Full wwPDB X-ray Structure Validation Report (i)

Dec 17, 2023 – 11:58 am GMT

PDB ID	:	2X89
Title	:	Structure of the Beta2_microglobulin involved in amyloidogenesis
Authors	:	Domanska, K.; Srinivasan, V.; Vanderhaegen, S.; Pardon, E.; Marquez, J.A.; Bellotti, V.; Wyns, L.; Steyaert, J.
Deposited on	:	2010-03-07
Resolution	:	2.16 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 (i) 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 (1)) were used in the production of this report:

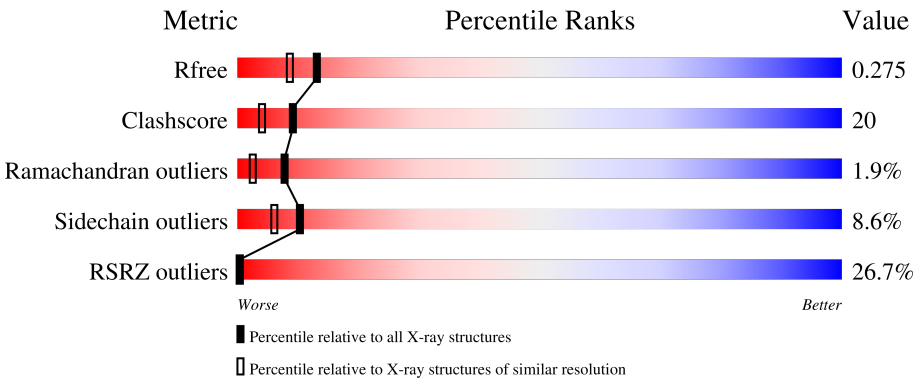
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.16 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}	130704	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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 <=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	128	23% 76% 20% • •
1	B	128	24% 67% 23% 9% •
1	C	128	22% 66% 23% 9% •
2	D	94	26% 60% 33% • • •
2	E	94	26% 74% 20% • •

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Mol	Chain	Length	Quality of chain
2	F	94	24% 59% 29% 7%
2	G	94	43% 51% 29% 11% 9%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 6225 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 ANTIBODY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	128	986	607	177	196	6	0	0	0
1	B	128	993	613	177	197	6	0	0	0
1	C	128	993	613	177	197	6	0	0	0

- Molecule 2 is a protein called BETA-2-MICROGLOBULIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	D	92	755	483	126	143	3	0	0	0
2	E	91	738	471	122	142	3	0	0	0
2	F	91	758	486	124	145	3	0	0	0
2	G	86	707	453	118	133	3	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	6	MET	-	expression tag	UNP P61769
E	6	MET	-	expression tag	UNP P61769
F	6	MET	-	expression tag	UNP P61769
G	6	MET	-	expression tag	UNP P61769

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	44	44	44	0	0

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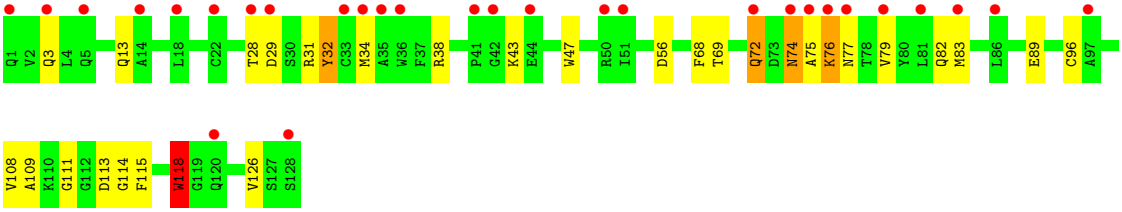
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	B	62	62	62	0	0
			Total	O		
3	C	43	43	43	0	0
			Total	O		
3	D	25	25	25	0	0
			Total	O		
3	E	44	44	44	0	0
			Total	O		
3	F	31	31	31	0	0
			Total	O		
3	G	46	46	46	0	0

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.

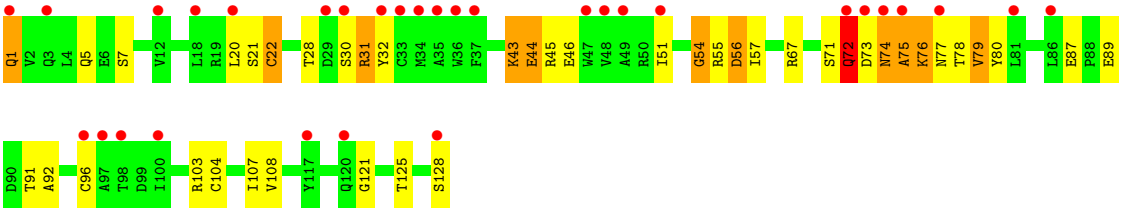
- Molecule 1: ANTIBODY

Chain A:



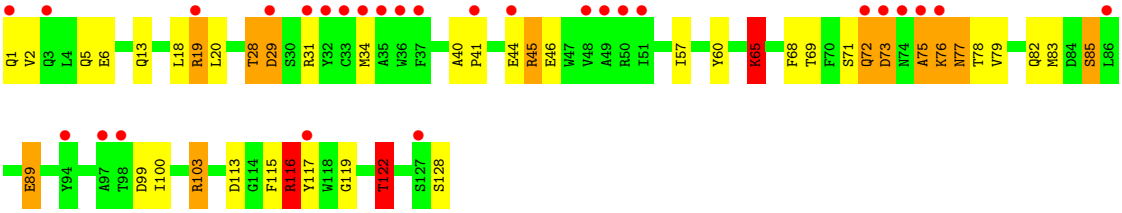
- Molecule 1: ANTIBODY

Chain B:



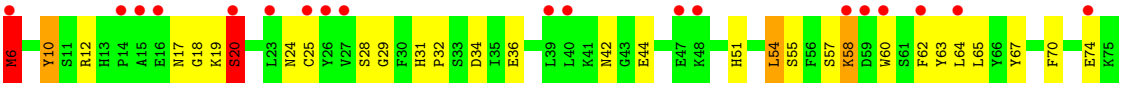
- Molecule 1: ANTIBODY

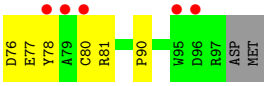
Chain C:



- Molecule 2: BETA-2-MICROGLOBULIN

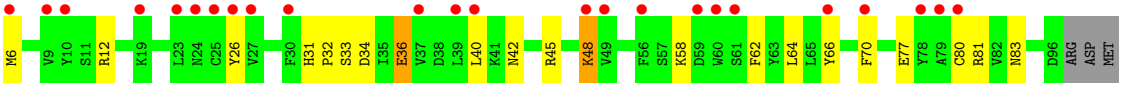
Chain D:





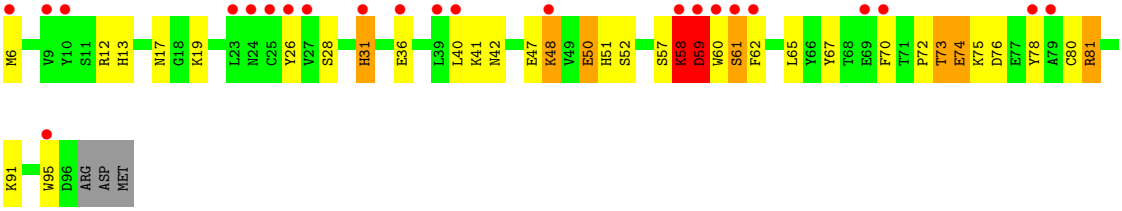
• Molecule 2: BETA-2-MICROGLOBULIN

Chain E:



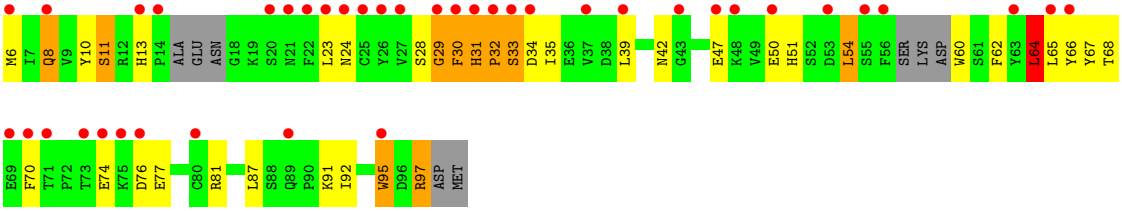
• Molecule 2: BETA-2-MICROGLOBULIN

Chain F:



• Molecule 2: BETA-2-MICROGLOBULIN

Chain G:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	77.49Å 100.86Å 83.74Å 90.00° 106.43° 90.00°	Depositor
Resolution (Å)	19.85 – 2.16 19.83 – 2.16	Depositor EDS
% Data completeness (in resolution range)	99.9 (19.85-2.16) 99.9 (19.83-2.16)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.04 (at 2.17Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.238 , 0.267 0.254 , 0.275	Depositor DCC
R_{free} test set	3356 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	48.1	Xtriage
Anisotropy	0.029	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 52.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	6225	wwPDB-VP
Average B, all atoms (Å^2)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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	1.15	4/1005 (0.4%)	1.03	4/1358 (0.3%)
1	B	1.43	13/1013 (1.3%)	1.15	4/1369 (0.3%)
1	C	1.45	11/1013 (1.1%)	1.33	12/1369 (0.9%)
2	D	1.18	5/777 (0.6%)	1.01	3/1052 (0.3%)
2	E	1.31	3/758 (0.4%)	1.11	2/1026 (0.2%)
2	F	1.28	1/780 (0.1%)	1.13	3/1057 (0.3%)
2	G	1.26	4/725 (0.6%)	1.21	9/980 (0.9%)
All	All	1.30	41/6071 (0.7%)	1.15	37/8211 (0.5%)

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	2
1	B	0	4
1	C	0	2
2	D	0	1
2	E	0	1
2	F	0	2
2	G	0	2
All	All	0	14

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	36	GLU	CG-CD	10.19	1.67	1.51
2	G	95	TRP	CE3-CZ3	-8.61	1.23	1.38
2	F	80	CYS	CB-SG	-8.56	1.67	1.82
2	D	80	CYS	CB-SG	-8.26	1.68	1.82
1	C	89	GLU	CB-CG	8.06	1.67	1.52
1	C	89	GLU	CD-OE2	7.76	1.34	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	89	GLU	CG-CD	7.60	1.63	1.51
1	B	55	ARG	C-N	-7.34	1.17	1.34
2	E	36	GLU	CD-OE1	6.91	1.33	1.25
2	G	30	PHE	N-CA	6.86	1.60	1.46
1	B	89	GLU	CG-CD	6.75	1.62	1.51
1	B	22	CYS	CB-SG	-6.67	1.71	1.82
2	E	80	CYS	CB-SG	-6.54	1.71	1.82
2	D	25	CYS	CB-SG	-6.50	1.71	1.82
1	B	104	CYS	CB-SG	-6.32	1.71	1.82
1	B	96	CYS	CA-CB	6.14	1.67	1.53
1	B	87	GLU	CG-CD	6.06	1.61	1.51
1	A	96	CYS	CB-SG	-5.92	1.72	1.81
1	C	122	THR	CB-CG2	-5.91	1.32	1.52
1	B	21	SER	CB-OG	-5.90	1.34	1.42
1	C	85	SER	CB-OG	-5.88	1.34	1.42
1	C	103	ARG	CG-CD	-5.78	1.37	1.51
1	B	54	GLY	C-N	-5.75	1.20	1.34
1	B	79	VAL	CB-CG2	5.73	1.64	1.52
1	C	89	GLU	CD-OE1	5.61	1.31	1.25
1	B	44	GLU	CG-CD	5.57	1.60	1.51
1	A	126	VAL	CB-CG2	5.52	1.64	1.52
2	D	44	GLU	CD-OE2	-5.44	1.19	1.25
1	B	46	GLU	CD-OE1	5.35	1.31	1.25
1	C	60	TYR	CE2-CZ	5.30	1.45	1.38
1	A	32	TYR	CD2-CE2	5.29	1.47	1.39
2	G	29	GLY	CA-C	5.25	1.60	1.51
2	D	10	TYR	CE2-CZ	-5.24	1.31	1.38
1	C	85	SER	CA-CB	5.22	1.60	1.52
1	B	87	GLU	CD-OE1	5.20	1.31	1.25
1	C	115	PHE	CE1-CZ	5.19	1.47	1.37
1	C	5	GLN	CG-CD	5.15	1.62	1.51
2	G	33	SER	N-CA	5.13	1.56	1.46
1	A	118	TRP	CE3-CZ3	-5.10	1.29	1.38
1	B	51	ILE	CB-CG2	5.04	1.68	1.52
2	D	78	TYR	CD2-CE2	5.02	1.46	1.39

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	103	ARG	NE-CZ-NH1	-10.47	115.06	120.30
1	C	65	LYS	CD-CE-NZ	-9.60	89.63	111.70
1	C	85	SER	CB-CA-C	9.55	128.25	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	56	ASP	CB-CG-OD1	9.29	126.66	118.30
1	C	19	ARG	CG-CD-NE	7.74	128.05	111.80
1	B	96	CYS	CA-CB-SG	-7.68	100.18	114.00
2	G	64	LEU	CA-CB-CG	7.46	132.45	115.30
2	E	81	ARG	NE-CZ-NH2	-7.31	116.64	120.30
1	C	29	ASP	CB-CG-OD1	-7.24	111.78	118.30
2	F	81	ARG	NE-CZ-NH2	-7.15	116.72	120.30
1	C	45	ARG	NE-CZ-NH2	-7.03	116.78	120.30
2	G	29	GLY	N-CA-C	7.00	130.59	113.10
1	B	72	GLN	CB-CA-C	-6.88	96.64	110.40
2	E	81	ARG	NE-CZ-NH1	6.85	123.72	120.30
2	G	30	PHE	N-CA-C	6.45	128.41	111.00
1	B	56	ASP	CB-CG-OD1	6.40	124.06	118.30
2	G	81	ARG	NE-CZ-NH2	-6.35	117.12	120.30
2	F	81	ARG	NE-CZ-NH1	6.33	123.47	120.30
1	A	38	ARG	NE-CZ-NH2	6.33	123.46	120.30
2	D	6	MET	CG-SD-CE	6.29	110.27	100.20
2	G	64	LEU	CB-CG-CD2	-6.14	100.55	111.00
2	F	59	ASP	N-CA-C	6.04	127.31	111.00
1	A	56	ASP	CB-CG-OD2	-5.83	113.06	118.30
1	C	103	ARG	CG-CD-NE	-5.80	99.62	111.80
1	C	116	ARG	NE-CZ-NH2	-5.63	117.48	120.30
2	G	33	SER	CA-C-N	-5.63	104.82	117.20
1	C	19	ARG	CB-CG-CD	-5.59	97.05	111.60
2	G	32	PRO	N-CA-C	5.50	126.40	112.10
1	A	83	MET	CG-SD-CE	5.42	108.87	100.20
2	G	34	ASP	CB-CG-OD2	-5.41	113.44	118.30
1	C	19	ARG	NE-CZ-NH1	-5.39	117.61	120.30
1	C	19	ARG	NE-CZ-NH2	5.36	122.98	120.30
1	C	103	ARG	NH1-CZ-NH2	5.31	125.24	119.40
2	D	81	ARG	NE-CZ-NH2	-5.20	117.70	120.30
2	G	95	TRP	CA-CB-CG	5.15	123.48	113.70
2	D	54	LEU	CB-CG-CD2	-5.12	102.30	111.00
1	B	67	ARG	NE-CZ-NH2	5.08	122.84	120.30

There are no chirality outliers.

All (14) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	74	ASN	Peptide
1	A	75	ALA	Peptide
1	B	7	SER	Peptide

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Mol	Chain	Res	Type	Group
1	B	71	SER	Peptide
1	B	75	ALA	Peptide
1	B	76	LYS	Peptide
1	C	119	GLY	Peptide
1	C	76	LYS	Peptide
2	D	20	SER	Peptide
2	E	48	LYS	Peptide
2	F	58	LYS	Peptide
2	F	61	SER	Peptide
2	G	31	HIS	Peptide
2	G	33	SER	Peptide

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	986	0	928	39	0
1	B	993	0	933	38	0
1	C	993	0	935	61	0
2	D	755	0	703	28	0
2	E	738	0	688	12	0
2	F	758	0	715	43	0
2	G	707	0	656	45	0
3	A	44	0	0	4	0
3	B	62	0	0	1	0
3	C	43	0	0	8	0
3	D	25	0	0	2	0
3	E	44	0	0	3	0
3	F	31	0	0	2	0
3	G	46	0	0	3	0
All	All	6225	0	5558	230	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 20.

All (230) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:59:ASP:C	2:F:61:SER:HA	1.28	1.54
1:C:72:GLN:OE1	1:C:79:VAL:HG22	1.27	1.27
1:A:72:GLN:NE2	1:A:79:VAL:HG22	1.63	1.13
1:C:103:ARG:NH2	2:G:8:GLN:HE21	1.48	1.10
1:C:28:THR:HG22	2:G:6:MET:SD	1.93	1.08
2:F:59:ASP:C	2:F:61:SER:CA	2.25	1.05
2:F:6:MET:HG3	2:F:28:SER:O	1.63	0.99
1:C:103:ARG:NH2	2:G:8:GLN:NE2	2.11	0.99
1:A:32:TYR:CE2	1:A:77:ASN:ND2	2.32	0.97
1:C:103:ARG:HH21	2:G:8:GLN:NE2	1.62	0.97
1:C:75:ALA:HB1	1:C:76:LYS:CA	1.95	0.97
1:B:30:SER:HB3	1:B:77:ASN:HD21	1.28	0.96
1:C:99:ASP:OD1	1:C:103:ARG:NH1	1.98	0.96
2:D:31:HIS:O	2:D:34:ASP:HB2	1.65	0.96
1:B:45:ARG:HD3	3:B:2022:HOH:O	1.66	0.96
1:A:72:GLN:OE1	1:A:79:VAL:HG13	1.64	0.96
2:F:60:TRP:N	2:F:61:SER:HA	1.81	0.95
2:G:32:PRO:HA	2:G:35:ILE:HG12	1.49	0.94
2:G:97:ARG:CG	2:G:97:ARG:HH11	1.81	0.94
1:C:28:THR:CG2	2:G:6:MET:SD	2.56	0.93
1:C:75:ALA:HB1	1:C:76:LYS:HA	1.50	0.93
1:C:75:ALA:CB	1:C:76:LYS:HA	1.99	0.93
1:C:72:GLN:OE1	1:C:79:VAL:CG2	2.17	0.92
1:A:32:TYR:HE2	1:A:77:ASN:ND2	1.66	0.92
2:F:59:ASP:CA	2:F:61:SER:HA	2.05	0.86
2:E:36:GLU:HG2	3:E:2017:HOH:O	1.76	0.86
2:F:59:ASP:O	2:F:61:SER:HA	1.76	0.85
1:B:30:SER:CB	1:B:77:ASN:HD21	1.90	0.84
2:E:12:ARG:HB2	3:E:2007:HOH:O	1.79	0.83
2:F:48:LYS:H	2:F:48:LYS:HD2	1.44	0.82
2:G:30:PHE:N	2:G:62:PHE:CZ	2.47	0.82
1:C:75:ALA:HB1	1:C:76:LYS:C	1.99	0.82
1:A:72:GLN:HE22	1:A:79:VAL:HG22	1.40	0.82
1:A:34:MET:HE3	1:A:79:VAL:CG1	2.11	0.81
2:G:29:GLY:C	2:G:62:PHE:CZ	2.54	0.81
2:E:31:HIS:HB2	2:E:34:ASP:OD2	1.81	0.81
1:C:103:ARG:HH21	2:G:8:GLN:HE21	0.81	0.81
1:A:34:MET:CE	1:A:79:VAL:HG11	2.12	0.80
1:B:43:LYS:HA	1:B:43:LYS:HE3	1.65	0.79
1:A:113:ASP:O	1:A:118:TRP:HH2	1.65	0.79
1:A:72:GLN:OE1	1:A:79:VAL:CG1	2.32	0.78
2:G:30:PHE:N	2:G:62:PHE:HZ	1.83	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:59:ASP:HA	2:F:61:SER:HB3	1.67	0.76
1:A:32:TYR:HE2	1:A:77:ASN:HD21	0.84	0.76
2:G:54:LEU:CD1	2:G:64:LEU:HG	2.16	0.76
1:A:77:ASN:HA	3:A:2031:HOH:O	1.86	0.75
1:C:72:GLN:HG2	3:C:2025:HOH:O	1.87	0.75
1:C:75:ALA:CB	1:C:76:LYS:CA	2.60	0.74
2:F:59:ASP:HA	2:F:61:SER:CB	2.17	0.74
1:C:89:GLU:CG	3:C:2031:HOH:O	2.36	0.74
1:B:76:LYS:O	1:B:76:LYS:HD3	1.86	0.74
1:C:73:ASP:O	1:C:75:ALA:O	2.04	0.73
2:F:59:ASP:HA	2:F:61:SER:HA	1.69	0.73
2:G:97:ARG:HH11	2:G:97:ARG:HG2	1.54	0.73
2:F:59:ASP:HA	2:F:61:SER:CA	2.19	0.72
2:F:73:THR:HG21	2:F:76:ASP:OD2	1.89	0.72
2:D:19:LYS:O	2:D:20:SER:HB2	1.89	0.72
1:C:28:THR:HB	2:G:6:MET:SD	2.30	0.72
2:G:42:ASN:ND2	2:G:77:GLU:H	1.89	0.71
1:B:30:SER:HB3	1:B:77:ASN:ND2	2.02	0.71
2:D:36:GLU:OE1	2:E:33:SER:HB2	1.91	0.71
1:C:89:GLU:HG3	3:C:2031:HOH:O	1.91	0.71
2:F:17:ASN:OD1	2:G:97:ARG:NH2	2.23	0.71
2:G:97:ARG:HH11	2:G:97:ARG:HG3	1.54	0.71
1:A:34:MET:HE3	1:A:79:VAL:HG11	1.72	0.70
1:A:72:GLN:HB3	3:A:2030:HOH:O	1.91	0.70
1:C:69:THR:HG22	3:C:2006:HOH:O	1.91	0.70
2:G:11:SER:OG	2:G:13:HIS:O	2.10	0.70
1:A:72:GLN:NE2	1:A:79:VAL:CG2	2.50	0.69
1:A:32:TYR:OH	1:A:77:ASN:ND2	2.26	0.69
2:F:6:MET:CG	2:F:28:SER:O	2.38	0.69
1:A:32:TYR:CZ	1:A:77:ASN:ND2	2.61	0.68
1:C:28:THR:HG21	2:G:28:SER:OG	1.95	0.67
1:B:1:GLN:N	1:B:1:GLN:CD	2.47	0.67
2:D:19:LYS:O	2:D:20:SER:CB	2.42	0.66
1:A:72:GLN:HE22	1:A:79:VAL:CG2	2.08	0.66
2:D:32:PRO:HD3	2:D:64:LEU:HD11	1.77	0.66
2:F:73:THR:HG22	2:F:76:ASP:H	1.60	0.66
1:A:72:GLN:NE2	1:A:72:GLN:HA	2.08	0.66
1:A:34:MET:CE	1:A:79:VAL:CG1	2.73	0.65
1:C:28:THR:CG2	2:G:28:SER:OG	2.45	0.65
1:B:31:ARG:HD2	2:D:58:LYS:HZ2	1.62	0.65
1:C:89:GLU:CB	3:C:2031:HOH:O	2.45	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:48:LYS:H	2:F:48:LYS:CD	2.10	0.64
1:A:72:GLN:CD	1:A:79:VAL:HG13	2.18	0.64
2:D:24:ASN:HD22	2:D:67:TYR:HB3	1.63	0.63
2:F:60:TRP:N	2:F:61:SER:CA	2.56	0.63
2:F:78:TYR:HB2	2:G:95:TRP:CD1	2.34	0.63
1:B:31:ARG:CZ	2:D:57:SER:HB2	2.29	0.62
2:F:58:LYS:O	2:F:59:ASP:HB2	1.99	0.62
2:D:51:HIS:HD2	3:D:2015:HOH:O	1.81	0.62
1:C:72:GLN:CG	1:C:73:ASP:O	2.49	0.61
2:F:48:LYS:HD2	2:F:48:LYS:N	2.12	0.60
2:G:29:GLY:C	2:G:62:PHE:HZ	2.01	0.60
1:C:2:VAL:HB	1:C:117:TYR:CE2	2.36	0.60
2:F:19:LYS:O	2:F:72:PRO:HD2	2.02	0.60
1:C:89:GLU:HB3	3:C:2031:HOH:O	2.01	0.59
1:C:34:MET:CE	1:C:79:VAL:HG11	2.32	0.59
1:C:72:GLN:HG3	1:C:73:ASP:O	2.02	0.59
1:C:34:MET:HE2	1:C:79:VAL:HG11	1.85	0.59
2:G:29:GLY:HA2	3:G:2016:HOH:O	2.02	0.59
1:A:34:MET:HE1	1:A:79:VAL:HG11	1.85	0.58
2:E:32:PRO:HG3	2:E:64:LEU:HD21	1.85	0.58
1:B:76:LYS:O	1:B:77:ASN:HB2	2.03	0.57
2:E:42:ASN:ND2	2:E:77:GLU:H	2.01	0.57
1:C:18:LEU:HB2	1:C:83:MET:HE3	1.87	0.57
2:F:31:HIS:ND1	2:F:31:HIS:N	2.53	0.56
1:B:28:THR:HA	2:D:6:MET:HE2	1.86	0.56
1:C:76:LYS:O	1:C:77:ASN:HB2	2.06	0.56
1:C:116:ARG:HG3	1:C:116:ARG:O	2.06	0.56
2:G:97:ARG:CG	2:G:97:ARG:NH1	2.52	0.55
1:C:45:ARG:NH2	1:C:113:ASP:OD2	2.40	0.55
1:A:115:PHE:O	1:A:118:TRP:HZ3	1.90	0.54
2:G:42:ASN:HD21	2:G:77:GLU:H	1.55	0.54
1:C:19:ARG:HH21	1:C:82:GLN:HE21	1.56	0.54
1:C:6:GLU:OE1	1:C:122:THR:HG22	2.07	0.54
1:A:69:THR:HG23	3:A:2026:HOH:O	2.08	0.53
1:B:1:GLN:CD	1:B:1:GLN:H3	2.10	0.53
1:C:75:ALA:HB1	1:C:76:LYS:O	2.08	0.53
2:F:60:TRP:O	2:F:60:TRP:CD1	2.61	0.53
2:F:78:TYR:O	2:G:95:TRP:HB3	2.09	0.53
1:A:72:GLN:NE2	1:A:79:VAL:HG13	2.23	0.53
1:C:99:ASP:OD2	1:C:103:ARG:CZ	2.57	0.53
2:F:73:THR:CG2	2:F:76:ASP:OD2	2.56	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:28:THR:CB	2:G:6:MET:SD	2.97	0.52
1:B:103:ARG:HD3	2:D:10:TYR:CD1	2.44	0.52
2:G:24:ASN:HB3	2:G:65:LEU:HD22	1.92	0.52
1:C:75:ALA:HB3	1:C:76:LYS:HA	1.88	0.52
1:B:103:ARG:HD3	2:D:10:TYR:CG	2.44	0.52
1:B:76:LYS:HD2	1:B:78:THR:H	1.73	0.52
2:F:58:LYS:O	2:F:59:ASP:CB	2.58	0.51
2:F:59:ASP:O	2:F:61:SER:CA	2.51	0.51
2:G:97:ARG:HG2	2:G:97:ARG:NH1	2.22	0.51
1:A:72:GLN:HE22	1:A:79:VAL:CB	2.23	0.51
1:B:1:GLN:H3	1:B:1:GLN:NE2	2.09	0.50
1:C:72:GLN:O	1:C:73:ASP:HB2	2.11	0.50
1:B:44:GLU:HG3	1:B:45:ARG:N	2.25	0.50
1:C:72:GLN:HB3	3:C:2026:HOH:O	2.11	0.50
2:D:6:MET:N	3:D:2001:HOH:O	2.45	0.50
2:F:74:GLU:HG3	2:F:75:LYS:HG2	1.93	0.50
1:B:28:THR:OG1	2:D:6:MET:HG3	2.11	0.50
1:C:20:LEU:HG	1:C:83:MET:CE	2.42	0.50
1:C:76:LYS:HB3	1:C:78:THR:H	1.75	0.50
2:G:60:TRP:N	3:G:2033:HOH:O	2.45	0.49
1:C:34:MET:HE1	1:C:79:VAL:CG1	2.42	0.49
1:C:116:ARG:NH2	2:G:10:TYR:HB3	2.26	0.49
2:G:31:HIS:CE1	2:G:62:PHE:CE1	3.01	0.49
1:C:72:GLN:CG	1:C:73:ASP:N	2.75	0.49
1:A:28:THR:O	1:A:29:ASP:OD1	2.31	0.48
1:A:113:ASP:O	1:A:118:TRP:CH2	2.56	0.48
1:B:76:LYS:HD2	1:B:78:THR:OG1	2.13	0.48
2:G:6:MET:HB3	2:G:28:SER:O	2.13	0.48
1:A:34:MET:CE	1:A:79:VAL:HB	2.44	0.48
2:E:32:PRO:HD2	3:E:2015:HOH:O	2.13	0.48
2:G:54:LEU:HD12	2:G:64:LEU:HG	1.92	0.48
1:A:108:VAL:HG12	2:D:76:ASP:OD1	2.14	0.48
1:C:72:GLN:HG2	1:C:73:ASP:O	2.13	0.48
2:G:50:GLU:HB2	2:G:67:TYR:CZ	2.49	0.48
2:F:73:THR:HG23	3:F:2024:HOH:O	2.13	0.47
2:F:50:GLU:HB2	2:F:67:TYR:CZ	2.49	0.47
1:A:72:GLN:HE22	1:A:79:VAL:HA	1.79	0.47
1:A:72:GLN:HG3	3:A:2030:HOH:O	2.14	0.47
2:D:24:ASN:ND2	2:D:67:TYR:HB3	2.30	0.47
2:D:29:GLY:C	2:D:62:PHE:CE2	2.88	0.47
2:F:36:GLU:OE2	2:G:87:LEU:HD11	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:ALA:H	2:D:42:ASN:ND2	2.14	0.46
1:B:32:TYR:CE2	1:B:77:ASN:OD1	2.68	0.46
1:B:72:GLN:OE1	1:B:79:VAL:HG13	2.14	0.46
2:F:57:SER:HA	2:F:62:PHE:HE2	1.80	0.46
2:F:61:SER:HB2	2:F:62:PHE:CD2	2.51	0.46
1:B:72:GLN:HG3	1:B:73:ASP:N	2.30	0.46
1:B:56:ASP:OD2	2:F:81:ARG:HD3	2.15	0.46
1:A:68:PHE:HA	1:A:82:GLN:O	2.16	0.46
2:G:42:ASN:HD21	2:G:76:ASP:HA	1.80	0.46
1:B:44:GLU:CG	1:B:45:ARG:N	2.78	0.46
2:F:36:GLU:HA	3:F:2012:HOH:O	2.15	0.46
1:A:72:GLN:HE22	1:A:79:VAL:HG13	1.80	0.46
1:B:72:GLN:OE1	1:B:79:VAL:HG22	2.16	0.46
2:D:54:LEU:HD13	2:D:64:LEU:HD21	1.98	0.45
2:F:51:HIS:HA	2:F:65:LEU:O	2.17	0.45
2:D:42:ASN:ND2	2:D:77:GLU:H	2.13	0.45
1:A:34:MET:CE	1:A:79:VAL:CB	2.95	0.45
1:C:99:ASP:CG	1:C:103:ARG:NH1	2.67	0.45
1:C:103:ARG:CZ	1:C:116:ARG:HH21	2.30	0.45
2:F:59:ASP:CA	2:F:61:SER:CB	2.93	0.45
1:B:28:THR:HA	2:D:6:MET:CE	2.47	0.45
2:D:28:SER:HA	2:D:63:TYR:HA	1.98	0.45
1:B:91:THR:O	1:B:92:ALA:HB2	2.16	0.44
1:B:57:ILE:HD11	2:F:40:LEU:HD13	1.98	0.44
1:C:28:THR:HG23	2:G:28:SER:OG	2.18	0.44
2:F:95:TRP:HB3	3:G:2039:HOH:O	2.17	0.44
1:C:57:ILE:HD11	2:E:40:LEU:HD13	2.00	0.44
1:C:103:ARG:HH22	2:G:8:GLN:NE2	2.11	0.44
2:G:54:LEU:HD13	2:G:64:LEU:HG	1.97	0.44
1:C:116:ARG:HH22	2:G:10:TYR:HB3	1.83	0.44
2:F:41:LYS:O	2:F:42:ASN:C	2.53	0.43
1:B:57:ILE:HG12	2:G:92:ILE:HG21	1.99	0.43
2:F:50:GLU:HB2	2:F:67:TYR:CE2	2.54	0.43
1:C:34:MET:CE	1:C:79:VAL:CG1	2.96	0.43
3:C:2015:HOH:O	2:D:90:PRO:CG	2.67	0.43
2:F:12:ARG:HG2	2:F:13:HIS:CD2	2.54	0.43
1:B:5:GLN:O	1:B:22:CYS:HA	2.19	0.42
2:D:54:LEU:CD1	2:D:64:LEU:HD21	2.49	0.42
2:E:31:HIS:CD2	2:E:62:PHE:CZ	3.07	0.42
1:C:72:GLN:O	1:C:73:ASP:CB	2.68	0.42
1:C:76:LYS:O	1:C:77:ASN:CB	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:90:PRO:HB3	2:E:83:ASN:OD1	2.18	0.42
1:B:76:LYS:HB3	1:B:78:THR:N	2.35	0.42
1:B:31:ARG:HH11	2:D:58:LYS:NZ	2.17	0.42
1:C:68:PHE:HA	1:C:82:GLN:O	2.20	0.42
1:C:72:GLN:HG2	1:C:73:ASP:H	1.85	0.41
1:A:47:TRP:CB	1:A:111:GLY:HA2	2.50	0.41
1:B:107:ILE:O	1:B:108:VAL:C	2.57	0.41
1:C:100:ILE:HD11	1:C:103:ARG:HG3	2.03	0.41
2:E:66:TYR:N	2:E:66:TYR:CD2	2.88	0.41
1:B:57:ILE:CD1	2:F:40:LEU:HD13	2.51	0.41
1:A:76:LYS:O	1:A:77:ASN:CB	2.69	0.41
1:A:74:ASN:CG	1:A:74:ASN:O	2.59	0.41
1:B:20:LEU:O	1:B:80:TYR:HA	2.21	0.41
1:B:74:ASN:HA	1:B:75:ALA:HA	1.45	0.41
1:C:40:ALA:HB1	1:C:41:PRO:HD2	2.03	0.41
2:D:51:HIS:HA	2:D:65:LEU:O	2.21	0.41
2:G:39:LEU:HD23	2:G:68:THR:HG22	2.03	0.41
2:E:40:LEU:HD23	2:E:45:ARG:HA	2.02	0.41
1:B:54:GLY:HA2	2:D:58:LYS:HE2	2.03	0.40
2:G:23:LEU:O	2:G:67:TYR:HA	2.21	0.40
1:B:32:TYR:HE2	1:B:77:ASN:OD1	2.02	0.40
1:C:65:LYS:HZ2	1:C:65:LYS:HG2	1.20	0.40
1:A:89:GLU:H	1:A:89:GLU:CD	2.24	0.40
2:G:51:HIS:HB3	2:G:66:TYR:CD1	2.56	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	126/128 (98%)	119 (94%)	6 (5%)	1 (1%)	19 12

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	126/128 (98%)	118 (94%)	6 (5%)	2 (2%)	9	4
1	C	126/128 (98%)	119 (94%)	2 (2%)	5 (4%)	3	0
2	D	90/94 (96%)	85 (94%)	2 (2%)	3 (3%)	4	0
2	E	89/94 (95%)	85 (96%)	2 (2%)	2 (2%)	6	2
2	F	89/94 (95%)	85 (96%)	3 (3%)	1 (1%)	14	8
2	G	80/94 (85%)	77 (96%)	3 (4%)	0	100	100
All	All	726/760 (96%)	688 (95%)	24 (3%)	14 (2%)	8	2

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	72	GLN
1	C	73	ASP
2	D	20	SER
2	E	58	LYS
2	F	59	ASP
1	C	77	ASN
2	D	17	ASN
1	C	72	GLN
1	C	75	ALA
2	D	18	GLY
1	C	29	ASP
2	E	48	LYS
1	B	121	GLY
1	A	114	GLY

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	102/103 (99%)	95 (93%)	7 (7%)	15	10
1	B	103/103 (100%)	97 (94%)	6 (6%)	20	15
1	C	103/103 (100%)	91 (88%)	12 (12%)	5	2

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	D	84/89 (94%)	77 (92%)	7 (8%)	11 6
2	E	82/89 (92%)	79 (96%)	3 (4%)	34 32
2	F	86/89 (97%)	75 (87%)	11 (13%)	4 1
2	G	78/89 (88%)	69 (88%)	9 (12%)	5 2
All	All	638/665 (96%)	583 (91%)	55 (9%)	10 6

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	13	GLN
1	A	31	ARG
1	A	43	LYS
1	A	72	GLN
1	A	76	LYS
1	A	118	TRP
1	B	1	GLN
1	B	31	ARG
1	B	43	LYS
1	B	74	ASN
1	B	125	THR
1	B	128	SER
1	C	1	GLN
1	C	13	GLN
1	C	28	THR
1	C	31	ARG
1	C	44	GLU
1	C	46	GLU
1	C	65	LYS
1	C	71	SER
1	C	85	SER
1	C	116	ARG
1	C	122	THR
1	C	128	SER
2	D	6	MET
2	D	12	ARG
2	D	55	SER
2	D	58	LYS
2	D	60	TRP
2	D	70	PHE
2	D	74	GLU

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Mol	Chain	Res	Type
2	E	6	MET
2	E	26	TYR
2	E	70	PHE
2	F	26	TYR
2	F	31	HIS
2	F	47	GLU
2	F	48	LYS
2	F	50	GLU
2	F	52	SER
2	F	58	LYS
2	F	70	PHE
2	F	73	THR
2	F	74	GLU
2	F	91	LYS
2	G	8	GLN
2	G	11	SER
2	G	47	GLU
2	G	54	LEU
2	G	64	LEU
2	G	70	PHE
2	G	74	GLU
2	G	91	LYS
2	G	97	ARG

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

Mol	Chain	Res	Type
1	A	77	ASN
1	A	82	GLN
1	A	123	GLN
1	B	1	GLN
1	B	77	ASN
1	C	82	GLN
1	C	123	GLN
2	D	24	ASN
2	D	42	ASN
2	E	42	ASN
2	F	42	ASN
2	F	84	HIS
2	G	8	GLN
2	G	24	ASN
2	G	31	HIS

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Mol	Chain	Res	Type
2	G	42	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 monosaccharides 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)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	55:ARG	C	56:ASP	N	1.17

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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	128/128 (100%)	1.22	29 (22%)	0	0	53, 59, 73, 88	0
1	B	128/128 (100%)	1.26	31 (24%)	0	0	23, 59, 72, 94	0
1	C	128/128 (100%)	1.37	28 (21%)	0	0	23, 58, 75, 94	0
2	D	92/94 (97%)	1.41	24 (26%)	0	0	24, 61, 74, 77	0
2	E	91/94 (96%)	1.19	24 (26%)	0	0	48, 60, 73, 75	0
2	F	91/94 (96%)	1.45	23 (25%)	0	0	49, 59, 76, 92	0
2	G	86/94 (91%)	2.06	40 (46%)	0	0	47, 60, 71, 79	0
All	All	744/760 (97%)	1.40	199 (26%)	0	0	23, 59, 75, 94	0

All (199) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	60	TRP	11.9
2	G	32	PRO	8.9
1	C	74	ASN	8.8
1	C	73	ASP	7.8
1	B	75	ALA	7.7
1	B	73	ASP	6.6
2	G	30	PHE	6.2
1	C	35	ALA	6.0
2	G	29	GLY	5.9
2	G	33	SER	5.9
1	A	29	ASP	5.9
2	F	59	ASP	5.8
2	F	10	TYR	5.5
1	B	74	ASN	5.5
2	F	23	LEU	5.5
1	C	34	MET	5.4
1	C	36	TRP	5.4

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Mol	Chain	Res	Type	RSRZ
1	C	75	ALA	5.3
2	F	61	SER	5.3
1	A	34	MET	5.3
2	E	61	SER	5.3
1	C	117	TYR	5.2
2	E	23	LEU	5.1
1	A	81	LEU	5.1
1	B	77	ASN	4.9
2	D	27	VAL	4.8
2	E	60	TRP	4.7
1	C	51	ILE	4.6
1	A	35	ALA	4.6
1	C	97	ALA	4.6
1	A	51	ILE	4.5
2	F	58	LYS	4.5
1	B	72	GLN	4.4
1	C	33	CYS	4.4
1	B	36	TRP	4.4
1	A	74	ASN	4.3
2	G	21	ASN	4.3
2	F	9	VAL	4.2
1	B	35	ALA	4.2
2	D	39	LEU	4.2
1	B	1	GLN	4.1
2	G	89	GLN	4.1
1	A	28	THR	4.1
2	D	20	SER	4.0
2	F	62	PHE	4.0
2	D	78	TYR	4.0
1	B	34	MET	4.0
2	D	23	LEU	3.9
1	B	97	ALA	3.9
1	B	96	CYS	3.9
2	G	27	VAL	3.8
2	F	79	ALA	3.8
1	C	1	GLN	3.8
2	G	74	GLU	3.7
2	G	22	PHE	3.7
1	C	86	LEU	3.7
1	C	98	THR	3.7
2	E	48	LYS	3.7
2	F	27	VAL	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	20	LEU	3.6
2	D	40	LEU	3.6
1	C	127	SER	3.6
1	A	1	GLN	3.6
2	G	24	ASN	3.6
2	D	26	TYR	3.6
1	B	3	GLN	3.6
2	G	73	THR	3.6
1	B	33	CYS	3.5
2	G	95	TRP	3.5
2	D	48	LYS	3.5
2	D	15	ALA	3.5
1	C	3	GLN	3.4
2	G	14	PRO	3.4
1	A	120	GLN	3.4
2	G	25	CYS	3.4
2	G	75	LYS	3.4
2	G	6	MET	3.4
2	D	95	TRP	3.4
2	E	40	LEU	3.4
1	C	32	TYR	3.3
2	G	23	LEU	3.3
2	F	25	CYS	3.3
2	E	25	CYS	3.3
2	D	6	MET	3.3
2	F	78	TYR	3.3
1	B	18	LEU	3.3
1	A	50	ARG	3.2
1	C	44	GLU	3.2
2	G	48	LYS	3.1
1	B	51	ILE	3.1
2	D	60	TRP	3.1
2	D	59	ASP	3.1
1	B	81	LEU	3.1
1	C	37	PHE	3.1
1	A	97	ALA	3.1
1	A	33	CYS	3.0
1	A	75	ALA	3.0
2	F	48	LYS	3.0
2	D	74	GLU	2.9
2	F	95	TRP	2.9
2	E	19	LYS	2.9

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Mol	Chain	Res	Type	RSRZ
2	G	47	GLU	2.9
2	F	26	TYR	2.9
2	E	56	PHE	2.9
1	C	29	ASP	2.9
1	A	44	GLU	2.9
2	G	31	HIS	2.9
1	A	5	GLN	2.8
1	B	37	PHE	2.8
1	C	76	LYS	2.8
2	D	14	PRO	2.8
2	F	24	ASN	2.8
2	G	63	TYR	2.8
2	G	43	GLY	2.8
2	E	10	TYR	2.8
1	A	128	SER	2.8
2	D	64	LEU	2.7
2	G	39	LEU	2.7
1	A	77	ASN	2.7
2	F	6	MET	2.7
1	A	18	LEU	2.7
2	G	70	PHE	2.7
1	B	29	ASP	2.7
1	B	48	VAL	2.7
2	G	50	GLU	2.7
2	D	58	LYS	2.6
2	D	47	GLU	2.6
1	A	36	TRP	2.6
1	C	94	TYR	2.6
2	D	96	ASP	2.6
1	A	14	ALA	2.6
2	E	37	VAL	2.6
1	C	31	ARG	2.6
1	C	50	ARG	2.6
2	E	6	MET	2.6
1	A	86	LEU	2.6
2	F	69	GLU	2.5
2	E	70	PHE	2.5
2	D	16	GLU	2.5
1	C	19	ARG	2.5
2	G	71	THR	2.5
1	A	41	PRO	2.5
1	C	48	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
2	G	37	VAL	2.5
2	E	80	CYS	2.5
1	B	32	TYR	2.4
2	E	24	ASN	2.4
1	B	49	ALA	2.4
1	B	120	GLN	2.4
2	E	79	ALA	2.4
1	A	42	GLY	2.4
1	A	22	CYS	2.4
2	D	62	PHE	2.4
1	C	49	ALA	2.4
1	B	47	TRP	2.4
2	G	56	PHE	2.3
2	G	80	CYS	2.3
1	A	72	GLN	2.3
1	B	117	TYR	2.3
1	B	30	SER	2.3
1	B	98	THR	2.3
2	E	78	TYR	2.3
1	A	79	VAL	2.3
1	A	76	LYS	2.3
1	C	41	PRO	2.3
2	D	79	ALA	2.2
2	E	39	LEU	2.2
2	G	26	TYR	2.2
2	F	31	HIS	2.2
2	G	76	ASP	2.2
1	B	12	VAL	2.2
2	G	69	GLU	2.2
1	A	3	GLN	2.2
2	F	39	LEU	2.2
1	B	86	LEU	2.2
2	G	65	LEU	2.2
1	B	128	SER	2.2
2	G	55	SER	2.2
2	E	9	VAL	2.2
2	G	66	TYR	2.2
2	G	13	HIS	2.2
2	D	80	CYS	2.1
2	G	34	ASP	2.1
2	E	27	VAL	2.1
2	E	66	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
2	D	25	CYS	2.1
2	E	49	VAL	2.1
2	F	36	GLU	2.1
2	E	30	PHE	2.1
2	G	20	SER	2.1
2	E	59	ASP	2.1
1	A	83	MET	2.1
2	E	26	TYR	2.1
2	G	8	GLN	2.1
1	B	100	ILE	2.1
2	F	40	LEU	2.0
2	F	70	PHE	2.0
2	G	53	ASP	2.0
1	C	72	GLN	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 monosaccharides 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.