



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 02:46 PM EDT

PDB ID : 3C5Z / pdb_00003c5z
Title : Crystal structure of mouse MHC class II I-Ab/3K peptide complexed with mouse TCR B3K506
Authors : Dai, S.; Kappler, J.
Deposited on : 2008-02-01
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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

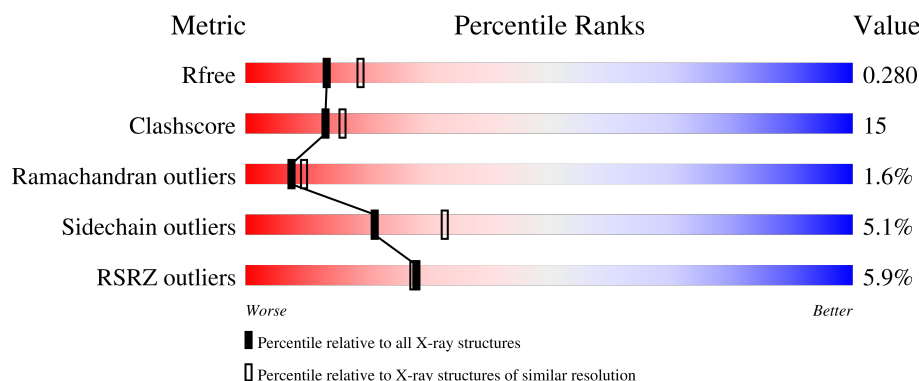
1 Overall quality at a glance

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}	180053	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	202	13% 72% 25% .
1	E	202	8% 67% 30% .
2	B	240	8% 66% 32% .
2	F	240	6% 68% 30% .
3	C	182	3% 64% 32% .

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Mol	Chain	Length	Quality of chain
3	G	182	4% 69% 28%
4	D	217	% 59% 32% 7%
4	H	217	2% 60% 29% 7%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 13288 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 TCR B3K506 Alpha Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	202	Total	C	N	O	S	0	0	0
			1581	990	261	323	7			
1	E	201	Total	C	N	O	S	0	0	0
			1573	985	260	321	7			

- Molecule 2 is a protein called TCR B3K506 Beta Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	240	Total	C	N	O	S	0	0	0
			1902	1196	330	370	6			
2	F	240	Total	C	N	O	S	0	0	0
			1902	1196	330	370	6			

- Molecule 3 is a protein called H-2 class II histocompatibility antigen, A-B alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	182	Total	C	N	O	S	0	0	0
			1459	944	230	282	3			
3	G	182	Total	C	N	O	S	0	0	0
			1459	944	230	282	3			

- Molecule 4 is a protein called 3K peptide, Linker, and H-2 class II histocompatibility antigen (A beta chain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	202	Total	C	N	O	S	0	0	0
			1673	1047	304	315	7			
4	H	202	Total	C	N	O	S	0	0	0
			1673	1047	304	315	7			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	14	GLY	-	linker	UNP P14483
D	15	GLY	-	linker	UNP P14483
D	16	GLY	-	linker	UNP P14483
D	17	GLY	-	linker	UNP P14483
D	18	SER	-	linker	UNP P14483
D	19	LEU	-	linker	UNP P14483
D	20	VAL	-	linker	UNP P14483
D	21	PRO	-	linker	UNP P14483
D	22	ARG	-	linker	UNP P14483
D	23	GLY	-	linker	UNP P14483
D	24	SER	-	linker	UNP P14483
D	25	GLY	-	linker	UNP P14483
D	26	GLY	-	linker	UNP P14483
D	27	GLY	-	linker	UNP P14483
D	28	GLY	-	linker	UNP P14483
D	216	LYS	ARG	engineered mutation	UNP P14483
H	14	GLY	-	linker	UNP P14483
H	15	GLY	-	linker	UNP P14483
H	16	GLY	-	linker	UNP P14483
H	17	GLY	-	linker	UNP P14483
H	18	SER	-	linker	UNP P14483
H	19	LEU	-	linker	UNP P14483
H	20	VAL	-	linker	UNP P14483
H	21	PRO	-	linker	UNP P14483
H	22	ARG	-	linker	UNP P14483
H	23	GLY	-	linker	UNP P14483
H	24	SER	-	linker	UNP P14483
H	25	GLY	-	linker	UNP P14483
H	26	GLY	-	linker	UNP P14483
H	27	GLY	-	linker	UNP P14483
H	28	GLY	-	linker	UNP P14483
H	216	LYS	ARG	engineered mutation	UNP P14483

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	6	Total O 6 6	0	0
5	B	6	Total O 6 6	0	0
5	C	10	Total O 10 10	0	0
5	D	10	Total O 10 10	0	0

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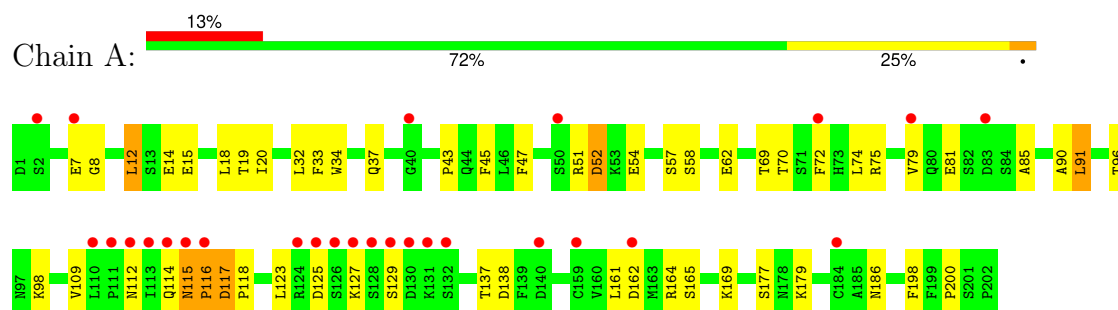
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	E	7	Total 7	O 7	0	0
5	F	4	Total 4	O 4	0	0
5	G	8	Total 8	O 8	0	0
5	H	15	Total 15	O 15	0	0

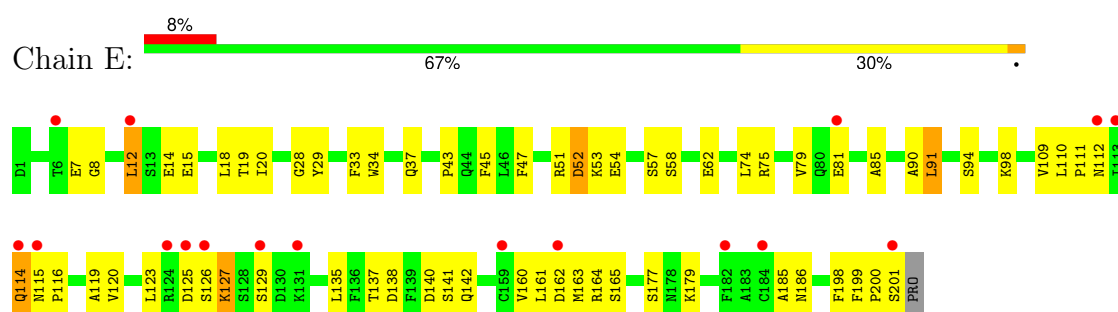
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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 and electron density. 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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

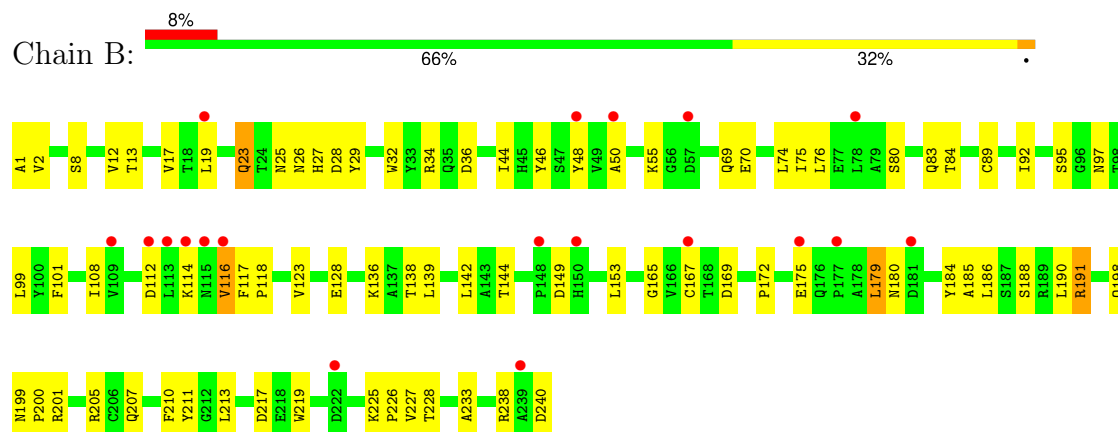
• Molecule 1: TCR B3K506 Alpha Chain



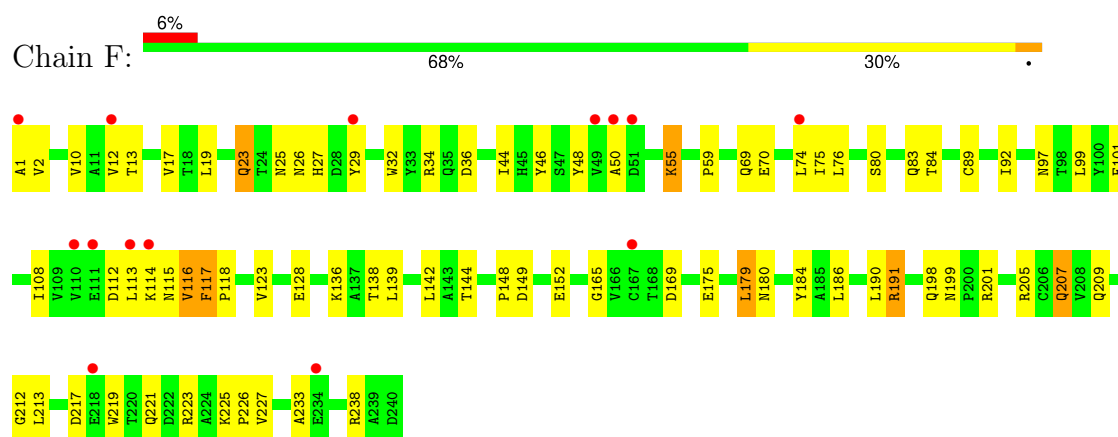
• Molecule 1: TCR B3K506 Alpha Chain



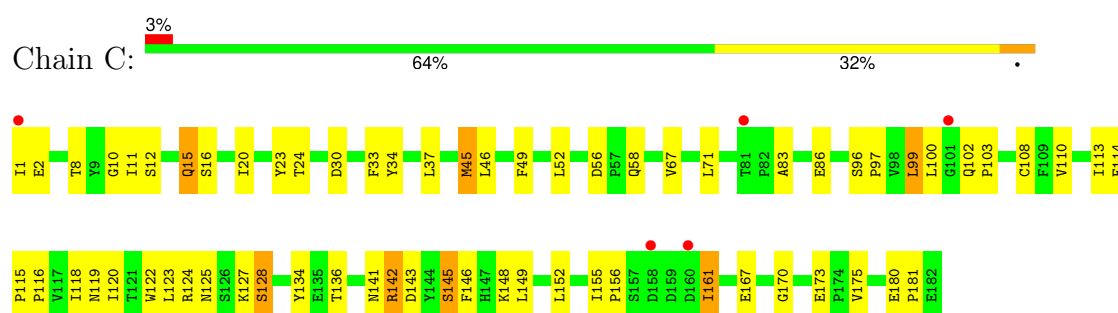
• Molecule 2: TCR B3K506 Beta Chain



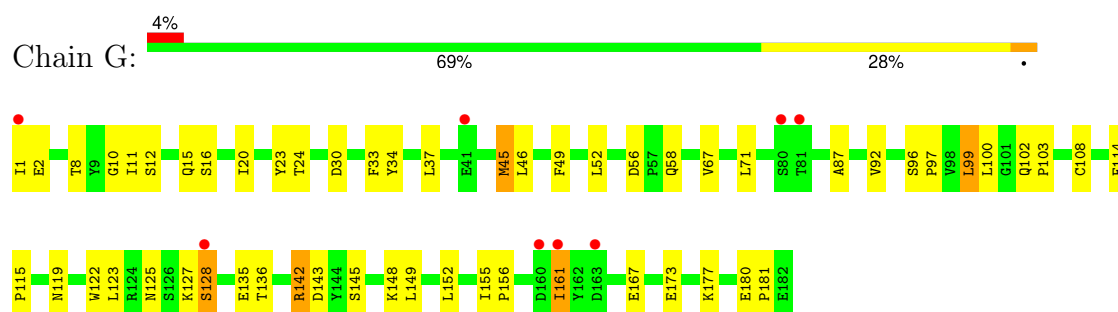
• Molecule 2: TCR B3K506 Beta Chain



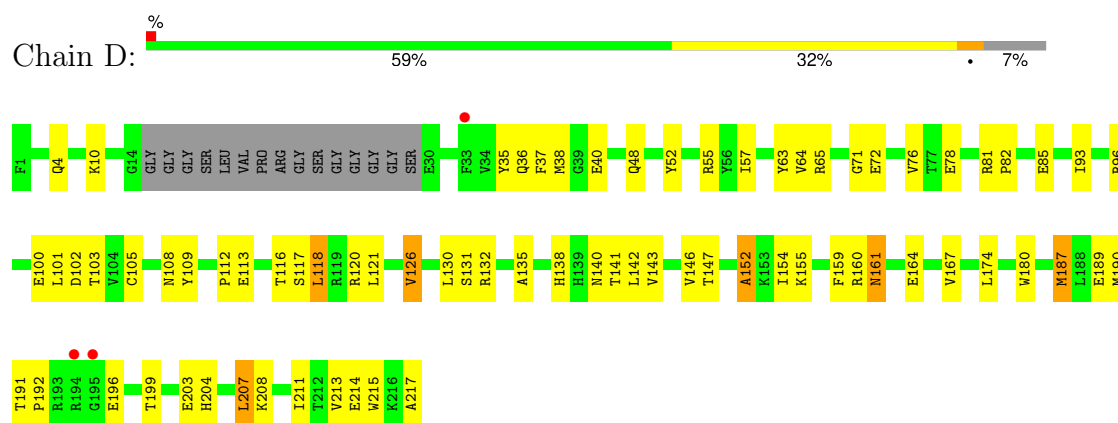
- Molecule 3: H-2 class II histocompatibility antigen, A-B alpha chain



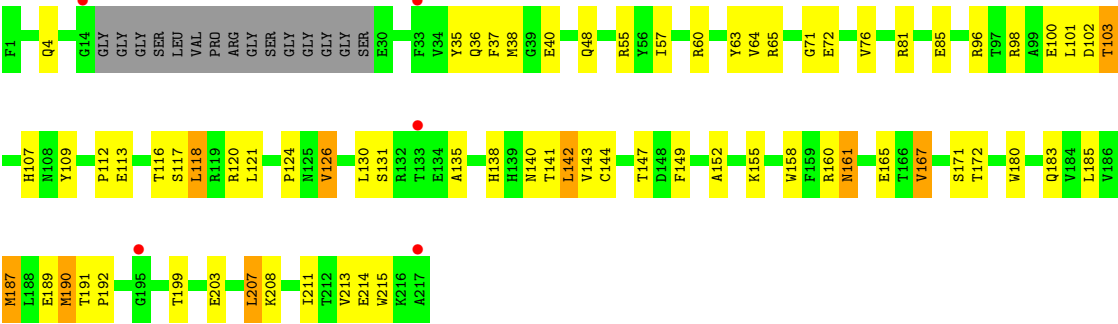
- Molecule 3: H-2 class II histocompatibility antigen, A-B alpha chain



- Molecule 4: 3K peptide, Linker, and H-2 class II histocompatibility antigen (A beta chain)



- Molecule 4: 3K peptide, Linker, and H-2 class II histocompatibility antigen (A beta chain)



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	65.30Å 117.80Å 136.20Å 90.00° 99.56° 90.00°	Depositor
Resolution (Å)	40.00 – 2.55 40.00 – 2.55	Depositor EDS
% Data completeness (in resolution range)	86.0 (40.00-2.55) 85.9 (40.00-2.55)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.54Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.246 , 0.281 0.246 , 0.280	Depositor DCC
R_{free} test set	2886 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	49.2	Xtriage
Anisotropy	0.821	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 39.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	13288	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% 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, 0.333 respectively for untwinned datasets, and 0.375, 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	0.51	0/1618	0.92	2/2195 (0.1%)
1	E	0.49	0/1609	0.98	3/2183 (0.1%)
2	B	0.54	0/1953	0.90	2/2660 (0.1%)
2	F	0.52	0/1953	0.91	2/2660 (0.1%)
3	C	0.58	0/1504	0.91	3/2054 (0.1%)
3	G	0.60	0/1504	0.92	5/2054 (0.2%)
4	D	0.61	1/1713 (0.1%)	0.99	5/2321 (0.2%)
4	H	0.64	1/1713 (0.1%)	0.99	5/2321 (0.2%)
All	All	0.56	2/13567 (0.0%)	0.94	27/18448 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	187	MET	SD-CE	6.30	1.95	1.79
4	D	187	MET	SD-CE	5.06	1.92	1.79

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	115	ASN	CA-C-N	13.00	136.90	120.51
1	E	115	ASN	C-N-CA	13.00	136.90	120.51
2	F	117	PHE	CA-C-N	6.78	124.61	119.66
2	F	117	PHE	C-N-CA	6.78	124.61	119.66
3	C	56	ASP	CA-C-N	6.36	126.85	119.47
3	C	56	ASP	C-N-CA	6.36	126.85	119.47
2	B	116	VAL	N-CA-C	6.23	117.69	109.21
1	A	115	ASN	CA-C-N	-6.11	112.20	119.84
1	A	115	ASN	C-N-CA	-6.11	112.20	119.84
3	C	145	SER	N-CA-C	-6.08	101.21	110.14
4	D	57	ILE	N-CA-C	5.95	116.44	108.11
4	D	152	ALA	N-CA-C	5.93	117.74	111.28
3	G	56	ASP	CA-C-N	5.83	126.23	119.47
3	G	56	ASP	C-N-CA	5.83	126.23	119.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	57	ILE	N-CA-C	5.77	116.18	107.75
4	H	152	ALA	N-CA-C	5.76	118.30	111.33
4	D	147	THR	N-CA-C	5.74	117.70	109.14
4	D	76	VAL	N-CA-C	-5.69	106.77	113.42
3	G	145	SER	N-CA-C	-5.69	100.98	109.96
4	H	208	LYS	N-CA-C	-5.68	106.19	113.23
2	B	116	VAL	CB-CA-C	-5.61	103.38	111.34
4	D	208	LYS	N-CA-C	-5.59	105.96	112.89
4	H	76	VAL	N-CA-C	-5.42	107.08	113.42
4	H	147	THR	N-CA-C	5.34	117.79	109.52
3	G	87	ALA	CA-C-N	5.08	125.47	119.99
3	G	87	ALA	C-N-CA	5.08	125.47	119.99
1	E	94	SER	N-CA-C	5.01	117.48	111.71

There are no chirality outliers.

There are no planarity outliers.

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	1581	0	1483	38	0
1	E	1573	0	1476	40	0
2	B	1902	0	1802	58	0
2	F	1902	0	1802	64	0
3	C	1459	0	1386	50	0
3	G	1459	0	1386	40	0
4	D	1673	0	1603	63	0
4	H	1673	0	1603	58	0
5	A	6	0	0	0	0
5	B	6	0	0	0	0
5	C	10	0	0	1	0
5	D	10	0	0	2	0
5	E	7	0	0	0	0
5	F	4	0	0	0	0
5	G	8	0	0	1	0
5	H	15	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	13288	0	12541	384	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 15.

All (384) 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)
4:D:112:PRO:O	4:D:116:THR:HG22	1.64	0.98
4:H:112:PRO:O	4:H:116:THR:HG22	1.65	0.94
4:D:199:THR:HG22	4:D:214:GLU:HG3	1.53	0.90
4:H:116:THR:HG23	4:H:117:SER:H	1.37	0.90
4:H:199:THR:HG22	4:H:214:GLU:HG3	1.53	0.89
4:D:116:THR:HG23	4:D:117:SER:H	1.38	0.89
2:B:179:LEU:HD23	2:B:180:ASN:H	1.48	0.79
3:C:161:ILE:HD13	3:C:161:ILE:H	1.47	0.77
3:G:161:ILE:HD13	3:G:161:ILE:H	1.46	0.77
1:E:114:GLN:O	1:E:116:PRO:HD3	1.86	0.75
3:G:161:ILE:HG22	3:G:180:GLU:HG3	1.69	0.74
3:G:102:GLN:NE2	3:G:103:PRO:HD2	2.04	0.72
2:B:28:ASP:OD1	4:D:10:LYS:HD3	1.89	0.72
2:B:19:LEU:HD12	2:B:74:LEU:HD23	1.72	0.71
3:C:96:SER:HB2	3:C:97:PRO:HD2	1.73	0.70
2:B:138:THR:OG1	2:B:191:ARG:HD3	1.91	0.70
2:F:19:LEU:HD12	2:F:74:LEU:HD23	1.73	0.70
3:G:96:SER:HB2	3:G:97:PRO:HD2	1.73	0.69
2:B:198:GLN:HG2	2:B:240:ASP:HB2	1.74	0.69
2:F:115:ASN:O	2:F:116:VAL:C	2.31	0.69
3:G:100:LEU:HD12	3:G:100:LEU:H	1.57	0.69
4:D:135:ALA:O	4:D:138:HIS:HB2	1.93	0.68
3:C:142:ARG:HH11	3:C:142:ARG:HB2	1.57	0.68
2:F:17:VAL:HG22	2:F:76:LEU:HB2	1.76	0.67
2:B:213:LEU:HD13	2:B:226:PRO:HG2	1.77	0.67
1:E:119:ALA:HB2	1:E:198:PHE:HB3	1.75	0.66
4:D:143:VAL:HG22	4:D:187:MET:HG3	1.77	0.66
2:B:17:VAL:HG22	2:B:76:LEU:HB2	1.77	0.66
2:F:205:ARG:HH11	2:F:207:GLN:HG3	1.60	0.66
4:H:100:GLU:HA	4:H:103:THR:HB	1.77	0.66
2:F:179:LEU:HD23	2:F:180:ASN:H	1.60	0.65
4:H:140:ASN:HB3	4:H:190:MET:HE1	1.78	0.65
2:B:198:GLN:HA	2:B:238:ARG:O	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:219:TRP:HB2	2:B:225:LYS:HD2	1.78	0.65
4:D:100:GLU:HA	4:D:103:THR:HB	1.78	0.65
4:D:140:ASN:HB3	4:D:190:MET:HE1	1.77	0.65
4:H:135:ALA:O	4:H:138:HIS:HB2	1.96	0.65
1:E:51:ARG:NH2	4:H:98:ARG:HG2	2.13	0.64
3:C:100:LEU:HD12	3:C:100:LEU:H	1.62	0.64
2:F:198:GLN:HA	2:F:238:ARG:O	1.98	0.64
4:H:126:VAL:HG22	4:H:213:VAL:HG21	1.80	0.64
2:B:205:ARG:HH11	2:B:207:GLN:HG3	1.63	0.64
4:D:63:TYR:CD2	4:D:64:VAL:HG23	2.33	0.64
1:A:125:ASP:HB3	1:A:129:SER:HA	1.79	0.63
4:D:164:GLU:O	4:H:155:LYS:HE2	1.98	0.63
3:C:142:ARG:HB2	3:C:142:ARG:NH1	2.13	0.63
3:C:161:ILE:HG22	3:C:180:GLU:HG3	1.81	0.63
4:H:63:TYR:CD2	4:H:64:VAL:HG23	2.34	0.62
4:H:120:ARG:O	4:H:121:LEU:HD12	1.99	0.62
3:C:10:GLY:HA3	5:D:221:HOH:O	1.98	0.62
4:D:36:GLN:NE2	5:D:219:HOH:O	2.33	0.62
2:B:219:TRP:CB	2:B:225:LYS:HD2	2.30	0.61
4:D:112:PRO:C	4:D:116:THR:HG22	2.26	0.61
1:E:111:PRO:HG3	1:E:160:VAL:HG13	1.81	0.60
1:E:114:GLN:C	1:E:116:PRO:HD3	2.26	0.60
1:A:198:PHE:CE2	1:A:200:PRO:HG3	2.36	0.60
4:D:155:LYS:HB2	4:D:203:GLU:HG3	1.84	0.60
3:C:102:GLN:H	3:C:155:ILE:HG23	1.66	0.60
3:G:142:ARG:HH11	3:G:142:ARG:HB2	1.65	0.60
1:E:161:LEU:C	1:E:161:LEU:HD23	2.26	0.59
3:G:142:ARG:HB2	3:G:142:ARG:NH1	2.16	0.59
1:A:115:ASN:CG	1:A:115:ASN:O	2.45	0.59
2:B:169:ASP:HB2	2:B:186:LEU:HD12	1.84	0.59
1:E:119:ALA:CB	1:E:198:PHE:HB3	2.32	0.59
4:H:112:PRO:C	4:H:116:THR:HG22	2.27	0.59
4:D:138:HIS:O	4:D:192:PRO:HD2	2.02	0.59
2:F:213:LEU:HD13	2:F:226:PRO:HG2	1.85	0.59
4:H:101:LEU:HD23	4:H:102:ASP:OD2	2.02	0.59
3:C:136:THR:O	3:C:148:LYS:NZ	2.36	0.59
4:D:101:LEU:HD23	4:D:102:ASP:OD2	2.02	0.59
3:G:45:MET:HE3	5:G:186:HOH:O	2.02	0.59
1:E:20:ILE:HD12	1:E:74:LEU:HD23	1.84	0.59
3:G:180:GLU:HG2	3:G:181:PRO:HD2	1.85	0.58
4:D:120:ARG:O	4:D:121:LEU:HD12	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:140:ASN:HB3	4:H:190:MET:CE	2.34	0.58
1:A:20:ILE:HD12	1:A:74:LEU:HD23	1.85	0.58
3:C:102:GLN:O	3:C:156:PRO:HD2	2.03	0.58
2:F:169:ASP:HB2	2:F:186:LEU:CD1	2.32	0.58
3:G:71:LEU:HD13	4:H:35:TYR:HB2	1.85	0.58
4:H:81:ARG:HB3	4:H:81:ARG:NH1	2.19	0.58
2:B:92:ILE:HD11	2:B:99:LEU:HG	1.86	0.58
4:D:131:SER:OG	4:D:141:THR:HB	2.04	0.58
4:H:155:LYS:HB2	4:H:203:GLU:HG3	1.84	0.58
3:C:71:LEU:HD13	4:D:35:TYR:HB2	1.84	0.58
4:D:81:ARG:NH1	4:D:81:ARG:HB3	2.18	0.58
2:F:123:VAL:HG23	2:F:233:ALA:HB3	1.85	0.58
1:E:125:ASP:HB3	1:E:129:SER:HA	1.87	0.57
2:F:92:ILE:HD11	2:F:99:LEU:HG	1.86	0.57
1:A:81:GLU:HA	1:A:109:VAL:HB	1.86	0.57
4:D:126:VAL:HG22	4:D:213:VAL:HG21	1.87	0.57
4:D:140:ASN:HB3	4:D:190:MET:CE	2.34	0.57
2:F:169:ASP:HB2	2:F:186:LEU:HD12	1.86	0.57
3:G:135:GLU:CG	3:G:148:LYS:HD3	2.34	0.56
3:C:123:LEU:HA	3:C:128:SER:HA	1.86	0.56
2:F:205:ARG:NH1	2:F:207:GLN:HG3	2.20	0.56
1:E:28:GLY:HA2	4:H:107:HIS:CE1	2.39	0.56
1:E:81:GLU:HA	1:E:109:VAL:HB	1.87	0.56
4:H:207:LEU:HG	4:H:211:ILE:HG13	1.87	0.56
1:A:98:LYS:HE3	2:B:46:TYR:CE2	2.41	0.56
2:B:169:ASP:HB2	2:B:186:LEU:CD1	2.36	0.56
3:C:141:ASN:HB2	3:C:145:SER:O	2.06	0.55
1:E:111:PRO:HG3	1:E:160:VAL:CG1	2.36	0.55
2:F:139:LEU:N	2:F:139:LEU:HD12	2.22	0.55
4:D:130:LEU:HD13	4:D:215:TRP:CZ2	2.42	0.55
4:D:138:HIS:C	4:D:191:THR:HG23	2.32	0.55
1:A:161:LEU:HD23	1:A:161:LEU:C	2.32	0.55
1:E:98:LYS:HE3	2:F:46:TYR:CE2	2.41	0.55
3:C:125:ASN:OD1	3:C:161:ILE:HD13	2.07	0.54
1:E:91:LEU:HD23	1:E:91:LEU:C	2.33	0.54
2:F:108:ILE:HD13	2:F:148:PRO:CB	2.37	0.54
1:A:198:PHE:CZ	1:A:200:PRO:HG3	2.42	0.54
1:E:37:GLN:O	1:E:85:ALA:HB1	2.07	0.54
2:B:34:ARG:HB3	2:B:44:ILE:HD11	1.90	0.54
2:F:34:ARG:HB3	2:F:44:ILE:HD11	1.90	0.54
4:H:155:LYS:HB2	4:H:203:GLU:CG	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:67:VAL:HG13	4:H:35:TYR:CD2	2.43	0.54
4:H:36:GLN:NE2	5:H:222:HOH:O	2.40	0.54
4:H:116:THR:HG23	4:H:117:SER:N	2.15	0.54
4:H:143:VAL:HG22	4:H:187:MET:HG3	1.88	0.54
2:B:95:SER:O	4:D:96:ARG:NH1	2.41	0.53
2:B:123:VAL:HG23	2:B:233:ALA:HB3	1.88	0.53
1:A:37:GLN:O	1:A:85:ALA:HB1	2.09	0.53
3:G:119:ASN:HB2	3:G:167:GLU:HB3	1.91	0.53
2:B:179:LEU:HD23	2:B:180:ASN:N	2.22	0.53
1:A:96:THR:HG22	3:C:58:GLN:HB3	1.89	0.53
2:F:114:LYS:HA	2:F:219:TRP:CZ3	2.43	0.53
3:C:102:GLN:NE2	3:C:103:PRO:HD2	2.24	0.53
2:F:138:THR:OG1	2:F:191:ARG:HD3	2.09	0.53
3:C:30:ASP:HB3	4:D:180:TRP:CE2	2.44	0.53
2:F:219:TRP:HB2	2:F:225:LYS:HD2	1.91	0.53
2:F:92:ILE:HG13	2:F:99:LEU:HD23	1.92	0.52
2:F:108:ILE:HD13	2:F:148:PRO:HB2	1.90	0.52
1:E:137:THR:O	1:E:138:ASP:HB2	2.10	0.52
2:F:219:TRP:CB	2:F:225:LYS:HD2	2.39	0.52
4:D:207:LEU:HG	4:D:211:ILE:HG13	1.91	0.52
1:A:91:LEU:C	1:A:91:LEU:HD23	2.34	0.52
3:C:83:ALA:HB1	3:C:114:PHE:HE1	1.74	0.52
2:B:205:ARG:NH1	2:B:207:GLN:HG3	2.24	0.52
2:B:217:ASP:O	2:B:225:LYS:HE3	2.10	0.52
4:D:161:ASN:HD22	4:D:196:GLU:HA	1.74	0.52
2:B:1:ALA:HB1	2:B:25:ASN:HD21	1.75	0.52
2:B:92:ILE:HG13	2:B:99:LEU:HD23	1.92	0.52
3:C:67:VAL:HG13	4:D:35:TYR:CD2	2.44	0.52
1:E:43:PRO:HG2	2:F:101:PHE:CD1	2.45	0.52
2:F:113:LEU:HB3	2:F:213:LEU:HD21	1.92	0.52
2:F:115:ASN:O	2:F:117:PHE:CD2	2.63	0.52
2:B:165:GLY:O	2:B:190:LEU:HA	2.10	0.51
3:C:108:CYS:HB2	3:C:122:TRP:CH2	2.45	0.51
3:G:102:GLN:H	3:G:155:ILE:HG23	1.75	0.51
3:G:148:LYS:HG2	3:G:149:LEU:N	2.25	0.51
4:H:199:THR:HG23	5:H:229:HOH:O	2.11	0.51
3:G:127:LYS:O	3:G:128:SER:HB3	2.11	0.51
3:G:100:LEU:HD12	3:G:100:LEU:N	2.25	0.51
2:F:165:GLY:O	2:F:190:LEU:HA	2.10	0.51
3:G:114:PHE:HA	3:G:115:PRO:C	2.35	0.51
4:D:116:THR:HG23	4:D:117:SER:N	2.17	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:225:LYS:HG2	2:F:227:VAL:HG13	1.93	0.51
3:C:33:PHE:CD1	3:C:33:PHE:C	2.90	0.50
4:D:138:HIS:O	4:D:191:THR:HG23	2.11	0.50
1:E:177:SER:OG	1:E:179:LYS:HG2	2.11	0.50
1:A:161:LEU:HD23	1:A:162:ASP:N	2.27	0.50
2:B:225:LYS:HG2	2:B:227:VAL:HG13	1.94	0.50
4:H:138:HIS:O	4:H:192:PRO:HD2	2.11	0.50
2:B:139:LEU:HD12	2:B:139:LEU:N	2.26	0.50
4:H:130:LEU:HD13	4:H:215:TRP:CZ2	2.45	0.50
4:D:207:LEU:HD11	4:D:211:ILE:HD11	1.94	0.49
1:A:43:PRO:HG2	2:B:101:PHE:CD1	2.47	0.49
2:B:95:SER:HB2	4:D:93:ILE:CD1	2.42	0.49
3:G:99:LEU:O	3:G:102:GLN:HB3	2.13	0.49
4:H:207:LEU:HD11	4:H:211:ILE:HD11	1.93	0.49
3:G:33:PHE:CD1	3:G:33:PHE:C	2.90	0.49
2:B:165:GLY:O	2:B:190:LEU:HD12	2.12	0.49
2:F:1:ALA:HB1	2:F:25:ASN:HD21	1.76	0.49
1:A:116:PRO:HB2	1:A:118:PRO:HD3	1.93	0.49
4:H:113:GLU:HA	4:H:116:THR:CG2	2.43	0.49
1:A:74:LEU:O	1:A:75:ARG:HG3	2.13	0.48
3:G:100:LEU:H	3:G:100:LEU:CD1	2.24	0.48
2:F:144:THR:HB	2:F:175:GLU:OE1	2.13	0.48
4:D:113:GLU:HA	4:D:116:THR:CG2	2.43	0.48
2:F:113:LEU:O	2:F:219:TRP:HZ3	1.96	0.48
3:C:142:ARG:HH11	3:C:142:ARG:CB	2.26	0.48
1:E:123:LEU:N	1:E:123:LEU:HD12	2.28	0.48
4:H:81:ARG:HB3	4:H:81:ARG:HH11	1.78	0.48
3:G:67:VAL:HG13	4:H:35:TYR:HD2	1.79	0.47
3:C:86:GLU:O	3:C:170:GLY:HA3	2.14	0.47
4:D:48:GLN:HA	4:D:48:GLN:NE2	2.30	0.47
4:D:81:ARG:HB3	4:D:81:ARG:HH11	1.78	0.47
4:D:48:GLN:HA	4:D:48:GLN:HE21	1.78	0.47
3:G:123:LEU:HA	3:G:128:SER:HA	1.96	0.47
4:H:138:HIS:C	4:H:191:THR:HG23	2.40	0.47
1:E:14:GLU:O	1:E:15:GLU:HB2	2.14	0.47
4:H:64:VAL:HG12	4:H:65:ARG:N	2.30	0.47
2:B:144:THR:HB	2:B:175:GLU:OE1	2.15	0.47
3:C:100:LEU:HD12	3:C:100:LEU:N	2.28	0.47
4:D:64:VAL:HG12	4:D:65:ARG:N	2.30	0.47
4:D:155:LYS:HB2	4:D:203:GLU:CG	2.43	0.47
1:E:126:SER:O	1:E:127:LYS:HB2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:149:ASP:HB3	2:B:184:TYR:CE2	2.49	0.47
1:E:29:TYR:HB3	4:H:103:THR:HA	1.97	0.47
2:F:32:TRP:CH2	2:F:89:CYS:HB2	2.50	0.47
3:G:8:THR:HG22	3:G:11:ILE:HG12	1.97	0.47
4:H:130:LEU:HD12	4:H:142:LEU:CD1	2.45	0.47
3:C:175:VAL:O	3:C:175:VAL:HG23	2.15	0.47
1:A:14:GLU:O	1:A:15:GLU:HB2	2.14	0.46
2:B:210:PHE:O	2:B:228:THR:HG23	2.15	0.46
4:D:40:GLU:CD	4:D:55:ARG:HE	2.23	0.46
2:F:80:SER:H	2:F:83:GLN:HE21	1.63	0.46
1:A:115:ASN:O	1:A:115:ASN:ND2	2.48	0.46
2:F:149:ASP:HB3	2:F:184:TYR:CD2	2.50	0.46
2:B:118:PRO:HD3	2:B:226:PRO:HB3	1.97	0.46
2:B:149:ASP:HB3	2:B:184:TYR:CD2	2.50	0.46
2:F:149:ASP:HB3	2:F:184:TYR:CE2	2.50	0.46
4:H:48:GLN:HE21	4:H:48:GLN:HA	1.81	0.46
2:B:199:ASN:ND2	2:B:201:ARG:HB2	2.31	0.46
1:E:74:LEU:O	1:E:75:ARG:HG3	2.16	0.46
4:H:130:LEU:HD12	4:H:142:LEU:HD13	1.97	0.46
1:A:137:THR:O	1:A:138:ASP:HB2	2.15	0.46
2:B:70:GLU:N	2:B:70:GLU:OE2	2.48	0.46
2:B:144:THR:HG22	2:B:185:ALA:HB1	1.98	0.46
1:E:33:PHE:HB2	1:E:90:ALA:HB3	1.97	0.46
2:F:112:ASP:C	2:F:114:LYS:N	2.72	0.46
4:H:144:CYS:HB2	4:H:158:TRP:CZ2	2.50	0.46
3:C:99:LEU:O	3:C:102:GLN:HB3	2.16	0.46
3:G:142:ARG:HH11	3:G:142:ARG:CB	2.29	0.46
3:C:119:ASN:HB2	3:C:167:GLU:HB3	1.97	0.46
1:E:110:LEU:HB3	1:E:141:SER:HB3	1.98	0.46
2:F:114:LYS:HG3	2:F:221:GLN:CG	2.45	0.46
4:H:37:PHE:CG	4:H:38:MET:N	2.84	0.46
2:B:23:GLN:OE1	2:B:27:HIS:N	2.49	0.46
2:F:136:LYS:HB3	2:F:191:ARG:HD2	1.97	0.46
3:G:136:THR:O	3:G:148:LYS:NZ	2.46	0.46
2:B:75:ILE:N	2:B:75:ILE:HD12	2.31	0.45
3:C:110:VAL:HG21	3:C:120:ILE:CD1	2.46	0.45
2:F:23:GLN:OE1	2:F:27:HIS:HB2	2.17	0.45
2:F:10:VAL:HG21	2:F:212:GLY:HA2	1.98	0.45
3:C:8:THR:HG22	3:C:11:ILE:HG12	1.99	0.45
4:D:37:PHE:CG	4:D:38:MET:N	2.84	0.45
4:H:124:PRO:HB3	4:H:149:PHE:HB3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:183:GLN:HB3	5:H:228:HOH:O	2.17	0.45
3:C:67:VAL:HG13	4:D:35:TYR:HD2	1.80	0.45
2:F:179:LEU:HD23	2:F:180:ASN:N	2.29	0.45
4:D:52:TYR:CD1	4:D:105:CYS:SG	3.09	0.45
2:F:118:PRO:HD3	2:F:226:PRO:HB3	1.98	0.45
2:F:217:ASP:O	2:F:225:LYS:HE3	2.17	0.45
4:D:96:ARG:HB3	4:D:96:ARG:CZ	2.46	0.45
3:C:146:PHE:CD1	3:C:146:PHE:N	2.85	0.45
3:G:114:PHE:CZ	4:H:60:ARG:HG3	2.52	0.45
1:A:12:LEU:HD13	1:A:79:VAL:HG11	1.99	0.45
2:F:23:GLN:OE1	2:F:27:HIS:N	2.48	0.45
2:F:80:SER:HB3	2:F:83:GLN:HG3	1.99	0.45
3:C:15:GLN:OE1	3:C:116:PRO:HG2	2.17	0.45
3:C:113:ILE:HA	5:C:185:HOH:O	2.15	0.45
1:E:12:LEU:HD13	1:E:79:VAL:HG11	1.99	0.45
2:F:114:LYS:HG3	2:F:221:GLN:HG3	1.98	0.45
3:C:100:LEU:H	3:C:100:LEU:CD1	2.27	0.45
4:D:160:ARG:O	4:D:161:ASN:C	2.59	0.45
2:F:23:GLN:O	2:F:70:GLU:HB2	2.17	0.45
2:F:75:ILE:HD12	2:F:75:ILE:N	2.31	0.45
2:B:153:LEU:HD13	2:B:188:SER:HB2	1.98	0.44
2:F:70:GLU:OE2	2:F:70:GLU:N	2.49	0.44
1:A:18:LEU:HD22	1:A:19:THR:H	1.82	0.44
1:A:33:PHE:HB2	1:A:90:ALA:HB3	1.98	0.44
2:B:80:SER:H	2:B:83:GLN:HE21	1.65	0.44
2:F:108:ILE:HG21	2:F:148:PRO:HB3	1.99	0.44
4:D:130:LEU:HD12	4:D:142:LEU:HD13	2.00	0.44
1:E:34:TRP:CZ2	1:E:74:LEU:HB2	2.52	0.44
2:F:48:TYR:CE1	3:G:58:GLN:NE2	2.85	0.44
3:C:180:GLU:HG2	3:C:181:PRO:HD2	1.98	0.44
2:F:114:LYS:O	2:F:223:ARG:NH2	2.43	0.44
1:E:18:LEU:HD22	1:E:19:THR:H	1.82	0.44
1:E:45:PHE:CZ	1:E:47:PHE:HA	2.52	0.44
1:A:18:LEU:HD22	1:A:19:THR:N	2.33	0.44
2:F:152:GLU:HB2	2:F:209:GLN:HB3	1.99	0.44
4:H:48:GLN:HA	4:H:48:GLN:NE2	2.32	0.44
4:D:146:VAL:HG11	4:D:154:ILE:HD11	2.00	0.44
2:B:23:GLN:OE1	2:B:27:HIS:HB2	2.18	0.44
2:B:23:GLN:O	2:B:70:GLU:HB2	2.18	0.44
1:E:120:VAL:HA	1:E:135:LEU:O	2.18	0.44
2:B:32:TRP:CH2	2:B:89:CYS:HB2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:148:LYS:HG2	3:C:149:LEU:N	2.33	0.43
4:D:132:ARG:H	4:D:132:ARG:HG2	1.57	0.43
4:D:215:TRP:CZ3	4:D:217:ALA:HA	2.52	0.43
1:E:18:LEU:HD22	1:E:19:THR:N	2.33	0.43
3:G:125:ASN:OD1	3:G:161:ILE:HD13	2.18	0.43
1:A:117:ASP:N	1:A:118:PRO:CD	2.81	0.43
2:B:112:ASP:C	2:B:114:LYS:N	2.74	0.43
1:E:161:LEU:HD23	1:E:162:ASP:N	2.32	0.43
1:E:164:ARG:HA	1:E:164:ARG:HD2	1.77	0.43
3:C:15:GLN:HE22	3:C:115:PRO:HB2	1.83	0.43
3:C:114:PHE:HA	3:C:115:PRO:C	2.43	0.43
1:A:34:TRP:CZ2	1:A:74:LEU:HB2	2.54	0.43
3:G:10:GLY:HA2	3:G:23:TYR:CZ	2.54	0.43
1:A:45:PHE:CZ	1:A:47:PHE:HA	2.53	0.43
1:A:52:ASP:C	1:A:54:GLU:H	2.26	0.43
2:B:136:LYS:HB3	2:B:191:ARG:HD2	2.01	0.43
1:E:29:TYR:CE1	4:H:102:ASP:HB3	2.54	0.43
3:G:46:LEU:O	3:G:49:PHE:HB2	2.18	0.43
4:H:40:GLU:CD	4:H:55:ARG:HE	2.25	0.43
1:E:52:ASP:C	1:E:54:GLU:H	2.27	0.43
1:A:117:ASP:O	1:A:117:ASP:CG	2.62	0.43
2:B:8:SER:HB3	2:B:211:TYR:CD2	2.54	0.43
3:C:118:ILE:HG12	3:C:119:ASN:N	2.34	0.43
3:C:127:LYS:O	3:C:128:SER:HB3	2.18	0.43
4:H:96:ARG:HB3	4:H:96:ARG:CZ	2.48	0.43
2:B:95:SER:CB	4:D:93:ILE:HG12	2.48	0.43
3:C:10:GLY:HA2	3:C:23:TYR:CZ	2.54	0.42
1:A:164:ARG:HA	1:A:164:ARG:HD2	1.76	0.42
2:B:207:GLN:HE21	2:B:207:GLN:HB3	1.65	0.42
3:C:12:SER:HB2	3:C:67:VAL:HG21	2.02	0.42
2:F:29:TYR:CE2	2:F:48:TYR:HB3	2.54	0.42
2:F:112:ASP:C	2:F:114:LYS:H	2.26	0.42
4:H:172:THR:HG21	4:H:185:LEU:HB2	2.01	0.42
3:G:108:CYS:HB2	3:G:122:TRP:CH2	2.54	0.42
4:H:131:SER:OG	4:H:141:THR:HB	2.20	0.42
1:A:57:SER:O	1:A:58:SER:HB2	2.19	0.42
3:C:155:ILE:HD12	3:C:155:ILE:N	2.34	0.42
2:B:84:THR:HG23	2:B:108:ILE:HA	2.02	0.42
1:E:47:PHE:CD1	1:E:47:PHE:C	2.98	0.42
4:D:109:TYR:CZ	4:D:118:LEU:HD21	2.55	0.42
2:F:128:GLU:H	2:F:128:GLU:HG2	1.61	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:29:TYR:CE2	2:B:48:TYR:HB3	2.55	0.42
2:B:95:SER:HB3	4:D:93:ILE:HG23	2.01	0.42
3:C:24:THR:HG22	3:C:34:TYR:HB3	2.01	0.42
3:G:135:GLU:HG3	3:G:148:LYS:HD3	2.02	0.42
1:A:117:ASP:H	1:A:118:PRO:CD	2.33	0.42
1:A:177:SER:OG	1:A:179:LYS:HG2	2.19	0.42
2:B:116:VAL:HG12	2:B:117:PHE:N	2.35	0.42
3:C:123:LEU:HD23	3:C:128:SER:HA	2.01	0.42
2:F:80:SER:N	2:F:83:GLN:HE21	2.17	0.42
2:F:169:ASP:HB2	2:F:186:LEU:HD11	2.02	0.42
3:G:37:LEU:H	3:G:37:LEU:HD22	1.85	0.42
4:H:71:GLY:O	4:H:72:GLU:HB3	2.20	0.42
4:D:78:GLU:OE2	4:D:81:ARG:HD2	2.20	0.41
4:D:204:HIS:HB3	4:D:207:LEU:CD2	2.50	0.41
4:H:109:TYR:CZ	4:H:118:LEU:HD21	2.54	0.41
2:B:149:ASP:HB3	2:B:184:TYR:CZ	2.55	0.41
3:C:46:LEU:O	3:C:49:PHE:HB2	2.19	0.41
1:E:140:ASP:OD2	1:E:142:GLN:HB2	2.21	0.41
2:F:114:LYS:HA	2:F:219:TRP:CH2	2.55	0.41
2:B:26:ASN:HA	2:B:69:GLN:HE22	1.85	0.41
3:C:161:ILE:H	3:C:161:ILE:CD1	2.25	0.41
4:D:130:LEU:HD12	4:D:142:LEU:CD1	2.50	0.41
2:F:26:ASN:HA	2:F:69:GLN:HE22	1.85	0.41
4:H:81:ARG:O	4:H:85:GLU:HG2	2.19	0.41
4:D:81:ARG:O	4:D:85:GLU:HG2	2.20	0.41
3:G:24:THR:HG22	3:G:34:TYR:HB3	2.02	0.41
1:A:81:GLU:HG2	1:A:169:LYS:HE3	2.02	0.41
4:D:38:MET:HE2	4:D:40:GLU:OE1	2.20	0.41
3:G:92:VAL:HG23	3:G:177:LYS:HG2	2.02	0.41
1:A:69:THR:O	1:A:70:THR:C	2.63	0.41
3:C:37:LEU:HD22	3:C:37:LEU:H	1.86	0.41
4:D:71:GLY:O	4:D:72:GLU:HB3	2.20	0.41
2:F:84:THR:HG23	2:F:108:ILE:HA	2.02	0.41
4:H:142:LEU:HD12	4:H:142:LEU:HA	1.87	0.41
4:H:172:THR:CG2	4:H:185:LEU:HB2	2.50	0.41
1:A:32:LEU:HB3	1:A:72:PHE:CD1	2.56	0.41
4:D:81:ARG:N	4:D:82:PRO:CD	2.84	0.41
1:E:185:ALA:HA	1:E:199:PHE:CE1	2.55	0.41
2:F:23:GLN:HE21	2:F:23:GLN:HB3	1.49	0.41
3:G:12:SER:HB2	3:G:67:VAL:HG21	2.02	0.41
4:H:165:GLU:CD	4:H:167:VAL:HG22	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:ASN:O	1:A:116:PRO:O	2.39	0.41
4:D:159:PHE:CE1	4:D:164:GLU:HB2	2.56	0.41
1:E:57:SER:O	1:E:58:SER:HB2	2.20	0.41
3:G:30:ASP:HB3	4:H:180:TRP:CE2	2.56	0.41
3:C:45:MET:HE2	3:C:134:TYR:OH	2.21	0.40
4:D:159:PHE:CZ	4:D:164:GLU:HB2	2.56	0.40
3:G:102:GLN:O	3:G:156:PRO:HD2	2.20	0.40
1:A:51:ARG:HH11	4:D:102:ASP:CG	2.29	0.40
2:B:199:ASN:HA	2:B:200:PRO:HD2	1.92	0.40
1:A:47:PHE:CD1	1:A:47:PHE:C	2.99	0.40
4:D:152:ALA:HB1	4:D:174:LEU:HD21	2.04	0.40
1:E:163:MET:SD	2:F:191:ARG:HG2	2.61	0.40
3:G:161:ILE:H	3:G:161:ILE:CD1	2.25	0.40
1:A:123:LEU:N	1:A:123:LEU:HD12	2.36	0.40
2:B:23:GLN:HE21	2:B:23:GLN:HB3	1.50	0.40
2:B:80:SER:N	2:B:83:GLN:HE21	2.19	0.40
4:D:108:ASN:HA	4:D:112:PRO:HD2	2.04	0.40
2:F:199:ASN:OD1	2:F:201:ARG:HB2	2.22	0.40
4:H:138:HIS:O	4:H:191:THR:HG23	2.22	0.40
4:H:160:ARG:O	4:H:161:ASN:C	2.64	0.40
3:C:124:ARG:O	3:C:127:LYS:HB2	2.21	0.40
2:F:55:LYS:HB2	2:F:59:PRO:HG3	2.03	0.40
4:H:38:MET:HE2	4:H:40:GLU:OE1	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	200/202 (99%)	174 (87%)	19 (10%)	7 (4%)	3 1
1	E	199/202 (98%)	176 (88%)	16 (8%)	7 (4%)	3 1

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	238/240 (99%)	212 (89%)	23 (10%)	3 (1%)	9	12
2	F	238/240 (99%)	214 (90%)	21 (9%)	3 (1%)	9	12
3	C	180/182 (99%)	166 (92%)	12 (7%)	2 (1%)	11	15
3	G	180/182 (99%)	167 (93%)	11 (6%)	2 (1%)	11	15
4	D	198/217 (91%)	184 (93%)	13 (7%)	1 (0%)	24	34
4	H	198/217 (91%)	183 (92%)	14 (7%)	1 (0%)	24	34
All	All	1631/1682 (97%)	1476 (90%)	129 (8%)	26 (2%)	7	9

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	117	ASP
1	A	165	SER
4	D	161	ASN
1	E	165	SER
1	A	7	GLU
1	A	52	ASP
1	A	116	PRO
2	B	50	ALA
3	C	128	SER
1	E	7	GLU
1	E	52	ASP
2	F	50	ALA
1	A	8	GLY
1	A	127	LYS
2	B	97	ASN
3	C	2	GLU
1	E	8	GLY
1	E	127	LYS
1	E	200	PRO
2	F	97	ASN
3	G	2	GLU
3	G	128	SER
4	H	161	ASN
2	B	172	PRO
1	E	53	LYS
2	F	116	VAL

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	179/179 (100%)	173 (97%)	6 (3%)	32	48
1	E	178/179 (99%)	171 (96%)	7 (4%)	28	43
2	B	208/208 (100%)	197 (95%)	11 (5%)	20	31
2	F	208/208 (100%)	198 (95%)	10 (5%)	23	35
3	C	163/163 (100%)	151 (93%)	12 (7%)	13	17
3	G	163/163 (100%)	151 (93%)	12 (7%)	13	17
4	D	182/189 (96%)	176 (97%)	6 (3%)	33	50
4	H	182/189 (96%)	172 (94%)	10 (6%)	19	29
All	All	1463/1478 (99%)	1389 (95%)	74 (5%)	21	33

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LEU
1	A	62	GLU
1	A	91	LEU
1	A	112	ASN
1	A	114	GLN
1	A	186	ASN
2	B	2	VAL
2	B	12	VAL
2	B	13	THR
2	B	23	GLN
2	B	36	ASP
2	B	55	LYS
2	B	128	GLU
2	B	142	LEU
2	B	167	CYS
2	B	179	LEU
2	B	191	ARG
3	C	1	ILE
3	C	15	GLN

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Mol	Chain	Res	Type
3	C	16	SER
3	C	20	ILE
3	C	45	MET
3	C	52	LEU
3	C	99	LEU
3	C	142	ARG
3	C	143	ASP
3	C	152	LEU
3	C	161	ILE
3	C	173	GLU
4	D	4	GLN
4	D	118	LEU
4	D	126	VAL
4	D	167	VAL
4	D	189	GLU
4	D	207	LEU
1	E	12	LEU
1	E	62	GLU
1	E	91	LEU
1	E	112	ASN
1	E	114	GLN
1	E	186	ASN
1	E	201	SER
2	F	2	VAL
2	F	12	VAL
2	F	13	THR
2	F	23	GLN
2	F	36	ASP
2	F	55	LYS
2	F	142	LEU
2	F	179	LEU
2	F	191	ARG
2	F	207	GLN
3	G	1	ILE
3	G	15	GLN
3	G	16	SER
3	G	20	ILE
3	G	45	MET
3	G	52	LEU
3	G	99	LEU
3	G	142	ARG
3	G	143	ASP

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Mol	Chain	Res	Type
3	G	152	LEU
3	G	161	ILE
3	G	173	GLU
4	H	4	GLN
4	H	103	THR
4	H	118	LEU
4	H	126	VAL
4	H	142	LEU
4	H	167	VAL
4	H	171	SER
4	H	189	GLU
4	H	190	MET
4	H	207	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	189	ASN
1	A	190	ASN
2	B	4	GLN
2	B	25	ASN
2	B	69	GLN
2	B	83	GLN
2	B	115	ASN
2	B	150	HIS
3	C	15	GLN
3	C	79	ASN
3	C	102	GLN
3	C	112	ASN
3	C	168	HIS
4	D	4	GLN
4	D	32	HIS
4	D	45	ASN
4	D	48	GLN
4	D	137	ASN
4	D	161	ASN
4	D	173	GLN
1	E	9	ASN
1	E	189	ASN
1	E	190	ASN
2	F	4	GLN

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Mol	Chain	Res	Type
2	F	25	ASN
2	F	69	GLN
2	F	83	GLN
2	F	115	ASN
2	F	133	HIS
2	F	202	ASN
2	F	221	GLN
3	G	58	GLN
3	G	79	ASN
3	G	102	GLN
3	G	147	HIS
3	G	168	HIS
4	H	4	GLN
4	H	32	HIS
4	H	36	GLN
4	H	45	ASN
4	H	48	GLN
4	H	137	ASN

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 oligosaccharides 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 ⓘ

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	202/202 (100%)	1.02	27 (13%) 7 5	38, 70, 100, 123	0
1	E	201/202 (99%)	0.90	17 (8%) 16 16	42, 69, 100, 124	0
2	B	240/240 (100%)	0.77	19 (7%) 18 18	40, 63, 95, 124	0
2	F	240/240 (100%)	0.74	14 (5%) 29 28	39, 64, 94, 110	0
3	C	182/182 (100%)	0.40	5 (2%) 56 57	30, 54, 91, 112	0
3	G	182/182 (100%)	0.35	8 (4%) 39 39	29, 54, 92, 111	0
4	D	202/217 (93%)	0.36	3 (1%) 72 73	35, 53, 86, 114	0
4	H	202/217 (93%)	0.32	5 (2%) 58 59	33, 52, 86, 113	0
All	All	1651/1682 (98%)	0.62	98 (5%) 28 27	29, 61, 95, 124	0

All (98) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	113	ILE	6.0
3	C	1	ILE	5.3
2	B	50	ALA	5.2
1	A	116	PRO	5.1
1	E	131	LYS	4.4
1	E	81	GLU	4.4
1	A	127	LYS	4.2
2	F	29	TYR	3.7
1	A	129	SER	3.5
2	B	112	ASP	3.4
2	B	114	LYS	3.3
1	E	126	SER	3.1
2	F	113	LEU	3.0
1	A	128	SER	3.0
3	G	1	ILE	3.0
1	A	2	SER	3.0

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Mol	Chain	Res	Type	RSRZ
2	F	1	ALA	2.9
1	A	126	SER	2.9
1	A	115	ASN	2.9
2	F	74	LEU	2.9
1	E	129	SER	2.8
1	E	201	SER	2.8
1	E	115	ASN	2.8
1	A	110	LEU	2.8
2	B	109	VAL	2.8
2	B	113	LEU	2.8
2	F	114	LYS	2.8
2	B	116	VAL	2.7
1	A	114	GLN	2.7
1	A	40	GLY	2.7
3	C	101	GLY	2.7
1	E	113	ILE	2.6
2	B	148	PRO	2.6
2	F	218	GLU	2.6
3	G	160	ASP	2.6
1	E	159	CYS	2.6
2	B	115	ASN	2.5
2	F	50	ALA	2.5
1	A	72	PHE	2.5
1	A	7	GLU	2.5
1	A	131	LYS	2.5
1	E	6	THR	2.5
4	H	133	THR	2.5
1	A	140	ASP	2.5
2	F	110	VAL	2.4
1	A	112	ASN	2.4
1	A	79	VAL	2.4
2	F	234	GLU	2.4
2	B	181	ASP	2.4
2	B	167	CYS	2.4
3	G	81	THR	2.3
3	G	161	ILE	2.3
3	C	160	ASP	2.3
2	B	177	PRO	2.3
3	C	81	THR	2.3
2	F	51	ASP	2.3
1	A	159	CYS	2.3
3	G	41	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
4	H	14	GLY	2.3
1	A	162	ASP	2.3
3	G	163	ASP	2.3
1	A	184	CYS	2.3
2	B	48	TYR	2.3
1	A	132	SER	2.3
3	G	80	SER	2.3
1	E	112	ASN	2.2
2	F	167	CYS	2.3
4	H	217	ALA	2.2
4	H	195	GLY	2.2
1	A	50	SER	2.2
1	E	12	LEU	2.2
4	H	33	PHE	2.2
2	F	49	VAL	2.2
2	B	19	LEU	2.2
2	F	111	GLU	2.2
1	E	182	PHE	2.2
2	F	12	VAL	2.2
1	A	130	ASP	2.2
1	E	162	ASP	2.2
2	B	222	ASP	2.2
4	D	195	GLY	2.1
1	A	125	ASP	2.1
3	C	158	ASP	2.1
2	B	175	GLU	2.1
1	E	184	CYS	2.1
2	B	239	ALA	2.1
1	A	124	ARG	2.1
1	A	111	PRO	2.1
2	B	150	HIS	2.1
1	E	114	GLN	2.1
1	E	124	ARG	2.1
4	D	194	ARG	2.1
2	B	57	ASP	2.1
3	G	128	SER	2.1
1	E	125	ASP	2.0
4	D	33	PHE	2.0
2	B	78	LEU	2.0
1	A	83	ASP	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 oligosaccharides 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.