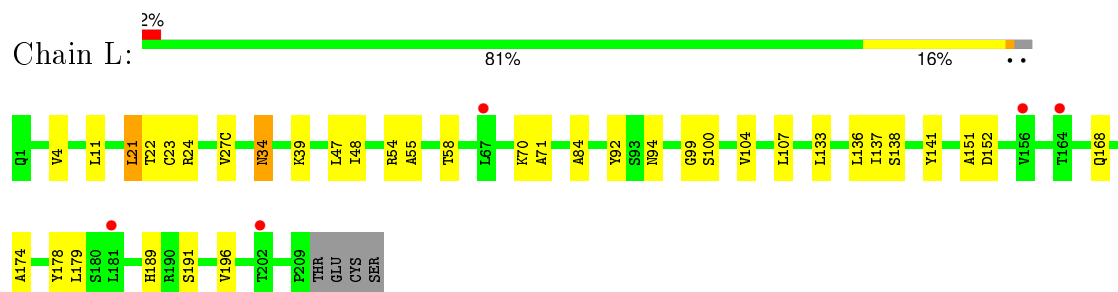
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	40	Total	O	0	0
			40	40		
4	M	55	Total	O	0	0
			55	55		

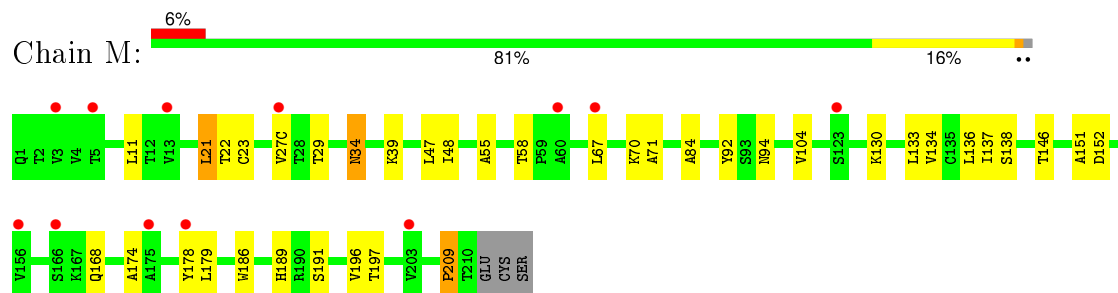
- Molecule 2: Antibody, Fab fragment, Heavy Chain



- Molecule 3: Antibody, Fab fragment, Light Chain



- Molecule 3: Antibody, Fab fragment, Light Chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	62.98Å 66.58Å 108.33Å 88.71° 94.30° 104.53°	Depositor
Resolution (Å)	47.33 – 2.55 47.33 – 2.55	Depositor EDS
% Data completeness (in resolution range)	98.7 (47.33-2.55) 79.1 (47.33-2.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.69 (at 2.54Å)	Xtriage
Refinement program	BUSTER 2.11.1	Depositor
R, R_{free}	0.242 , 0.275 0.257 , 0.297	Depositor DCC
R_{free} test set	2772 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	43.1	Xtriage
Anisotropy	0.637	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 54657 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	11749	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.14% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

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 > 5$	RMSZ	# $ Z > 5$
1	A	0.52	0/2665	0.75	3/3615 (0.1%)
1	B	0.54	1/2637 (0.0%)	0.74	0/3579
2	H	0.56	0/1575	0.78	0/2148
2	I	0.50	0/1587	0.77	0/2165
3	L	0.55	0/1623	0.74	0/2222
3	M	0.59	0/1630	0.75	1/2232 (0.0%)
All	All	0.54	1/11717 (0.0%)	0.75	4/15961 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	206	PRO	N-CD	-5.43	1.40	1.47

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	107	ARG	NE-CZ-NH2	-7.07	116.76	120.30
1	A	107	ARG	NE-CZ-NH1	6.30	123.45	120.30
1	A	167	ARG	CD-NE-CZ	5.99	131.99	123.60
3	M	209	PRO	C-N-CA	5.32	135.00	121.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	167	ARG	Sidechain
1	B	257	ASP	Sidechain

5.2 Too-close contacts [i](#)

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	2605	0	2536	43	0
1	B	2576	0	2500	37	0
2	H	1535	0	1502	22	0
2	I	1547	0	1514	16	0
3	L	1582	0	1541	25	0
3	M	1589	0	1548	27	0
4	A	64	0	0	0	0
4	B	74	0	0	0	0
4	H	46	0	0	0	0
4	I	36	0	0	0	0
4	L	40	0	0	1	0
4	M	55	0	0	1	0
All	All	11749	0	11141	158	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 7.

All (158) 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)
1:B:317:ILE:HD11	1:B:340:GLY:HA2	1.52	0.92
2:I:181:VAL:HG11	3:M:136:LEU:HD13	1.54	0.89
1:A:383:LEU:HG	1:A:383:LEU:O	1.76	0.85
2:H:75:VAL:HG13	2:H:77:THR:HG22	1.66	0.77
2:H:181:VAL:HG11	3:L:136:LEU:HD13	1.66	0.76
1:A:383:LEU:HD21	3:L:54:ARG:H	1.51	0.74
2:I:75:VAL:HG13	2:I:77:THR:HG22	1.70	0.73
1:A:169:THR:HG22	1:A:222:VAL:HG21	1.70	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:47:LEU:HD22	3:L:58:THR:HG22	1.76	0.67
2:I:181:VAL:CG1	3:M:136:LEU:HD13	2.24	0.67
1:A:317:ILE:HD11	1:A:340:GLY:HA2	1.77	0.67
2:H:67:PHE:N	2:H:67:PHE:CD1	2.63	0.65
3:M:47:LEU:HD22	3:M:58:THR:HG22	1.77	0.65
1:B:338:CYS:SG	1:B:387:PRO:HB2	2.37	0.65
2:H:61:ASP:OD1	2:H:64:LYS:NZ	2.28	0.64
3:M:134:VAL:HG12	3:M:136:LEU:HD21	1.78	0.64
1:A:105:ASP:O	1:A:109:ALA:HB3	1.99	0.63
3:L:100:SER:HB2	4:L:225:HOH:O	1.97	0.62
2:H:61:ASP:HA	2:H:64:LYS:HD2	1.80	0.62
1:B:196:ASP:O	1:B:262:HIS:HD2	1.83	0.62
1:A:169:THR:HG22	1:A:222:VAL:CG2	2.31	0.60
1:B:84:ARG:HB2	1:B:125:LEU:HD21	1.83	0.60
1:A:196:ASP:O	1:A:262:HIS:HD2	1.84	0.60
3:M:151:ALA:O	3:M:191:SER:O	2.21	0.59
3:L:151:ALA:O	3:L:191:SER:O	2.20	0.59
3:M:23:CYS:HB3	3:M:71:ALA:HB3	1.84	0.59
3:L:23:CYS:HB3	3:L:71:ALA:HB3	1.84	0.58
3:L:168:GLN:HE21	3:L:174:ALA:HB2	1.68	0.58
3:M:168:GLN:HE21	3:M:174:ALA:HB2	1.68	0.58
1:B:222:VAL:CG2	1:B:309:LEU:HB3	2.34	0.58
1:A:383:LEU:HD21	3:L:54:ARG:N	2.19	0.57
3:M:11:LEU:HD12	3:M:21:LEU:HD22	1.86	0.57
1:A:322:CYS:HB2	1:A:332:ILE:HD11	1.86	0.57
1:B:341:TYR:CD2	3:M:29:THR:HG21	2.40	0.57
1:A:105:ASP:O	1:A:109:ALA:CB	2.53	0.57
1:B:322:CYS:HB2	1:B:332:ILE:HD11	1.87	0.57
2:H:181:VAL:CG1	3:L:136:LEU:HD13	2.33	0.57
3:L:11:LEU:HD12	3:L:21:LEU:HD22	1.87	0.56
1:A:222:VAL:CG2	1:A:309:LEU:HB3	2.37	0.55
3:L:55:ALA:HB3	3:L:58:THR:HG23	1.89	0.55
1:A:142:ARG:HG2	1:A:144:ARG:HG3	1.89	0.55
3:M:152:ASP:OD2	3:M:189:HIS:HD2	1.91	0.54
3:M:55:ALA:HB3	3:M:58:THR:HG23	1.89	0.54
1:A:333:LYS:HB3	1:A:334:PRO:HD2	1.90	0.54
1:A:128:THR:HG21	1:A:134:VAL:HG11	1.89	0.54
2:I:197:ASN:HB3	2:I:206:LYS:HE2	1.90	0.53
3:M:134:VAL:CG1	3:M:136:LEU:HD21	2.38	0.53
1:B:333:LYS:HB3	1:B:334:PRO:HD2	1.90	0.53
3:M:138:SER:HB2	3:M:168:GLN:HE22	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:27:TYR:CZ	2:H:94:ARG:HD2	2.44	0.52
2:H:146:PHE:HB2	2:H:175:LEU:HD23	1.90	0.52
1:A:373:GLY:HA2	1:A:391:THR:O	2.10	0.52
2:H:67:PHE:H	2:H:67:PHE:HD1	1.56	0.52
2:H:197:ASN:HB3	2:H:206:LYS:HE2	1.91	0.52
1:A:167:ARG:O	1:A:310:GLN:HA	2.10	0.52
1:B:385:GLN:HB3	3:M:29:THR:HG23	1.92	0.51
2:I:27:TYR:CZ	2:I:94:ARG:HD2	2.45	0.51
1:B:222:VAL:HG22	1:B:309:LEU:HB3	1.93	0.51
2:I:146:PHE:HB2	2:I:175:LEU:HD23	1.92	0.51
3:M:137:ILE:HG12	3:M:196:VAL:HG21	1.92	0.51
3:M:92:TYR:O	3:M:94:ASN:N	2.41	0.51
1:A:326:ASP:O	2:H:35:ARG:NH2	2.43	0.51
1:B:373:GLY:HA2	1:B:391:THR:O	2.11	0.51
1:B:341:TYR:O	1:B:342:PRO:C	2.48	0.50
3:M:22:THR:HB	3:M:70:LYS:HD2	1.92	0.50
2:I:30:THR:HA	2:I:52(A):THR:HG22	1.93	0.50
1:B:57:SER:HB2	1:B:62:ARG:HG2	1.94	0.50
1:B:212:LEU:HD12	1:B:412:GLY:HA2	1.94	0.50
1:B:128:THR:HG21	1:B:134:VAL:HG11	1.93	0.50
1:B:157:LYS:HB2	1:B:271:GLU:HA	1.94	0.50
1:B:326:ASP:O	2:I:35:ARG:NH2	2.44	0.50
2:I:87:THR:HA	2:I:109:VAL:O	2.12	0.50
1:A:125:LEU:N	1:A:126:PRO:HD2	2.27	0.50
3:L:22:THR:HB	3:L:70:LYS:HD2	1.93	0.50
3:L:152:ASP:OD2	3:L:189:HIS:HD2	1.95	0.49
3:L:138:SER:HB2	3:L:168:GLN:HE22	1.77	0.49
1:A:57:SER:HB2	1:A:62:ARG:HG2	1.94	0.49
1:B:317:ILE:HD12	1:B:387:PRO:HB3	1.95	0.49
2:I:143:LYS:NZ	3:M:130:LYS:HD2	2.27	0.49
2:H:87:THR:HA	2:H:109:VAL:O	2.13	0.48
1:A:245:PRO:HA	1:A:248:ASP:O	2.12	0.48
3:L:107:LEU:HD23	3:L:141:TYR:HE1	1.78	0.48
2:H:184:VAL:HG11	2:H:194:TYR:CE1	2.47	0.48
2:I:184:VAL:HG11	2:I:194:TYR:CE1	2.48	0.48
1:A:383:LEU:O	1:A:384:ALA:HB2	2.13	0.48
2:H:11:LEU:HD11	2:H:146:PHE:HZ	1.78	0.48
2:H:30:THR:HA	2:H:52(A):THR:HG22	1.94	0.48
2:H:70:SER:OG	2:H:79:TYR:HB2	2.13	0.48
1:A:212:LEU:HD12	1:A:412:GLY:HA2	1.94	0.48
1:A:341:TYR:HE2	1:A:385:GLN:HE21	1.61	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:137:ILE:HG12	3:L:196:VAL:HG21	1.95	0.47
1:B:245:PRO:HA	1:B:248:ASP:O	2.15	0.47
3:L:133:LEU:HD12	3:L:179:LEU:HD23	1.97	0.47
1:B:81:SER:HB2	1:B:84:ARG:HD3	1.96	0.47
3:L:92:TYR:O	3:L:94:ASN:N	2.43	0.47
1:B:285:LEU:HG	1:B:316:ILE:HD13	1.97	0.47
2:I:11:LEU:HD11	2:I:146:PHE:HZ	1.80	0.46
1:B:341:TYR:HE2	1:B:385:GLN:HE21	1.63	0.46
2:H:142:VAL:HG11	2:H:150:VAL:HG11	1.97	0.46
1:A:222:VAL:HG23	1:A:309:LEU:HB3	1.97	0.46
1:B:196:ASP:HB3	1:B:266:PRO:HA	1.98	0.46
1:A:222:VAL:HG22	1:A:309:LEU:HB3	1.98	0.46
3:L:47:LEU:HA	3:L:58:THR:HG21	1.98	0.46
1:B:317:ILE:CD1	1:B:340:GLY:HA2	2.36	0.46
1:A:115:SER:HB2	1:B:70:GLU:O	2.15	0.46
1:B:253:GLU:HG3	1:B:255:SER:H	1.81	0.45
3:L:39:LYS:HG2	3:L:84:ALA:HB2	1.98	0.45
1:A:196:ASP:HB3	1:A:266:PRO:HA	1.99	0.45
1:B:222:VAL:HG23	1:B:309:LEU:HB3	1.97	0.45
1:A:253:GLU:HG3	1:A:255:SER:N	2.31	0.45
1:A:157:LYS:HB2	1:A:271:GLU:HA	1.98	0.45
1:B:199:LEU:HD21	1:B:374:ILE:HD11	1.98	0.45
1:B:125:LEU:N	1:B:126:PRO:HD2	2.32	0.44
1:B:199:LEU:CD2	1:B:374:ILE:HD11	2.46	0.44
1:A:67:ASP:OD2	1:A:70:GLU:HB2	2.17	0.44
2:I:142:VAL:HG11	2:I:150:VAL:HG11	2.00	0.44
2:I:159:LEU:HD21	2:I:182:VAL:HG21	2.00	0.44
1:A:105:ASP:HB2	1:A:144:ARG:HH11	1.83	0.44
1:A:383:LEU:CG	1:A:383:LEU:O	2.52	0.44
1:A:285:LEU:HG	1:A:316:ILE:HD13	1.98	0.44
3:L:24:ARG:HB2	3:L:70:LYS:HD3	2.00	0.44
3:M:186:TRP:O	3:M:209:PRO:HG3	2.17	0.44
1:B:253:GLU:HG3	1:B:255:SER:N	2.32	0.43
1:A:199:LEU:CD2	1:A:374:ILE:HD11	2.48	0.43
2:H:48:MET:HG2	2:H:63:PHE:CZ	2.53	0.43
1:A:142:ARG:CG	1:A:144:ARG:HG3	2.49	0.43
1:A:254:ASN:O	1:A:257:ASP:HB2	2.17	0.43
1:A:317:ILE:HD12	1:A:387:PRO:HB3	2.01	0.43
2:H:40:ALA:O	2:H:43:GLN:HB2	2.18	0.43
1:B:179:SER:HB3	1:B:218:PHE:HB3	2.01	0.43
1:B:217:VAL:HG21	1:B:263:LEU:CD2	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:217:VAL:HG21	1:B:263:LEU:HD22	2.00	0.43
3:M:133:LEU:HD12	3:M:179:LEU:HD23	2.01	0.42
2:I:44:GLY:HA3	4:M:216:HOH:O	2.19	0.42
3:M:133:LEU:O	3:M:178:TYR:HA	2.19	0.42
1:B:226:SER:HA	1:B:227:PRO:HD3	1.90	0.42
3:M:39:LYS:HG2	3:M:84:ALA:HB2	2.01	0.42
2:H:159:LEU:HD21	2:H:182:VAL:HG21	2.00	0.42
1:A:226:SER:HA	1:A:227:PRO:HD3	1.92	0.42
1:A:171:LEU:HD22	1:A:272:TYR:CD1	2.55	0.42
3:L:133:LEU:O	3:L:178:TYR:HA	2.19	0.42
2:H:56:SER:HA	2:H:57:PRO:HD3	1.97	0.42
3:M:47:LEU:HA	3:M:58:THR:HG21	2.02	0.42
1:B:331:GLN:HB2	1:B:331:GLN:HE21	1.69	0.42
3:M:34:ASN:HA	3:M:48:ILE:O	2.20	0.42
2:H:116:THR:HA	2:H:146:PHE:O	2.20	0.42
1:A:57:SER:O	1:A:61:ALA:HA	2.19	0.41
1:A:217:VAL:HG21	1:A:263:LEU:CD2	2.50	0.41
3:L:27(C):VAL:HG11	3:L:71:ALA:HB2	2.01	0.41
1:B:57:SER:O	1:B:61:ALA:HA	2.21	0.41
1:A:217:VAL:HG21	1:A:263:LEU:HD22	2.03	0.41
3:L:34:ASN:HA	3:L:48:ILE:O	2.20	0.41
3:M:152:ASP:OD2	3:M:189:HIS:CD2	2.72	0.41
3:M:146:THR:HB	3:M:197:THR:HB	2.02	0.41
1:A:199:LEU:HD21	1:A:374:ILE:HD11	2.02	0.41
3:L:4:VAL:O	3:L:99:GLY:HA2	2.22	0.40
3:M:27(C):VAL:HG11	3:M:71:ALA:HB2	2.02	0.40
2:I:116:THR:HA	2:I:146:PHE:O	2.21	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 X-ray entries followed by that with respect to entries of similar resolution.

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	329/372 (88%)	319 (97%)	10 (3%)	0	100	100
1	B	325/372 (87%)	316 (97%)	9 (3%)	0	100	100
2	H	199/225 (88%)	193 (97%)	6 (3%)	0	100	100
2	I	201/225 (89%)	194 (96%)	7 (4%)	0	100	100
3	L	209/215 (97%)	203 (97%)	6 (3%)	0	100	100
3	M	210/215 (98%)	200 (95%)	10 (5%)	0	100	100
All	All	1473/1624 (91%)	1425 (97%)	48 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	280/302 (93%)	275 (98%)	5 (2%)	66	87
1	B	276/302 (91%)	268 (97%)	8 (3%)	50	75
2	H	172/191 (90%)	168 (98%)	4 (2%)	58	82
2	I	173/191 (91%)	169 (98%)	4 (2%)	58	82
3	L	175/179 (98%)	172 (98%)	3 (2%)	68	88
3	M	176/179 (98%)	172 (98%)	4 (2%)	58	82
All	All	1252/1344 (93%)	1224 (98%)	28 (2%)	60	83

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	SER
1	A	169	THR
1	A	188	CYS
1	A	250	ASN
1	A	319	ASN
1	B	58	SER
1	B	60	ASP

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Mol	Chain	Res	Type
1	B	70	GLU
1	B	188	CYS
1	B	250	ASN
1	B	319	ASN
1	B	338	CYS
1	B	377	TRP
2	H	67	PHE
2	H	150	VAL
2	H	169	VAL
2	H	205	THR
2	I	150	VAL
2	I	169	VAL
2	I	186	SER
2	I	187	SER
3	L	21	LEU
3	L	34	ASN
3	L	104	VAL
3	M	21	LEU
3	M	34	ASN
3	M	67	LEU
3	M	104	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	250	ASN
1	A	262	HIS
1	B	256	ASN
1	B	262	HIS
2	H	1	GLN
2	H	43	GLN
2	H	199	ASN
2	I	43	GLN
2	I	199	ASN
3	L	53	ASN
3	L	79	GLN
3	L	168	GLN
3	L	189	HIS
3	M	79	GLN
3	M	168	GLN
3	M	189	HIS

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.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	339/372 (91%)	0.67	27 (7%)	15 17	25, 48, 72, 93	1 (0%)
1	B	335/372 (90%)	0.71	24 (7%)	18 21	28, 49, 72, 102	1 (0%)
2	H	203/225 (90%)	0.53	8 (3%)	43 49	25, 45, 68, 88	0
2	I	205/225 (91%)	0.50	7 (3%)	49 55	25, 44, 67, 88	0
3	L	211/215 (98%)	0.35	5 (2%)	62 67	24, 44, 61, 79	0
3	M	212/215 (98%)	0.41	12 (5%)	27 32	30, 43, 58, 64	0
All	All	1505/1624 (92%)	0.55	83 (5%)	29 33	24, 46, 69, 102	2 (0%)

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	222	VAL	6.4
1	B	123	GLY	5.5
1	A	383	LEU	4.9
1	B	239	TYR	4.1
2	H	137	ALA	3.8
1	B	59	ALA	3.8
3	M	156	VAL	3.7
1	B	230	LEU	3.7
2	I	187	SER	3.7
1	A	159	PRO	3.6
2	H	159	LEU	3.6
1	A	125	LEU	3.6
1	A	141	PRO	3.5
1	A	250	ASN	3.5
1	A	168	ASP	3.5
1	A	56	VAL	3.4
1	B	224	GLN	3.4
3	M	67	LEU	3.3
1	B	252	GLU	3.3

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Mol	Chain	Res	Type	RSRZ
2	H	195	ILE	3.3
1	A	286	VAL	3.3
3	M	203	VAL	3.3
2	I	185	PRO	3.3
1	B	148	ALA	3.1
2	H	194	TYR	3.1
1	B	95	PHE	3.1
1	A	105	ASP	2.9
2	I	136	ALA	2.9
1	B	265	SER	2.9
3	M	5	THR	2.8
1	A	218	PHE	2.8
2	H	187	SER	2.8
1	A	229	GLY	2.7
2	I	29	PHE	2.7
1	B	141	PRO	2.7
1	B	66	PHE	2.7
2	H	42	GLY	2.7
1	A	127	HIS	2.7
1	B	332	ILE	2.6
1	B	107	ARG	2.5
2	I	183	THR	2.5
1	A	129	GLN	2.5
3	M	60	ALA	2.5
1	A	60	ASP	2.5
1	A	263	LEU	2.5
1	A	167	ARG	2.4
3	L	164	THR	2.4
3	M	13	VAL	2.4
3	L	67	LEU	2.4
3	M	166	SER	2.4
3	L	181	LEU	2.4
3	L	156	VAL	2.4
2	H	82	ILE	2.4
1	B	149	ILE	2.4
1	A	259	ALA	2.4
1	A	342	PRO	2.3
1	A	308	VAL	2.3
1	A	414	VAL	2.3
1	B	144	ARG	2.3
3	M	178	TYR	2.3
1	B	169	THR	2.3

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Mol	Chain	Res	Type	RSRZ
2	I	181	VAL	2.3
1	B	139	ASP	2.2
1	B	246	PHE	2.2
1	B	341	TYR	2.2
1	A	294	THR	2.2
3	M	175	ALA	2.2
1	A	169	THR	2.2
1	B	286	VAL	2.2
2	I	194	TYR	2.2
1	A	341	TYR	2.1
1	B	105	ASP	2.1
3	M	3	VAL	2.1
1	A	246	PHE	2.1
1	A	310	GLN	2.1
2	H	77	THR	2.1
1	B	217	VAL	2.1
1	B	222	VAL	2.1
1	A	230	LEU	2.1
3	L	202	THR	2.1
3	M	123	SER	2.0
1	B	248	ASP	2.0
3	M	27(C)	VAL	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.