



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:09 AM UTC

PDB ID : 3X2R / pdb_00003x2r
Title : Structure of the nonameric bacterial amyloid secretion channel CsgG
Authors : Huang, Y.; Cao, B.; Zhao, Y.; Kou, Y.; Ni, D.; Zhang, X.C.
Deposited on : 2014-12-29
Resolution : 2.90 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

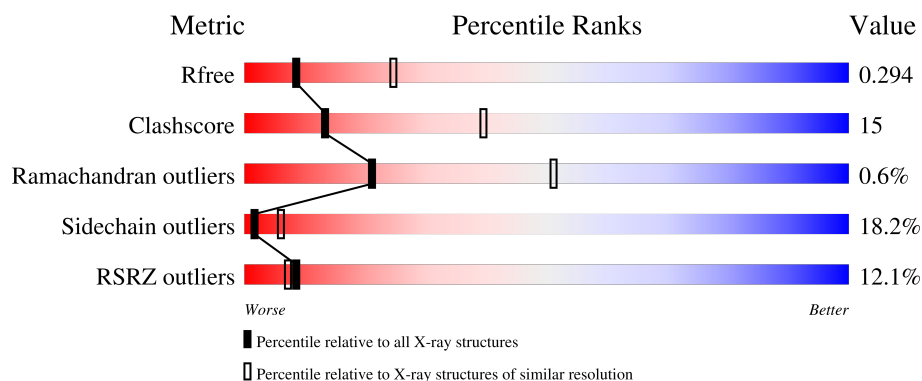
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	277	8% 41% 23% 6% 29%
1	B	277	6% 44% 21% • 31%
1	C	277	10% 42% 24% 5% 29%
1	D	277	10% 41% 22% 5% 32%
1	E	277	9% 42% 26% • 29%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	277	6% 42% 21% 5% 32%
1	G	277	12% 39% 25% 5% 31%
1	H	277	6% 39% 24% 5% 31%
1	I	277	9% 42% 23% • 31%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 13042 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.

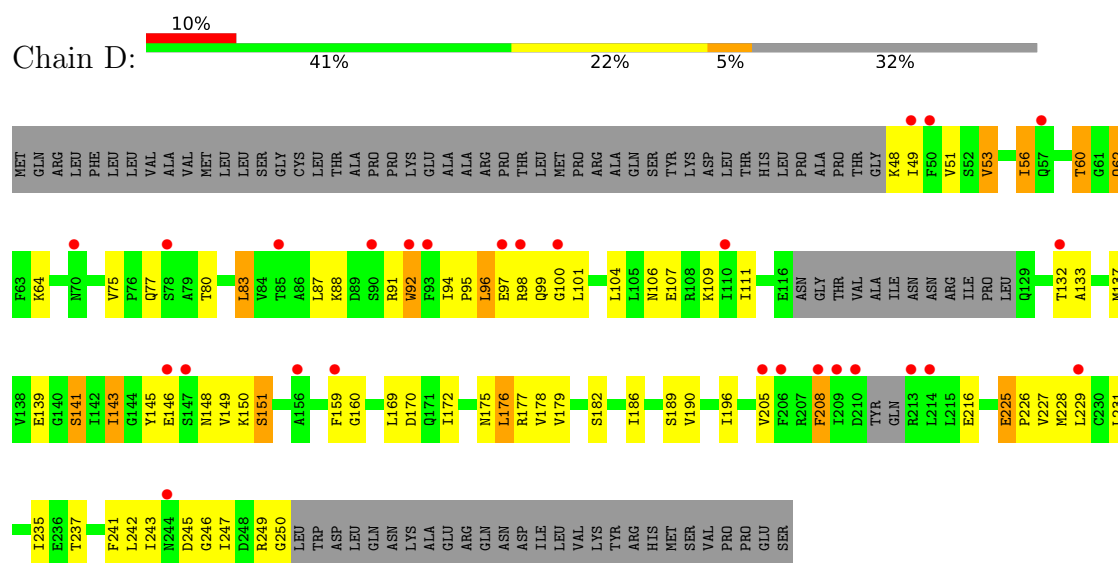
- Molecule 1 is a protein called CsgG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	189	Total	C	N	O	S	0	0	0
			1424	904	236	279	5			
1	A	196	Total	C	N	O	S	0	0	0
			1467	932	243	287	5			
1	B	190	Total	C	N	O	S	0	0	0
			1427	904	236	282	5			
1	C	198	Total	C	N	O	S	0	0	0
			1498	954	248	291	5			
1	E	198	Total	C	N	O	S	0	0	0
			1498	954	248	291	5			
1	F	189	Total	C	N	O	S	0	0	0
			1424	904	236	279	5			
1	G	190	Total	C	N	O	S	0	0	0
			1440	918	237	280	5			
1	H	191	Total	C	N	O	S	0	0	0
			1441	915	238	283	5			
1	I	190	Total	C	N	O	S	0	0	0
			1423	902	236	280	5			

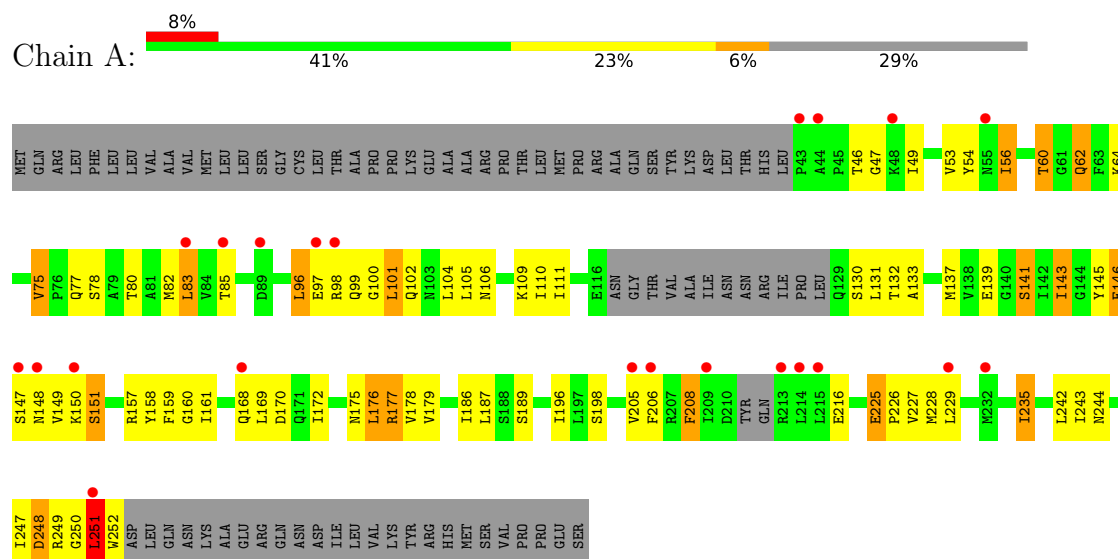
3 Residue-property plots [i](#)

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.

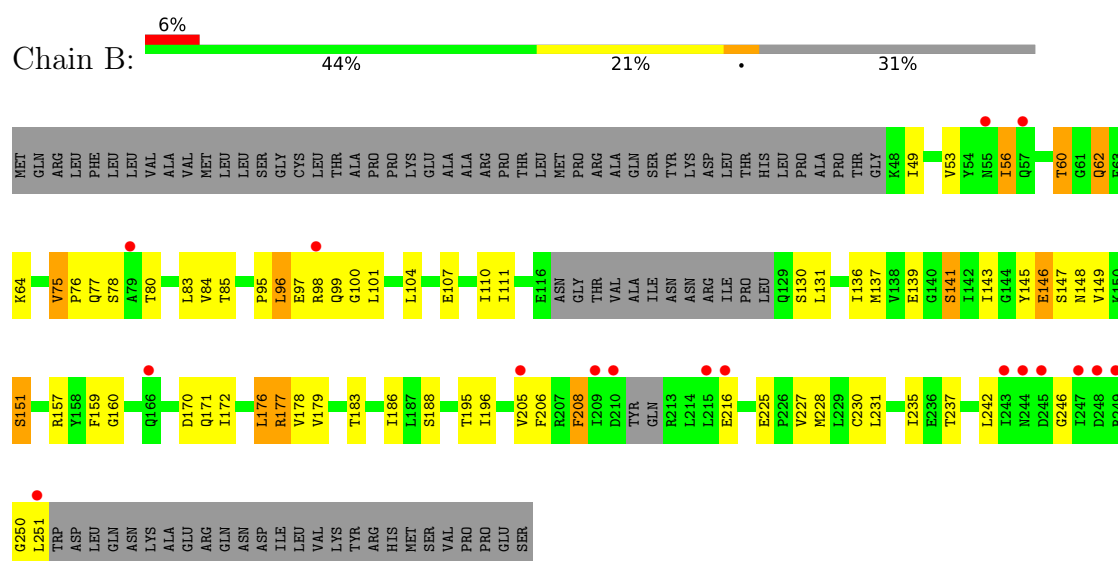
• Molecule 1: CsgG



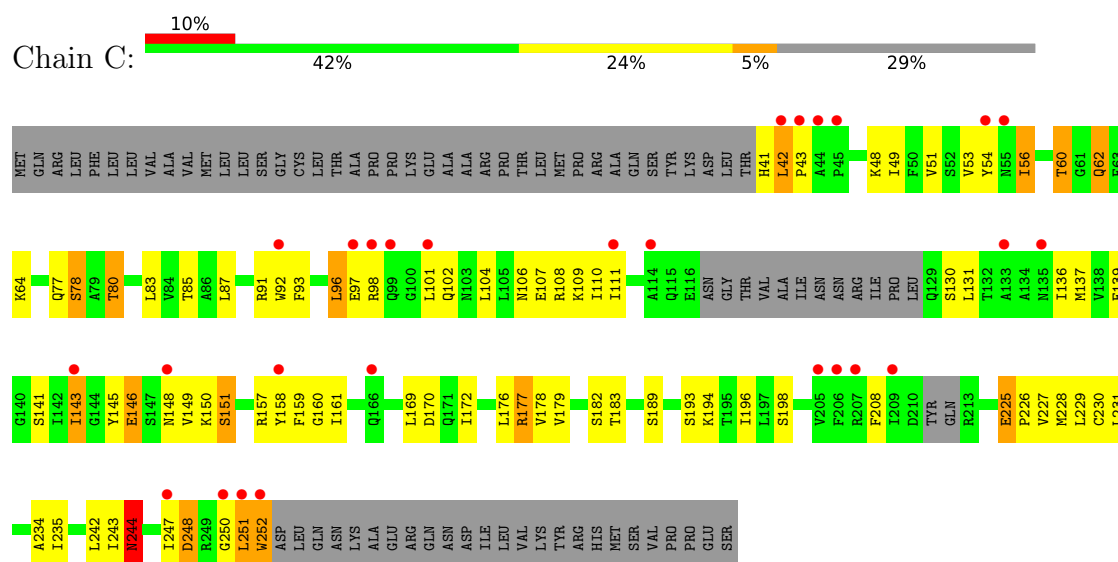
• Molecule 1: CsgG



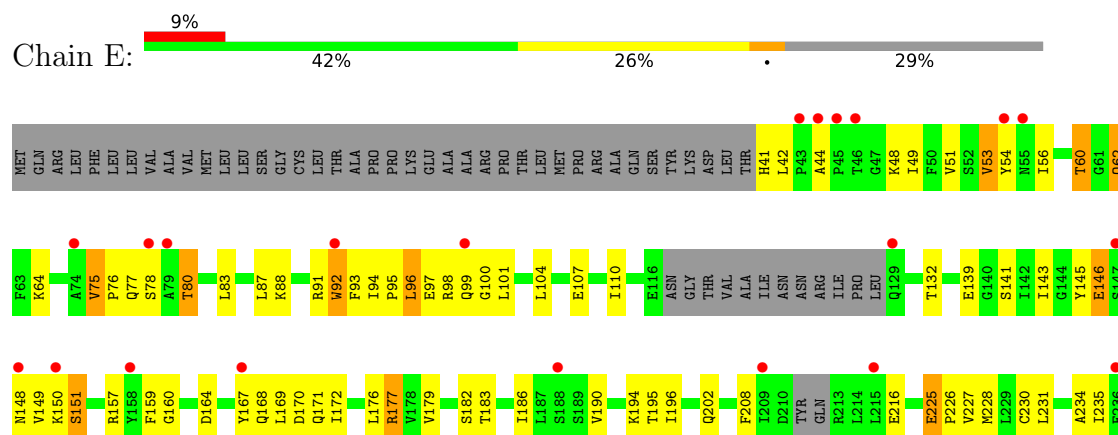
• Molecule 1: CsgG

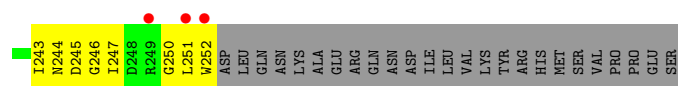


- Molecule 1: CsgG

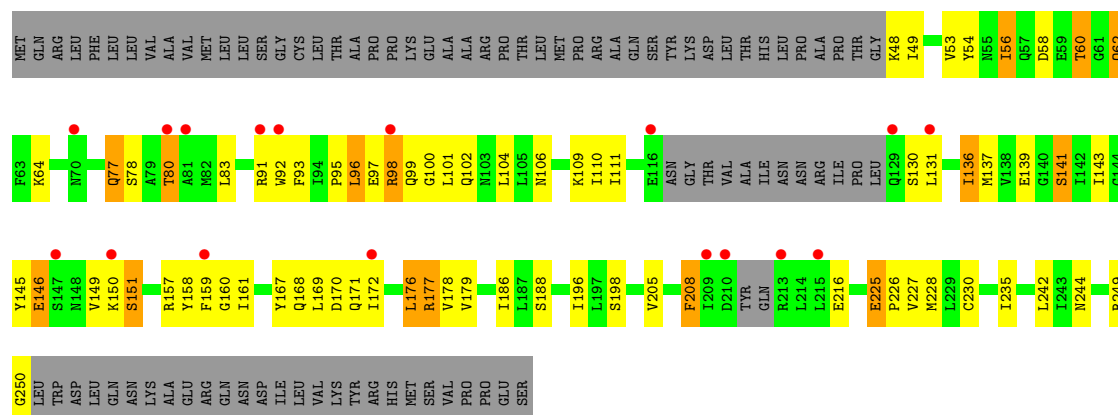


- Molecule 1: CsgG

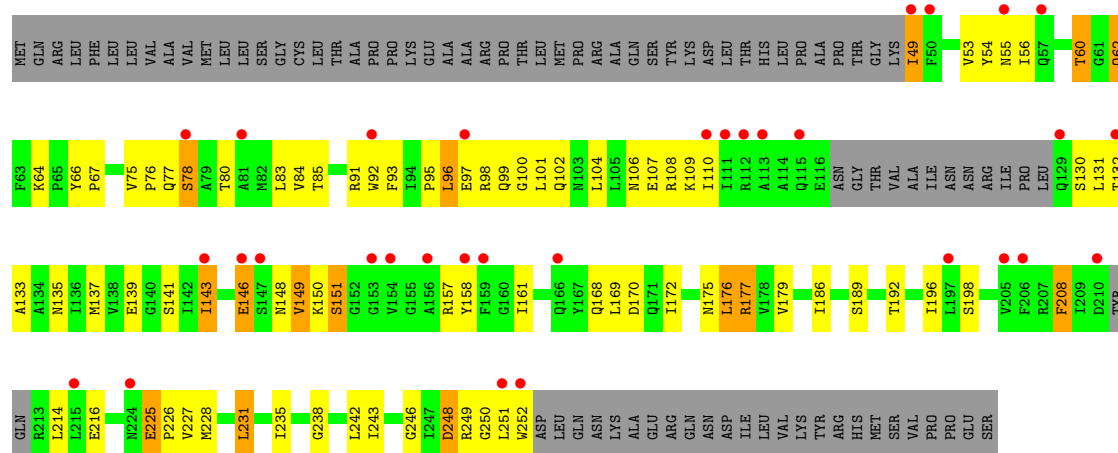




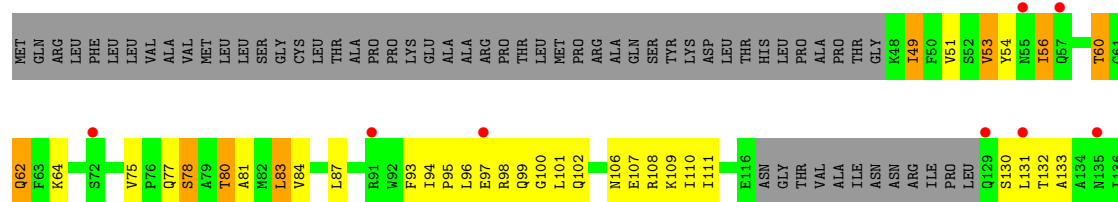
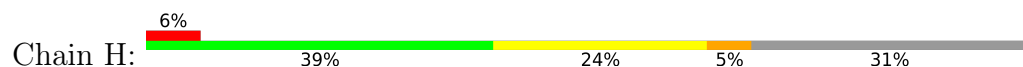
• Molecule 1: CsgG

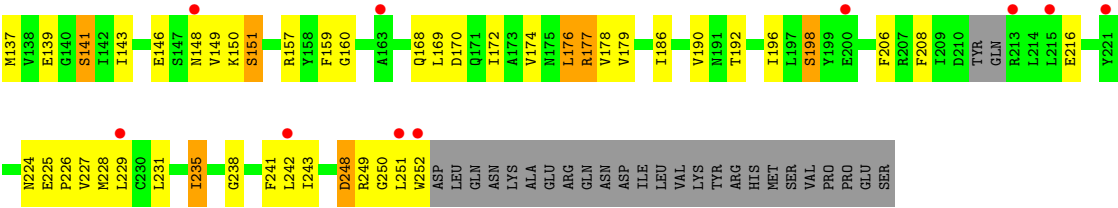


• Molecule 1: CsgG

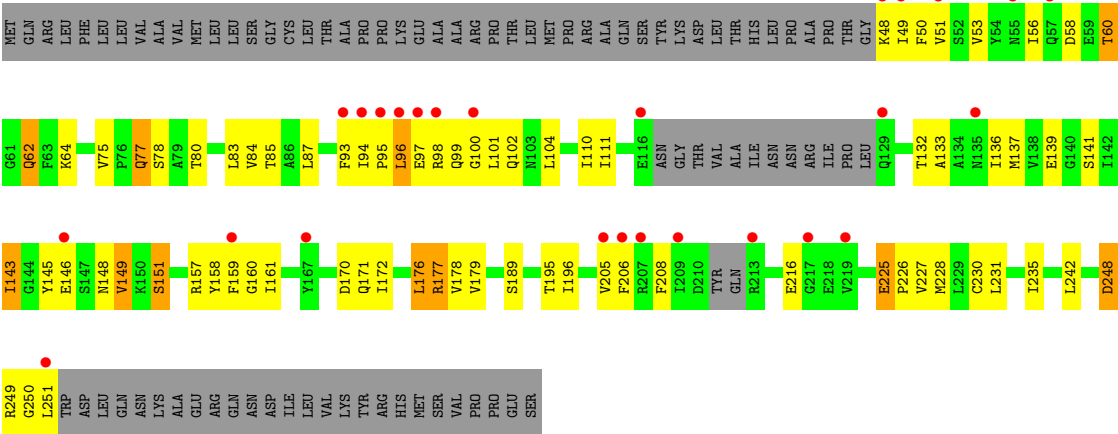


• Molecule 1: CsgG





● Molecule 1: CsgG



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	171.88Å 176.72Å 103.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.07 – 2.90 48.07 – 2.90	Depositor EDS
% Data completeness (in resolution range)	86.8 (48.07-2.90) 77.2 (48.07-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.244 , 0.292 0.247 , 0.294	Depositor DCC
R_{free} test set	3428 reflections (4.35%)	wwPDB-VP
Wilson B-factor (Å ²)	58.7	Xtriage
Anisotropy	0.943	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 48.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.012 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	13042	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.18% 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	0.46	0/1488	0.96	5/2019 (0.2%)
1	B	0.44	0/1444	0.93	3/1956 (0.2%)
1	C	0.55	0/1522	1.00	4/2067 (0.2%)
1	D	0.47	0/1443	0.93	4/1956 (0.2%)
1	E	0.49	0/1522	0.97	6/2067 (0.3%)
1	F	0.44	0/1443	0.89	3/1956 (0.2%)
1	G	0.45	0/1461	0.91	4/1983 (0.2%)
1	H	0.44	0/1460	0.93	6/1979 (0.3%)
1	I	0.44	0/1440	0.92	2/1951 (0.1%)
All	All	0.47	0/13223	0.94	37/17934 (0.2%)

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	G	0	1

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	225	GLU	CA-C-N	8.30	130.21	119.84
1	I	225	GLU	C-N-CA	8.30	130.21	119.84
1	A	146	GLU	N-CA-C	7.61	118.65	108.38
1	C	225	GLU	CA-C-N	7.52	129.24	119.84
1	C	225	GLU	C-N-CA	7.52	129.24	119.84
1	C	146	GLU	N-CA-C	7.05	117.89	108.38
1	E	225	GLU	CA-C-N	6.92	128.49	119.84
1	E	225	GLU	C-N-CA	6.92	128.49	119.84
1	B	225	GLU	CA-C-N	6.76	128.30	119.84

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	225	GLU	C-N-CA	6.76	128.30	119.84
1	D	146	GLU	N-CA-C	6.75	117.50	108.38
1	D	225	GLU	CA-C-N	6.58	128.07	119.84
1	D	225	GLU	C-N-CA	6.58	128.07	119.84
1	A	225	GLU	CA-C-N	6.46	127.91	119.84
1	A	225	GLU	C-N-CA	6.46	127.91	119.84
1	F	146	GLU	N-CA-C	6.36	116.96	108.38
1	G	225	GLU	CA-C-N	6.32	127.74	119.84
1	G	225	GLU	C-N-CA	6.32	127.74	119.84
1	F	225	GLU	CA-C-N	6.17	127.55	119.84
1	F	225	GLU	C-N-CA	6.17	127.55	119.84
1	E	146	GLU	N-CA-C	6.14	116.67	108.38
1	H	225	GLU	CA-C-N	6.06	127.41	119.84
1	H	225	GLU	C-N-CA	6.06	127.41	119.84
1	B	146	GLU	N-CA-C	5.86	116.29	108.38
1	H	78	SER	N-CA-C	5.74	120.12	113.18
1	E	44	ALA	CA-C-N	5.46	126.67	119.84
1	E	44	ALA	C-N-CA	5.46	126.67	119.84
1	C	244	ASN	N-CA-CB	5.40	117.83	110.01
1	A	75	VAL	CA-C-N	5.37	125.31	119.78
1	A	75	VAL	C-N-CA	5.37	125.31	119.78
1	G	78	SER	N-CA-C	5.36	119.67	113.18
1	E	78	SER	N-CA-C	5.21	119.49	113.18
1	H	146	GLU	N-CA-C	5.21	115.42	108.38
1	H	94	ILE	CA-C-N	5.19	125.49	119.93
1	H	94	ILE	C-N-CA	5.19	125.49	119.93
1	D	96	LEU	N-CA-C	5.01	119.08	112.92
1	G	146	GLU	N-CA-C	5.00	115.14	108.38

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	G	55	ASN	Peptide

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	1467	0	1451	58	0
1	B	1427	0	1415	44	0
1	C	1498	0	1480	53	1
1	D	1424	0	1408	46	0
1	E	1498	0	1480	54	1
1	F	1424	0	1408	45	0
1	G	1440	0	1425	56	0
1	H	1441	0	1425	55	0
1	I	1423	0	1411	55	0
All	All	13042	0	12903	384	1

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)
1:E:92:TRP:HE3	1:E:92:TRP:H	1.22	0.83
1:C:150:LYS:HB3	1:C:169:LEU:HD12	1.61	0.82
1:E:196:ILE:HD13	1:E:227:VAL:HG22	1.64	0.80
1:G:49:ILE:HG21	1:G:92:TRP:HB2	1.65	0.79
1:E:48:LYS:HG2	1:E:92:TRP:HB2	1.66	0.77
1:G:170:ASP:HB2	1:G:196:ILE:HB	1.65	0.76
1:I:176:LEU:HB3	1:I:242:LEU:HD11	1.68	0.76
1:B:176:LEU:HB3	1:B:242:LEU:HD11	1.67	0.75
1:A:46:THR:HB	1:A:47:GLY:HA3	1.67	0.74
1:H:150:LYS:HB3	1:H:169:LEU:HD12	1.71	0.73
1:A:137:MET:HE1	1:B:104:LEU:HA	1.71	0.72
1:D:48:LYS:HD3	1:D:94:ILE:HD11	1.70	0.71
1:D:176:LEU:HB3	1:D:242:LEU:HD11	1.70	0.71
1:F:48:LYS:HA	1:F:92:TRP:HB2	1.70	0.71
1:H:172:ILE:HD11	1:H:227:VAL:HG13	1.72	0.71
1:D:170:ASP:HB2	1:D:196:ILE:HB	1.73	0.71
1:E:177:ARG:NH1	1:H:97:GLU:OE2	2.24	0.70
1:G:177:ARG:NH1	1:I:97:GLU:OE2	2.25	0.69
1:E:150:LYS:HB3	1:E:169:LEU:HD12	1.73	0.69
1:G:137:MET:HE1	1:I:104:LEU:HA	1.75	0.69
1:H:176:LEU:HB3	1:H:242:LEU:HD11	1.76	0.67
1:D:97:GLU:OE2	1:F:177:ARG:NH1	2.27	0.67
1:G:172:ILE:HD11	1:G:227:VAL:HG13	1.76	0.66
1:H:56:ILE:HG13	1:H:141:SER:HA	1.76	0.66
1:E:172:ILE:HD11	1:E:227:VAL:HG13	1.77	0.65

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:84:VAL:HG13	1:H:95:PRO:HB3	1.77	0.65
1:I:84:VAL:HG13	1:I:95:PRO:HB3	1.78	0.65
1:A:83:LEU:HD23	1:A:235:ILE:HG12	1.79	0.65
1:F:172:ILE:HD11	1:F:227:VAL:HG13	1.79	0.65
1:F:178:VAL:HG23	1:F:242:LEU:HD22	1.79	0.65
1:F:196:ILE:HD13	1:F:227:VAL:HG22	1.78	0.64
1:G:150:LYS:HB3	1:G:169:LEU:HD12	1.77	0.64
1:C:172:ILE:HD11	1:C:227:VAL:HG13	1.79	0.64
1:F:176:LEU:HB3	1:F:242:LEU:HD11	1.80	0.64
1:I:157:ARG:HH22	1:I:206:PHE:HB2	1.63	0.64
1:E:194:LYS:HE3	1:E:234:ALA:HA	1.79	0.64
1:G:176:LEU:HB3	1:G:242:LEU:HD11	1.80	0.63
1:F:104:LEU:HA	1:H:137:MET:HE1	1.80	0.63
1:C:136:ILE:HG21	1:C:251:LEU:HD11	1.80	0.63
1:C:148:ASN:HB3	1:C:151:SER:HB3	1.81	0.63
1:C:196:ILE:HD13	1:C:227:VAL:HG22	1.81	0.62
1:D:104:LEU:HA	1:F:137:MET:HE1	1.81	0.62
1:C:182:SER:O	1:G:108:ARG:NH1	2.32	0.62
1:F:97:GLU:OE2	1:H:177:ARG:NH1	2.33	0.61
1:B:186:ILE:HD12	1:E:95:PRO:HB2	1.82	0.61
1:A:98:ARG:HH22	1:B:110:ILE:HG21	1.65	0.61
1:G:196:ILE:HD13	1:G:227:VAL:HG22	1.83	0.61
1:I:172:ILE:HD11	1:I:227:VAL:HG13	1.82	0.61
1:A:172:ILE:HD11	1:A:227:VAL:HG13	1.81	0.61
1:F:60:THR:HG23	1:F:62:GLN:H	1.65	0.61
1:F:110:ILE:HD13	1:H:98:ARG:NH2	2.15	0.61
1:A:157:ARG:HH22	1:A:206:PHE:HB2	1.66	0.60
1:G:60:THR:HG23	1:G:62:GLN:H	1.66	0.60
1:E:98:ARG:HH22	1:H:110:ILE:HG21	1.66	0.60
1:A:196:ILE:HD13	1:A:227:VAL:HG22	1.83	0.60
1:C:48:LYS:HG2	1:C:92:TRP:HB2	1.81	0.60
1:H:196:ILE:HD13	1:H:227:VAL:HG22	1.83	0.60
1:D:107:GLU:OE2	1:F:98:ARG:NH1	2.35	0.60
1:C:194:LYS:HE3	1:C:234:ALA:HA	1.84	0.60
1:E:60:THR:HG23	1:E:62:GLN:H	1.65	0.60
1:A:177:ARG:NH1	1:B:97:GLU:OE2	2.34	0.59
1:D:150:LYS:HB3	1:D:169:LEU:HD12	1.84	0.59
1:G:84:VAL:HG13	1:G:95:PRO:HB3	1.85	0.59
1:D:60:THR:HG23	1:D:62:GLN:HB2	1.83	0.59
1:C:170:ASP:HB2	1:C:196:ILE:HB	1.83	0.59
1:G:98:ARG:HH22	1:I:110:ILE:HG21	1.67	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:148:ASN:HB3	1:H:151:SER:HB3	1.85	0.59
1:H:170:ASP:HB2	1:H:196:ILE:HB	1.83	0.59
1:B:137:MET:HE1	1:E:104:LEU:HA	1.85	0.59
1:E:170:ASP:HB2	1:E:196:ILE:HB	1.85	0.59
1:A:248:ASP:OD1	1:A:248:ASP:N	2.36	0.59
1:B:178:VAL:HG23	1:B:242:LEU:HD22	1.84	0.58
1:G:98:ARG:NH2	1:I:110:ILE:HD13	2.19	0.58
1:G:186:ILE:HD12	1:I:95:PRO:HB2	1.86	0.58
1:A:170:ASP:HB2	1:A:196:ILE:HB	1.86	0.57
1:B:60:THR:HG23	1:B:62:GLN:HB2	1.86	0.57
1:F:150:LYS:HB3	1:F:169:LEU:HD12	1.85	0.57
1:I:248:ASP:N	1:I:248:ASP:OD1	2.37	0.57
1:F:110:ILE:HG21	1:H:98:ARG:HH22	1.69	0.57
1:B:170:ASP:HB2	1:B:196:ILE:HB	1.85	0.57
1:B:171:GLN:HG2	1:B:195:THR:HG22	1.87	0.57
1:A:98:ARG:NH1	1:B:107:GLU:OE2	2.37	0.57
1:H:53:VAL:HG21	1:H:83:LEU:HD12	1.86	0.56
1:C:60:THR:HG23	1:C:62:GLN:H	1.70	0.56
1:H:60:THR:HG23	1:H:62:GLN:HB2	1.86	0.56
1:D:98:ARG:NH1	1:C:107:GLU:OE2	2.39	0.56
1:A:104:LEU:HA	1:I:137:MET:HE1	1.88	0.56
1:I:60:THR:HG23	1:I:62:GLN:HB2	1.87	0.56
1:A:60:THR:HG23	1:A:62:GLN:HB2	1.87	0.56
1:D:196:ILE:HD13	1:D:227:VAL:HG22	1.88	0.56
1:E:53:VAL:HG22	1:E:80:THR:HG23	1.88	0.55
1:A:60:THR:HG23	1:A:62:GLN:H	1.72	0.55
1:C:248:ASP:OD1	1:C:248:ASP:N	2.40	0.55
1:G:54:TYR:CZ	1:G:98:ARG:HG2	2.40	0.55
1:G:248:ASP:N	1:G:248:ASP:OD1	2.38	0.55
1:F:170:ASP:HB2	1:F:196:ILE:HB	1.87	0.55
1:I:60:THR:HG23	1:I:62:GLN:H	1.72	0.55
1:H:157:ARG:HH22	1:H:206:PHE:HB2	1.71	0.55
1:E:54:TYR:CZ	1:E:98:ARG:HG2	2.42	0.55
1:F:60:THR:HG23	1:F:62:GLN:HB2	1.88	0.55
1:H:248:ASP:OD1	1:H:248:ASP:N	2.40	0.54
1:C:145:TYR:OH	1:C:170:ASP:OD2	2.18	0.54
1:A:249:ARG:N	1:A:250:GLY:HA2	2.22	0.54
1:D:137:MET:HE1	1:C:104:LEU:HA	1.90	0.54
1:B:196:ILE:HD13	1:B:227:VAL:HG22	1.89	0.54
1:C:92:TRP:H	1:C:92:TRP:HE3	1.54	0.54
1:E:51:VAL:HG11	1:E:87:LEU:HD13	1.89	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:148:ASN:HB3	1:E:151:SER:HB3	1.90	0.54
1:D:60:THR:HG23	1:D:62:GLN:H	1.73	0.53
1:E:60:THR:HG23	1:E:62:GLN:HB2	1.90	0.53
1:E:186:ILE:HD12	1:H:95:PRO:HB2	1.89	0.53
1:B:157:ARG:HH22	1:B:206:PHE:HB2	1.73	0.53
1:I:178:VAL:HG23	1:I:242:LEU:HD22	1.90	0.53
1:I:196:ILE:HD13	1:I:227:VAL:HG22	1.90	0.53
1:E:98:ARG:NH2	1:H:110:ILE:HD13	2.23	0.53
1:B:56:ILE:HG13	1:B:141:SER:HA	1.91	0.53
1:A:178:VAL:HG23	1:A:242:LEU:HD22	1.90	0.52
1:A:147:SER:HB3	1:A:170:ASP:CG	2.34	0.52
1:E:182:SER:O	1:H:108:ARG:NH1	2.41	0.52
1:G:146:GLU:OE1	1:I:226:PRO:HB3	2.09	0.52
1:G:148:ASN:HB3	1:G:151:SER:HB3	1.91	0.52
1:I:145:TYR:HA	1:I:171:GLN:O	2.10	0.52
1:E:171:GLN:HG2	1:E:195:THR:HG22	1.91	0.51
1:A:133:ALA:HB2	1:B:110:ILE:HG22	1.92	0.51
1:C:60:THR:HG23	1:C:62:GLN:HB2	1.91	0.51
1:B:98:ARG:HH22	1:E:110:ILE:HG21	1.76	0.51
1:F:99:GLN:HB3	1:F:100:GLY:HA3	1.92	0.51
1:D:95:PRO:HB2	1:F:186:ILE:HD12	1.93	0.51
1:A:98:ARG:NH2	1:B:110:ILE:HD13	2.26	0.51
1:C:194:LYS:HG3	1:C:230:CYS:O	2.11	0.51
1:F:111:ILE:HG12	1:H:133:ALA:HB1	1.93	0.51
1:G:133:ALA:HB2	1:I:110:ILE:HG22	1.92	0.51
1:G:143:ILE:HA	1:I:78:SER:HB3	1.93	0.50
1:D:186:ILE:HG12	1:C:97:GLU:HG2	1.93	0.50
1:G:99:GLN:HB3	1:G:100:GLY:HA3	1.92	0.50
1:D:178:VAL:HG23	1:D:242:LEU:HD22	1.93	0.50
1:A:97:GLU:OE2	1:I:177:ARG:NH1	2.45	0.50
1:I:58:ASP:HB3	1:I:77:GLN:NE2	2.26	0.50
1:E:194:LYS:HE3	1:E:234:ALA:CA	2.41	0.50
1:C:51:VAL:HG11	1:C:87:LEU:HD13	1.93	0.49
1:A:110:ILE:HD13	1:I:98:ARG:NH2	2.26	0.49
1:D:99:GLN:HB3	1:D:100:GLY:HA3	1.94	0.49
1:D:143:ILE:HG12	1:C:78:SER:HB2	1.94	0.49
1:E:145:TYR:OH	1:E:170:ASP:OD2	2.18	0.49
1:H:151:SER:HB2	1:H:168:GLN:HG2	1.94	0.49
1:D:133:ALA:HB1	1:C:111:ILE:HG12	1.95	0.49
1:D:148:ASN:HB3	1:D:151:SER:HB3	1.95	0.49
1:C:98:ARG:HH22	1:G:110:ILE:HG21	1.78	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:246:GLY:HA3	1:G:252:TRP:CZ3	2.47	0.49
1:H:54:TYR:O	1:H:80:THR:HG21	2.12	0.49
1:H:99:GLN:HB3	1:H:100:GLY:HA3	1.94	0.49
1:A:176:LEU:HB3	1:A:242:LEU:HD11	1.93	0.49
1:A:99:GLN:HB3	1:A:100:GLY:HA3	1.95	0.48
1:E:98:ARG:NH2	1:H:110:ILE:HG21	2.26	0.48
1:A:243:ILE:O	1:A:247:ILE:HG13	2.11	0.48
1:B:60:THR:HG23	1:B:62:GLN:H	1.78	0.48
1:B:99:GLN:HB3	1:B:100:GLY:HA3	1.96	0.48
1:B:98:ARG:CZ	1:E:110:ILE:HD13	2.44	0.48
1:B:148:ASN:HB3	1:B:151:SER:HB3	1.95	0.48
1:F:60:THR:HG23	1:F:62:GLN:N	2.28	0.48
1:G:246:GLY:O	1:G:250:GLY:HA2	2.13	0.48
1:H:250:GLY:C	1:H:252:TRP:H	2.20	0.48
1:E:93:PHE:HE2	1:E:243:ILE:HG12	1.78	0.48
1:C:137:MET:HE1	1:G:104:LEU:HA	1.95	0.48
1:I:148:ASN:HB3	1:I:151:SER:HB3	1.95	0.48
1:C:159:PHE:HA	1:C:160:GLY:HA2	1.52	0.48
1:D:51:VAL:HG11	1:D:87:LEU:HD13	1.96	0.48
1:A:54:TYR:CZ	1:A:98:ARG:HG2	2.49	0.48
1:B:145:TYR:HD2	1:B:172:ILE:HG12	1.77	0.48
1:A:148:ASN:HB3	1:A:151:SER:HB3	1.96	0.47
1:D:190:VAL:HG11	1:D:241:PHE:HD2	1.79	0.47
1:D:205:VAL:O	1:D:216:GLU:HA	2.15	0.47
1:C:250:GLY:O	1:C:252:TRP:NE1	2.46	0.47
1:D:249:ARG:HB3	1:D:250:GLY:O	2.14	0.47
1:G:130:SER:OG	1:G:131:LEU:N	2.47	0.47
1:D:172:ILE:HD11	1:D:227:VAL:HG13	1.96	0.47
1:A:110:ILE:HG21	1:I:98:ARG:HH22	1.79	0.47
1:E:183:THR:HB	1:H:132:THR:HG23	1.95	0.47
1:E:194:LYS:HG3	1:E:230:CYS:O	2.14	0.47
1:D:56:ILE:HG13	1:D:141:SER:HA	1.96	0.47
1:A:186:ILE:HG12	1:B:97:GLU:HG2	1.97	0.47
1:F:54:TYR:CZ	1:F:98:ARG:HG2	2.50	0.47
1:H:60:THR:HG23	1:H:62:GLN:H	1.79	0.47
1:C:157:ARG:HB3	1:G:216:GLU:HG3	1.97	0.47
1:E:60:THR:HG23	1:E:62:GLN:N	2.29	0.47
1:F:136:ILE:HD12	1:F:178:VAL:HG13	1.96	0.47
1:C:183:THR:HB	1:G:132:THR:HG23	1.97	0.47
1:F:95:PRO:HB2	1:H:186:ILE:HD12	1.96	0.47
1:G:98:ARG:NH2	1:I:110:ILE:HG21	2.30	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:159:PHE:HA	1:D:160:GLY:HA2	1.52	0.47
1:A:56:ILE:HG13	1:A:141:SER:HA	1.97	0.46
1:I:146:GLU:HG2	1:I:149:VAL:HG22	1.97	0.46
1:A:159:PHE:HA	1:A:160:GLY:HA2	1.52	0.46
1:C:143:ILE:HA	1:G:78:SER:HB3	1.98	0.46
1:D:48:LYS:HA	1:D:92:TRP:HB3	1.96	0.46
1:A:98:ARG:NH2	1:B:110:ILE:HG21	2.29	0.46
1:A:249:ARG:H	1:A:250:GLY:HA2	1.80	0.46
1:G:49:ILE:HG22	1:G:93:PHE:HA	1.98	0.46
1:G:60:THR:HG23	1:G:62:GLN:HB2	1.98	0.46
1:I:58:ASP:HB3	1:I:77:GLN:HE22	1.80	0.46
1:F:196:ILE:HG12	1:F:230:CYS:SG	2.56	0.46
1:I:249:ARG:H	1:I:250:GLY:HA2	1.81	0.46
1:D:145:TYR:HD2	1:D:172:ILE:HG12	1.81	0.46
1:G:60:THR:HG23	1:G:62:GLN:N	2.30	0.46
1:A:78:SER:OG	1:A:82:MET:HE3	2.16	0.46
1:A:151:SER:HB2	1:A:168:GLN:HG2	1.98	0.46
1:E:97:GLU:HB3	1:E:98:ARG:H	1.54	0.46
1:E:164:ASP:OD1	1:E:202:GLN:HB2	2.16	0.46
1:I:159:PHE:HA	1:I:160:GLY:HA2	1.47	0.46
1:D:98:ARG:HH22	1:C:110:ILE:HG21	1.81	0.45
1:D:182:SER:O	1:C:108:ARG:NH1	2.49	0.45
1:G:250:GLY:C	1:G:252:TRP:H	2.24	0.45
1:I:145:TYR:HD2	1:I:172:ILE:HG12	1.81	0.45
1:H:243:ILE:HG23	1:H:252:TRP:HH2	1.81	0.45
1:D:98:ARG:NH2	1:C:110:ILE:HD13	2.32	0.45
1:B:246:GLY:O	1:B:250:GLY:HA2	2.16	0.45
1:E:225:GLU:HA	1:E:226:PRO:HD3	1.71	0.45
1:H:198:SER:HB3	1:H:224:ASN:HD22	1.81	0.45
1:D:143:ILE:HD12	1:D:175:ASN:HB2	1.98	0.45
1:B:147:SER:HB3	1:B:170:ASP:CG	2.41	0.45
1:C:130:SER:OG	1:C:131:LEU:N	2.50	0.45
1:H:54:TYR:CZ	1:H:98:ARG:HG2	2.52	0.45
1:I:51:VAL:HG11	1:I:87:LEU:HD13	1.97	0.45
1:A:145:TYR:HD2	1:A:172:ILE:HG12	1.81	0.45
1:E:246:GLY:O	1:E:250:GLY:HA2	2.17	0.45
1:B:186:ILE:HG12	1:E:97:GLU:HG2	1.99	0.45
1:H:249:ARG:H	1:H:250:GLY:HA2	1.81	0.45
1:I:50:PHE:HD1	1:I:94:ILE:HB	1.82	0.45
1:B:159:PHE:HA	1:B:160:GLY:HA2	1.51	0.45
1:C:178:VAL:HG23	1:C:242:LEU:HD13	1.98	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:194:LYS:HE3	1:C:234:ALA:CA	2.46	0.45
1:E:194:LYS:HE3	1:E:234:ALA:N	2.32	0.45
1:A:101:LEU:HD22	1:A:105:LEU:HG	1.98	0.45
1:D:91:ARG:HA	1:D:92:TRP:C	2.42	0.44
1:F:225:GLU:HA	1:F:226:PRO:HD3	1.69	0.44
1:I:99:GLN:HB3	1:I:100:GLY:HA3	1.97	0.44
1:A:187:LEU:HD13	1:A:251:LEU:HD13	1.99	0.44
1:C:225:GLU:HA	1:C:226:PRO:HD3	1.70	0.44
1:D:88:LYS:HE2	1:F:188:SER:HA	2.00	0.44
1:F:205:VAL:O	1:F:216:GLU:HA	2.18	0.44
1:E:151:SER:HB2	1:E:168:GLN:HG2	1.99	0.44
1:A:97:GLU:HB3	1:A:98:ARG:H	1.58	0.44
1:B:107:GLU:O	1:B:111:ILE:HG13	2.17	0.44
1:C:56:ILE:HG23	1:C:80:THR:OG1	2.18	0.44
1:B:98:ARG:NH2	1:E:110:ILE:HD13	2.33	0.44
1:E:99:GLN:HB3	1:E:100:GLY:HA3	2.00	0.44
1:F:96:LEU:HD22	1:F:96:LEU:HA	1.74	0.44
1:I:205:VAL:O	1:I:216:GLU:HA	2.18	0.44
1:D:189:SER:OG	1:C:85:THR:HG22	2.17	0.44
1:B:84:VAL:HG13	1:B:95:PRO:HB3	2.00	0.44
1:B:130:SER:OG	1:B:131:LEU:N	2.51	0.44
1:E:159:PHE:HA	1:E:160:GLY:HA2	1.51	0.44
1:F:159:PHE:HA	1:F:160:GLY:HA2	1.50	0.44
1:H:190:VAL:HG21	1:H:241:PHE:HD2	1.83	0.44
1:H:192:THR:HG21	1:H:238:GLY:HA2	2.00	0.44
1:I:196:ILE:HG12	1:I:230:CYS:SG	2.57	0.44
1:A:54:TYR:CE1	1:A:98:ARG:HG2	2.53	0.44
1:C:60:THR:HG23	1:C:62:GLN:N	2.32	0.44
1:D:243:ILE:O	1:D:247:ILE:HG13	2.17	0.43
1:C:242:LEU:HD23	1:C:242:LEU:HA	1.78	0.43
1:F:249:ARG:H	1:F:250:GLY:HA2	1.82	0.43
1:D:226:PRO:O	1:D:229:LEU:N	2.51	0.43
1:B:196:ILE:HG12	1:B:230:CYS:SG	2.57	0.43
1:G:225:GLU:HA	1:G:226:PRO:HD3	1.71	0.43
1:I:158:TYR:O	1:I:161:ILE:N	2.50	0.43
1:G:243:ILE:HA	1:G:252:TRP:CZ3	2.53	0.43
1:H:49:ILE:HG22	1:H:93:PHE:CD1	2.53	0.43
1:H:243:ILE:HG23	1:H:252:TRP:CH2	2.53	0.43
1:I:146:GLU:N	1:I:146:GLU:OE1	2.51	0.43
1:C:243:ILE:O	1:C:247:ILE:HG13	2.19	0.43
1:G:192:THR:HG21	1:G:238:GLY:HA2	1.99	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:178:VAL:HG23	1:H:242:LEU:HD22	2.01	0.43
1:I:170:ASP:HB2	1:I:196:ILE:HB	1.99	0.43
1:I:171:GLN:HG2	1:I:195:THR:HG22	1.99	0.43
1:A:130:SER:OG	1:A:131:LEU:N	2.50	0.43
1:B:157:ARG:HB3	1:E:216:GLU:HG3	2.01	0.43
1:F:111:ILE:HG12	1:H:133:ALA:CB	2.48	0.43
1:D:225:GLU:HA	1:D:226:PRO:HD3	1.66	0.43
1:F:91:ARG:HA	1:F:93:PHE:N	2.33	0.43
1:F:106:ASN:O	1:F:109:LYS:HB2	2.19	0.43
1:G:157:ARG:HB3	1:I:216:GLU:HG3	2.01	0.43
1:H:174:VAL:HG11	1:H:235:ILE:HG13	2.00	0.43
1:A:250:GLY:C	1:A:252:TRP:H	2.26	0.43
1:B:183:THR:HB	1:E:132:THR:HG23	2.00	0.43
1:E:145:TYR:HD2	1:E:172:ILE:HG12	1.83	0.43
1:G:151:SER:HB2	1:G:168:GLN:HG2	2.01	0.43
1:A:158:TYR:O	1:A:161:ILE:N	2.51	0.43
1:B:75:VAL:HA	1:B:76:PRO:HD3	1.92	0.43
1:C:226:PRO:O	1:C:229:LEU:N	2.51	0.43
1:G:189:SER:OG	1:I:85:THR:HG22	2.19	0.43
1:I:48:LYS:HD3	1:I:94:ILE:HD11	2.01	0.43
1:B:205:VAL:O	1:B:216:GLU:HA	2.18	0.42
1:H:172:ILE:HG13	1:H:196:ILE:HD11	1.99	0.42
1:A:96:LEU:HD22	1:A:96:LEU:HA	1.74	0.42
1:C:42:LEU:HA	1:C:43:PRO:HD3	1.91	0.42
1:E:48:LYS:HD3	1:E:94:ILE:HD11	2.01	0.42
1:C:189:SER:OG	1:G:85:THR:HG22	2.18	0.42
1:G:60:THR:CG2	1:G:62:GLN:H	2.32	0.42
1:G:76:PRO:HD2	1:G:231:LEU:HD12	2.00	0.42
1:H:53:VAL:HG22	1:H:80:THR:HG23	2.00	0.42
1:A:244:ASN:O	1:A:247:ILE:HB	2.18	0.42
1:H:226:PRO:HG2	1:H:229:LEU:HD12	2.01	0.42
1:I:50:PHE:CD1	1:I:94:ILE:HB	2.54	0.42
1:B:98:ARG:NH1	1:E:107:GLU:OE2	2.53	0.42
1:F:208:PHE:HD1	1:F:208:PHE:HA	1.68	0.42
1:H:106:ASN:O	1:H:109:LYS:HB2	2.19	0.42
1:H:249:ARG:N	1:H:250:GLY:HA2	2.35	0.42
1:D:216:GLU:CG	1:F:157:ARG:HB3	2.50	0.42
1:C:106:ASN:O	1:C:109:LYS:HB2	2.20	0.42
1:E:96:LEU:HD22	1:E:96:LEU:HA	1.66	0.42
1:F:151:SER:HB2	1:F:168:GLN:HG2	2.01	0.42
1:C:194:LYS:HE3	1:C:234:ALA:N	2.34	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:78:SER:O	1:H:81:ALA:HB3	2.20	0.42
1:I:96:LEU:HD22	1:I:96:LEU:HA	1.72	0.42
1:D:246:GLY:O	1:D:250:GLY:HA2	2.19	0.42
1:A:60:THR:HG23	1:A:62:GLN:N	2.35	0.42
1:C:96:LEU:HA	1:C:96:LEU:HD22	1.73	0.42
1:C:193:SER:O	1:C:193:SER:OG	2.38	0.42
1:F:150:LYS:HE3	1:F:167:TYR:OH	2.19	0.42
1:H:159:PHE:HA	1:H:160:GLY:HA2	1.52	0.42
1:E:150:LYS:HE3	1:E:167:TYR:OH	2.20	0.42
1:G:97:GLU:HB3	1:G:98:ARG:H	1.54	0.42
1:I:60:THR:HG23	1:I:62:GLN:N	2.35	0.42
1:A:78:SER:HB2	1:I:143:ILE:HG12	2.01	0.41
1:B:188:SER:HA	1:E:88:LYS:HE2	2.02	0.41
1:I:49:ILE:HG23	1:I:93:PHE:CD1	2.55	0.41
1:A:85:THR:HG22	1:I:189:SER:OG	2.19	0.41
1:A:146:GLU:HB2	1:A:147:SER:H	1.69	0.41
1:E:157:ARG:HB3	1:H:216:GLU:HG3	2.03	0.41
1:H:51:VAL:HG11	1:H:87:LEU:HD13	2.02	0.41
1:D:53:VAL:HG21	1:D:83:LEU:HD12	2.02	0.41
1:B:96:LEU:HA	1:B:96:LEU:HD22	1.76	0.41
1:B:177:ARG:NH1	1:E:97:GLU:OE2	2.52	0.41
1:F:97:GLU:HG2	1:H:186:ILE:HG12	2.02	0.41
1:D:107:GLU:O	1:D:111:ILE:HG13	2.20	0.41
1:A:226:PRO:O	1:A:229:LEU:N	2.52	0.41
1:B:208:PHE:HD1	1:B:208:PHE:HA	1.69	0.41
1:F:130:SER:OG	1:F:131:LEU:N	2.54	0.41
1:G:175:ASN:ND2	1:I:85:THR:HG21	2.35	0.41
1:G:249:ARG:H	1:G:250:GLY:HA2	1.86	0.41
1:I:145:TYR:OH	1:I:170:ASP:OD2	2.30	0.41
1:I:225:GLU:HA	1:I:226:PRO:HD3	1.62	0.41
1:A:111:ILE:HG12	1:I:133:ALA:HB1	2.03	0.41
1:C:91:ARG:HA	1:C:93:PHE:N	2.34	0.41
1:F:54:TYR:O	1:F:80:THR:HG21	2.20	0.41
1:A:143:ILE:CD1	1:A:175:ASN:HB2	2.50	0.41
1:A:205:VAL:O	1:A:216:GLU:HA	2.21	0.41
1:C:54:TYR:CZ	1:C:98:ARG:HG2	2.55	0.41
1:E:250:GLY:C	1:E:252:TRP:H	2.28	0.41
1:G:66:TYR:CG	1:G:67:PRO:HA	2.56	0.41
1:G:133:ALA:HB1	1:I:111:ILE:HG12	2.01	0.41
1:H:107:GLU:O	1:H:111:ILE:HG13	2.19	0.41
1:D:60:THR:HG23	1:D:62:GLN:N	2.35	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:ILE:HG12	1:I:133:ALA:CB	2.51	0.41
1:A:225:GLU:HA	1:A:226:PRO:HD3	1.73	0.41
1:B:97:GLU:HB3	1:B:98:ARG:H	1.50	0.41
1:F:56:ILE:HG13	1:F:141:SER:HA	2.03	0.41
1:F:158:TYR:O	1:F:161:ILE:N	2.53	0.41
1:G:106:ASN:O	1:G:109:LYS:HB2	2.21	0.41
1:I:249:ARG:N	1:I:250:GLY:HA2	2.36	0.41
1:D:97:GLU:HB3	1:D:98:ARG:H	1.54	0.41
1:F:58:ASP:HB3	1:F:77:GLN:NE2	2.36	0.41
1:D:208:PHE:HD1	1:D:208:PHE:HA	1.69	0.41
1:A:189:SER:OG	1:B:85:THR:HG22	2.21	0.41
1:C:137:MET:HE3	1:C:137:MET:HB2	1.78	0.41
1:C:177:ARG:NH1	1:G:97:GLU:OE2	2.54	0.41
1:E:75:VAL:HA	1:E:76:PRO:HD3	1.95	0.41
1:G:96:LEU:HD22	1:G:96:LEU:HA	1.76	0.41
1:G:158:TYR:O	1:G:161:ILE:N	2.54	0.41
1:F:110:ILE:HD13	1:H:98:ARG:CZ	2.51	0.41
1:H:198:SER:HB3	1:H:224:ASN:ND2	2.36	0.41
1:D:106:ASN:O	1:D:109:LYS:HB2	2.21	0.40
1:C:158:TYR:O	1:C:161:ILE:N	2.55	0.40
1:C:98:ARG:NH1	1:G:107:GLU:OE2	2.54	0.40
1:E:91:ARG:HA	1:E:93:PHE:N	2.36	0.40
1:A:208:PHE:HD1	1:A:208:PHE:HA	1.68	0.40
1:G:91:ARG:HA	1:G:93:PHE:N	2.36	0.40
1:G:146:GLU:HG3	1:G:149:VAL:CG2	2.51	0.40
1:E:243:ILE:O	1:E:247:ILE:HG13	2.21	0.40
1:G:208:PHE:HD1	1:G:208:PHE:HA	1.68	0.40
1:A:106:ASN:O	1:A:109:LYS:HB2	2.21	0.40
1:A:150:LYS:HD3	1:A:169:LEU:HD12	2.03	0.40
1:F:145:TYR:HA	1:F:171:GLN:O	2.22	0.40
1:H:130:SER:OG	1:H:131:LEU:N	2.55	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:244:ASN:OD1	1:E:244:ASN:ND2[4_555]	2.06	0.14

5.3 Torsion angles

5.3.1 Protein backbone

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	190/277 (69%)	176 (93%)	12 (6%)	2 (1%)	11	36
1	B	184/277 (66%)	175 (95%)	8 (4%)	1 (0%)	24	54
1	C	192/277 (69%)	176 (92%)	15 (8%)	1 (0%)	24	54
1	D	183/277 (66%)	173 (94%)	9 (5%)	1 (0%)	24	54
1	E	192/277 (69%)	176 (92%)	15 (8%)	1 (0%)	24	54
1	F	183/277 (66%)	173 (94%)	9 (5%)	1 (0%)	24	54
1	G	184/277 (66%)	173 (94%)	10 (5%)	1 (0%)	24	54
1	H	185/277 (67%)	174 (94%)	10 (5%)	1 (0%)	24	54
1	I	184/277 (66%)	175 (95%)	8 (4%)	1 (0%)	24	54
All	All	1677/2493 (67%)	1571 (94%)	96 (6%)	10 (1%)	21	51

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	149	VAL
1	A	149	VAL
1	B	149	VAL
1	C	149	VAL
1	E	149	VAL
1	F	149	VAL
1	G	149	VAL
1	H	149	VAL
1	I	149	VAL
1	A	251	LEU

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	157/236 (66%)	130 (83%)	27 (17%)	2	7
1	B	154/236 (65%)	126 (82%)	28 (18%)	2	6
1	C	161/236 (68%)	129 (80%)	32 (20%)	1	4
1	D	153/236 (65%)	126 (82%)	27 (18%)	2	6
1	E	161/236 (68%)	131 (81%)	30 (19%)	1	5
1	F	153/236 (65%)	125 (82%)	28 (18%)	2	6
1	G	155/236 (66%)	126 (81%)	29 (19%)	1	5
1	H	155/236 (66%)	128 (83%)	27 (17%)	2	7
1	I	153/236 (65%)	126 (82%)	27 (18%)	2	6
All	All	1402/2124 (66%)	1147 (82%)	255 (18%)	2	6

All (255) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	49	ILE
1	D	53	VAL
1	D	56	ILE
1	D	60	THR
1	D	62	GLN
1	D	64	LYS
1	D	75	VAL
1	D	77	GLN
1	D	80	THR
1	D	83	LEU
1	D	92	TRP
1	D	96	LEU
1	D	101	LEU
1	D	132	THR
1	D	139	GLU
1	D	141	SER
1	D	143	ILE
1	D	151	SER
1	D	176	LEU
1	D	177	ARG
1	D	179	VAL
1	D	208	PHE
1	D	228	MET

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	231	LEU
1	D	235	ILE
1	D	237	THR
1	D	245	ASP
1	A	49	ILE
1	A	53	VAL
1	A	56	ILE
1	A	60	THR
1	A	62	GLN
1	A	64	LYS
1	A	75	VAL
1	A	77	GLN
1	A	80	THR
1	A	83	LEU
1	A	96	LEU
1	A	101	LEU
1	A	102	GLN
1	A	132	THR
1	A	139	GLU
1	A	141	SER
1	A	143	ILE
1	A	151	SER
1	A	176	LEU
1	A	177	ARG
1	A	179	VAL
1	A	198	SER
1	A	208	PHE
1	A	228	MET
1	A	235	ILE
1	A	248	ASP
1	A	251	LEU
1	B	49	ILE
1	B	53	VAL
1	B	56	ILE
1	B	60	THR
1	B	62	GLN
1	B	64	LYS
1	B	75	VAL
1	B	77	GLN
1	B	78	SER
1	B	80	THR
1	B	83	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	96	LEU
1	B	101	LEU
1	B	136	ILE
1	B	139	GLU
1	B	141	SER
1	B	143	ILE
1	B	146	GLU
1	B	151	SER
1	B	176	LEU
1	B	177	ARG
1	B	179	VAL
1	B	208	PHE
1	B	228	MET
1	B	231	LEU
1	B	235	ILE
1	B	237	THR
1	B	251	LEU
1	C	41	HIS
1	C	42	LEU
1	C	49	ILE
1	C	53	VAL
1	C	56	ILE
1	C	60	THR
1	C	62	GLN
1	C	64	LYS
1	C	77	GLN
1	C	78	SER
1	C	80	THR
1	C	83	LEU
1	C	96	LEU
1	C	101	LEU
1	C	102	GLN
1	C	139	GLU
1	C	141	SER
1	C	143	ILE
1	C	146	GLU
1	C	151	SER
1	C	176	LEU
1	C	177	ARG
1	C	179	VAL
1	C	198	SER
1	C	208	PHE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	228	MET
1	C	231	LEU
1	C	235	ILE
1	C	244	ASN
1	C	248	ASP
1	C	251	LEU
1	C	252	TRP
1	E	41	HIS
1	E	42	LEU
1	E	49	ILE
1	E	53	VAL
1	E	56	ILE
1	E	60	THR
1	E	62	GLN
1	E	64	LYS
1	E	75	VAL
1	E	77	GLN
1	E	80	THR
1	E	83	LEU
1	E	92	TRP
1	E	96	LEU
1	E	101	LEU
1	E	139	GLU
1	E	141	SER
1	E	143	ILE
1	E	146	GLU
1	E	151	SER
1	E	176	LEU
1	E	177	ARG
1	E	179	VAL
1	E	190	VAL
1	E	208	PHE
1	E	228	MET
1	E	231	LEU
1	E	235	ILE
1	E	245	ASP
1	E	251	LEU
1	F	49	ILE
1	F	53	VAL
1	F	56	ILE
1	F	60	THR
1	F	62	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	F	64	LYS
1	F	77	GLN
1	F	78	SER
1	F	80	THR
1	F	83	LEU
1	F	96	LEU
1	F	98	ARG
1	F	101	LEU
1	F	102	GLN
1	F	136	ILE
1	F	139	GLU
1	F	141	SER
1	F	143	ILE
1	F	146	GLU
1	F	151	SER
1	F	176	LEU
1	F	177	ARG
1	F	179	VAL
1	F	198	SER
1	F	208	PHE
1	F	228	MET
1	F	235	ILE
1	F	244	ASN
1	G	49	ILE
1	G	53	VAL
1	G	56	ILE
1	G	60	THR
1	G	62	GLN
1	G	64	LYS
1	G	75	VAL
1	G	77	GLN
1	G	80	THR
1	G	83	LEU
1	G	96	LEU
1	G	101	LEU
1	G	102	GLN
1	G	135	ASN
1	G	139	GLU
1	G	141	SER
1	G	143	ILE
1	G	151	SER
1	G	176	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	G	177	ARG
1	G	179	VAL
1	G	198	SER
1	G	208	PHE
1	G	214	LEU
1	G	228	MET
1	G	231	LEU
1	G	235	ILE
1	G	248	ASP
1	G	251	LEU
1	H	49	ILE
1	H	53	VAL
1	H	56	ILE
1	H	60	THR
1	H	62	GLN
1	H	64	LYS
1	H	75	VAL
1	H	77	GLN
1	H	80	THR
1	H	83	LEU
1	H	96	LEU
1	H	101	LEU
1	H	102	GLN
1	H	139	GLU
1	H	141	SER
1	H	143	ILE
1	H	151	SER
1	H	176	LEU
1	H	177	ARG
1	H	179	VAL
1	H	198	SER
1	H	208	PHE
1	H	228	MET
1	H	231	LEU
1	H	235	ILE
1	H	248	ASP
1	H	251	LEU
1	I	53	VAL
1	I	56	ILE
1	I	60	THR
1	I	62	GLN
1	I	64	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	I	75	VAL
1	I	77	GLN
1	I	80	THR
1	I	83	LEU
1	I	96	LEU
1	I	101	LEU
1	I	102	GLN
1	I	132	THR
1	I	136	ILE
1	I	139	GLU
1	I	141	SER
1	I	143	ILE
1	I	151	SER
1	I	176	LEU
1	I	177	ARG
1	I	179	VAL
1	I	208	PHE
1	I	228	MET
1	I	231	LEU
1	I	235	ILE
1	I	248	ASP
1	I	251	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	GLN
1	A	99	GLN
1	A	103	ASN
1	C	62	GLN
1	C	103	ASN
1	F	62	GLN
1	F	103	ASN
1	H	57	GLN
1	H	224	ASN
1	I	77	GLN

5.3.3 RNA ⓘ

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 [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	196/277 (70%)	0.82	22 (11%) 10 9	49, 76, 153, 185	0
1	B	190/277 (68%)	0.69	17 (8%) 15 13	46, 76, 158, 191	0
1	C	198/277 (71%)	0.75	27 (13%) 7 5	47, 70, 156, 199	0
1	D	189/277 (68%)	0.82	27 (14%) 6 5	48, 74, 149, 188	0
1	E	198/277 (71%)	0.81	24 (12%) 8 7	47, 73, 156, 191	0
1	F	189/277 (68%)	0.63	17 (8%) 15 13	50, 77, 145, 192	0
1	G	190/277 (68%)	0.95	32 (16%) 4 3	50, 78, 168, 187	0
1	H	191/277 (68%)	0.76	18 (9%) 14 12	47, 79, 157, 188	0
1	I	190/277 (68%)	0.86	26 (13%) 6 5	45, 77, 157, 186	0
All	All	1731/2493 (69%)	0.79	210 (12%) 8 7	45, 76, 157, 199	0

All (210) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	45	PRO	5.7
1	G	115	GLN	5.6
1	I	49	ILE	5.5
1	G	110	ILE	5.5
1	A	43	PRO	5.1
1	E	92	TRP	4.8
1	D	93	PHE	4.5
1	C	148	ASN	4.5
1	C	44	ALA	4.4
1	A	148	ASN	4.3
1	I	135	ASN	4.3
1	G	111	ILE	4.3
1	E	45	PRO	4.3
1	G	251	LEU	4.1
1	F	210	ASP	3.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	H	148	ASN	3.9
1	D	49	ILE	3.8
1	E	251	LEU	3.8
1	A	44	ALA	3.8
1	B	251	LEU	3.7
1	C	247	ILE	3.7
1	H	251	LEU	3.7
1	E	55	ASN	3.7
1	E	252	TRP	3.7
1	F	92	TRP	3.7
1	C	251	LEU	3.6
1	I	129	GLN	3.6
1	E	150	LYS	3.6
1	A	251	LEU	3.6
1	B	55	ASN	3.6
1	D	146	GLU	3.5
1	H	129	GLN	3.5
1	B	247	ILE	3.5
1	F	80	THR	3.5
1	G	129	GLN	3.4
1	I	206	PHE	3.3
1	A	147	SER	3.3
1	E	44	ALA	3.3
1	G	146	GLU	3.3
1	F	209	ILE	3.3
1	I	251	LEU	3.3
1	G	50	PHE	3.2
1	F	116	GLU	3.2
1	G	92	TRP	3.2
1	G	215	LEU	3.2
1	A	213	ARG	3.2
1	I	55	ASN	3.1
1	D	50	PHE	3.1
1	A	85	THR	3.0
1	F	150	LYS	3.0
1	G	49	ILE	3.0
1	I	146	GLU	3.0
1	I	205	VAL	3.0
1	C	54	TYR	3.0
1	I	48	LYS	3.0
1	A	97	GLU	3.0
1	D	92	TRP	2.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	148	ASN	2.9
1	B	98	ARG	2.9
1	B	166	GLN	2.9
1	C	158	TYR	2.9
1	A	89	ASP	2.9
1	G	113	ALA	2.9
1	C	133	ALA	2.9
1	G	78	SER	2.9
1	D	209	ILE	2.9
1	I	51	VAL	2.8
1	E	78	SER	2.8
1	D	213	ARG	2.8
1	G	252	TRP	2.8
1	G	143	ILE	2.8
1	C	135	ASN	2.8
1	D	78	SER	2.8
1	D	205	VAL	2.7
1	A	215	LEU	2.7
1	D	97	GLU	2.7
1	D	214	LEU	2.7
1	A	206	PHE	2.7
1	F	147	SER	2.7
1	B	248	ASP	2.7
1	A	150	LYS	2.7
1	I	213	ARG	2.6
1	B	209	ILE	2.6
1	D	147	SER	2.6
1	G	147	SER	2.6
1	E	158	TYR	2.6
1	D	70	ASN	2.6
1	A	232	MET	2.6
1	C	99	GLN	2.6
1	D	90	SER	2.6
1	D	98	ARG	2.5
1	C	207	ARG	2.5
1	F	91	ARG	2.5
1	C	92	TRP	2.5
1	H	252	TRP	2.5
1	E	215	LEU	2.5
1	E	79	ALA	2.5
1	F	159	PHE	2.5
1	G	210	ASP	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	216	GLU	2.5
1	D	156	ALA	2.5
1	I	219	VAL	2.5
1	I	116	GLU	2.5
1	A	168	GLN	2.5
1	D	244	ASN	2.5
1	A	209	ILE	2.4
1	A	98	ARG	2.4
1	C	143	ILE	2.4
1	A	48	LYS	2.4
1	F	81	ALA	2.4
1	I	57	GLN	2.4
1	D	132	THR	2.4
1	C	42	LEU	2.4
1	F	131	LEU	2.4
1	H	229	LEU	2.4
1	H	55	ASN	2.4
1	E	43	PRO	2.4
1	E	236	GLU	2.4
1	G	197	LEU	2.4
1	B	249	ARG	2.4
1	C	98	ARG	2.4
1	I	159	PHE	2.4
1	I	209	ILE	2.4
1	C	43	PRO	2.4
1	C	55	ASN	2.3
1	I	98	ARG	2.3
1	D	159	PHE	2.3
1	D	210	ASP	2.3
1	C	97	GLU	2.3
1	B	57	GLN	2.3
1	F	213	ARG	2.3
1	H	135	ASN	2.3
1	C	206	PHE	2.3
1	E	99	GLN	2.3
1	F	129	GLN	2.3
1	G	57	GLN	2.3
1	E	188	SER	2.3
1	B	244	ASN	2.3
1	E	249	ARG	2.3
1	G	112	ARG	2.3
1	D	206	PHE	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	G	159	PHE	2.3
1	C	114	ALA	2.3
1	F	215	LEU	2.3
1	D	85	THR	2.3
1	E	54	TYR	2.3
1	G	97	GLU	2.2
1	B	215	LEU	2.2
1	H	57	GLN	2.2
1	I	207	ARG	2.2
1	G	154	VAL	2.2
1	I	217	GLY	2.2
1	C	209	ILE	2.2
1	B	210	ASP	2.2
1	I	93	PHE	2.2
1	D	229	LEU	2.2
1	E	74	ALA	2.2
1	H	131	LEU	2.2
1	B	243	ILE	2.2
1	I	94	ILE	2.2
1	G	158	TYR	2.2
1	G	206	PHE	2.2
1	A	229	LEU	2.2
1	D	110	ILE	2.2
1	B	205	VAL	2.2
1	D	100	GLY	2.2
1	I	100	GLY	2.2
1	E	167	TYR	2.2
1	I	167	TYR	2.2
1	E	147	SER	2.2
1	H	215	LEU	2.2
1	C	111	ILE	2.1
1	D	57	GLN	2.1
1	F	70	ASN	2.1
1	G	55	ASN	2.1
1	C	205	VAL	2.1
1	G	205	VAL	2.1
1	C	250	GLY	2.1
1	G	132	THR	2.1
1	C	252	TRP	2.1
1	G	81	ALA	2.1
1	H	97	GLU	2.1
1	A	55	ASN	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	214	LEU	2.1
1	F	98	ARG	2.1
1	H	91	ARG	2.1
1	H	221	TYR	2.1
1	H	72	SER	2.1
1	H	200	GLU	2.1
1	C	101	LEU	2.1
1	G	153	GLY	2.1
1	E	46	THR	2.1
1	E	209	ILE	2.1
1	C	166	GLN	2.1
1	G	224	ASN	2.1
1	H	242	LEU	2.1
1	I	96	LEU	2.1
1	D	208	PHE	2.1
1	B	245	ASP	2.1
1	H	213	ARG	2.1
1	G	166	GLN	2.0
1	A	205	VAL	2.0
1	A	83	LEU	2.0
1	F	172	ILE	2.0
1	H	163	ALA	2.0
1	E	129	GLN	2.0
1	I	95	PRO	2.0
1	B	79	ALA	2.0
1	G	156	ALA	2.0
1	I	97	GLU	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.