Full wwPDB X-ray Structure Validation Report (i)

Nov 28, 2023 – 07:37 pm GMT

PDB ID	:	1DXX
Title	:	N-terminal Actin-binding Domain of Human Dystrophin
Authors	:	Norwood, F.L.; Sutherland-Smith, A.J.; Keep, N.H.; Kendrick-Jones, J.
Deposited on	:	2000-01-20
Resolution	:	2.60 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:

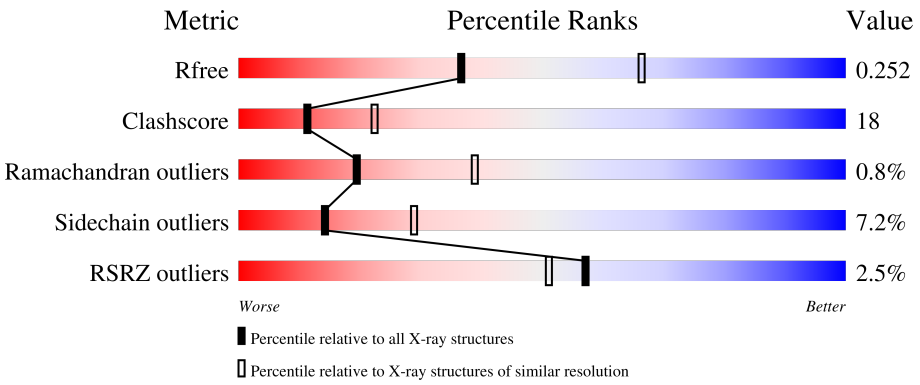
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	246	2% 60% 30% 6% • •
1	B	246	3% 61% 29% 5% • •
1	C	246	3% 63% 28% 5% • •
1	D	246	2% 59% 31% 5% • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 7622 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 DYSTROPHIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	238	1887	1197	332	355	3	0	0	0
1	B	238	1885	1194	333	355	3	0	0	0
1	C	238	1891	1199	333	356	3	0	0	0
1	D	238	1885	1193	333	356	3	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	10	SER	CYS	engineered mutation	UNP P11532
B	10	SER	CYS	engineered mutation	UNP P11532
C	10	SER	CYS	engineered mutation	UNP P11532
D	10	SER	CYS	engineered mutation	UNP P11532
A	188	SER	CYS	engineered mutation	UNP P11532
B	188	SER	CYS	engineered mutation	UNP P11532
C	188	SER	CYS	engineered mutation	UNP P11532
D	188	SER	CYS	engineered mutation	UNP P11532

- Molecule 2 is water.

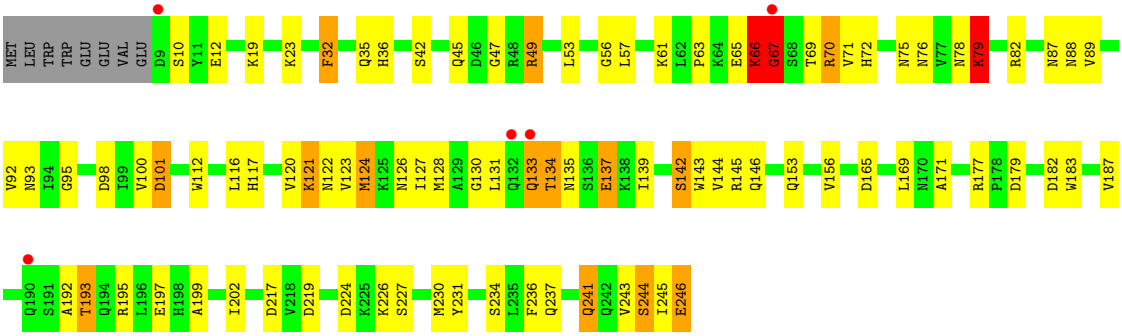
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	26	26	26	0	0
2	B	19	19	19	0	0
2	C	14	14	14	0	0
2	D	15	15	15	0	0

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.

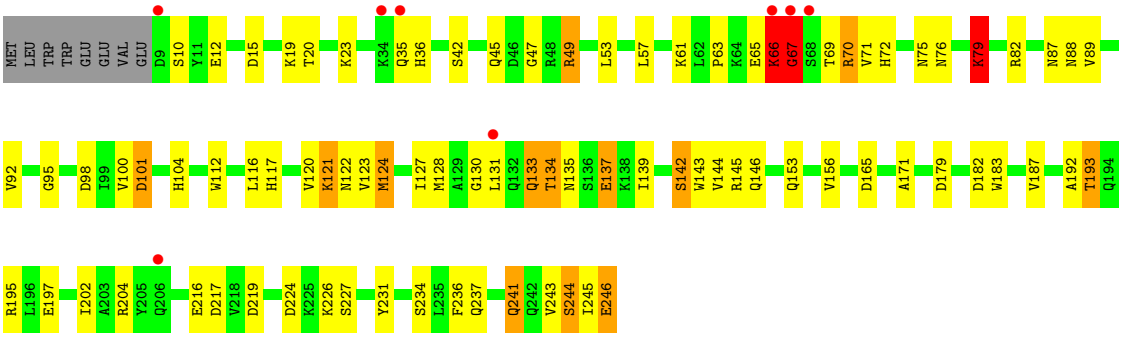
- Molecule 1: DYSTROPHIN

Chain A:



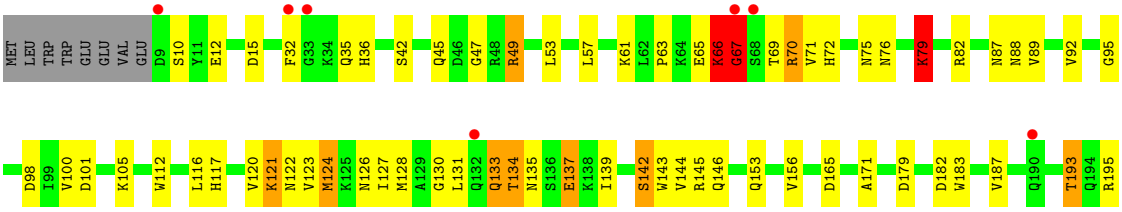
- Molecule 1: DYSTROPHIN

Chain B:



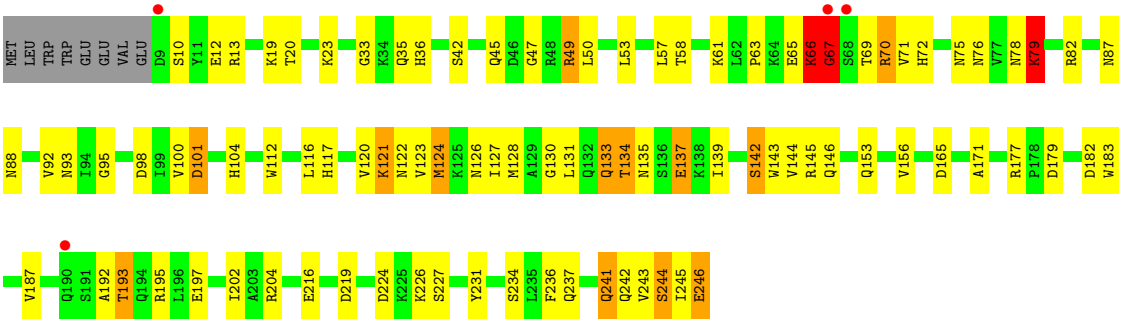
- Molecule 1: DYSTROPHIN

Chain C:



• Molecule 1: DYSTROPHIN

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	59.69Å 79.33Å 81.95Å 61.08° 78.22° 70.54°	Depositor
Resolution (Å)	20.00 – 2.60 39.93 – 2.60	Depositor EDS
% Data completeness (in resolution range)	95.0 (20.00-2.60) 95.2 (39.93-2.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.29 (at 2.61Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.232 , 0.268 0.223 , 0.252	Depositor DCC
R_{free} test set	1799 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å^2)	38.7	Xtriage
Anisotropy	0.348	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 53.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.130 for -h,-k,-k+l	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	7622	wwPDB-VP
Average B, all atoms (Å^2)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 26.35 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.6765e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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	0.64	0/1923	1.79	34/2614 (1.3%)
1	B	0.65	0/1920	1.77	34/2609 (1.3%)
1	C	0.64	0/1927	1.80	35/2619 (1.3%)
1	D	0.65	0/1920	1.77	32/2610 (1.2%)
All	All	0.65	0/7690	1.78	135/10452 (1.3%)

There are no bond length outliers.

All (135) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	145	ARG	NE-CZ-NH2	-20.56	110.02	120.30
1	A	145	ARG	NE-CZ-NH2	-19.73	110.44	120.30
1	B	145	ARG	NE-CZ-NH2	-19.64	110.48	120.30
1	D	145	ARG	NE-CZ-NH2	-18.56	111.02	120.30
1	C	145	ARG	NE-CZ-NH1	18.22	129.41	120.30
1	D	145	ARG	NE-CZ-NH1	17.74	129.17	120.30
1	D	219	ASP	CB-CG-OD2	-17.52	102.54	118.30
1	B	145	ARG	NE-CZ-NH1	17.51	129.05	120.30
1	B	219	ASP	CB-CG-OD2	-17.37	102.67	118.30
1	A	145	ARG	NE-CZ-NH1	17.34	128.97	120.30
1	C	219	ASP	CB-CG-OD2	-16.82	103.17	118.30
1	A	219	ASP	CB-CG-OD2	-16.55	103.41	118.30
1	A	32	PHE	CB-CG-CD2	-14.46	110.68	120.80
1	C	12	GLU	CA-CB-CG	10.74	137.04	113.40
1	D	12	GLU	CA-CB-CG	10.70	136.94	113.40
1	A	12	GLU	CA-CB-CG	10.67	136.87	113.40
1	C	49	ARG	NE-CZ-NH1	10.66	125.63	120.30
1	B	12	GLU	CA-CB-CG	10.55	136.60	113.40
1	D	49	ARG	NE-CZ-NH1	10.28	125.44	120.30
1	B	49	ARG	NE-CZ-NH1	10.16	125.38	120.30
1	C	32	PHE	CB-CG-CD1	-10.12	113.71	120.80
1	A	32	PHE	CB-CG-CD1	9.69	127.58	120.80
1	A	49	ARG	NE-CZ-NH1	9.67	125.13	120.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	82	ARG	CD-NE-CZ	9.62	137.07	123.60
1	A	145	ARG	CD-NE-CZ	9.49	136.88	123.60
1	D	145	ARG	CD-NE-CZ	9.45	136.82	123.60
1	C	145	ARG	CD-NE-CZ	9.39	136.75	123.60
1	B	145	ARG	CD-NE-CZ	9.37	136.72	123.60
1	C	32	PHE	CB-CG-CD2	9.34	127.34	120.80
1	A	82	ARG	CD-NE-CZ	9.15	136.41	123.60
1	C	82	ARG	CD-NE-CZ	9.07	136.30	123.60
1	B	82	ARG	CD-NE-CZ	8.98	136.17	123.60
1	C	195	ARG	CD-NE-CZ	8.93	136.09	123.60
1	A	195	ARG	CD-NE-CZ	8.67	135.74	123.60
1	B	195	ARG	CD-NE-CZ	8.67	135.74	123.60
1	D	195	ARG	CD-NE-CZ	8.45	135.44	123.60
1	D	67	GLY	CA-C-O	7.87	134.77	120.60
1	B	70	ARG	NE-CZ-NH1	-7.81	116.39	120.30
1	A	67	GLY	CA-C-O	7.80	134.63	120.60
1	C	67	GLY	CA-C-O	7.68	134.42	120.60
1	B	67	GLY	CA-C-O	7.65	134.37	120.60
1	C	101	ASP	CB-CG-OD1	7.62	125.16	118.30
1	B	195	ARG	NE-CZ-NH1	7.57	124.08	120.30
1	D	101	ASP	CB-CG-OD1	7.53	125.08	118.30
1	D	195	ARG	NE-CZ-NH1	7.45	124.03	120.30
1	A	195	ARG	NE-CZ-NH1	7.44	124.02	120.30
1	D	137	GLU	OE1-CD-OE2	7.42	132.20	123.30
1	C	195	ARG	NE-CZ-NH1	7.37	123.98	120.30
1	C	70	ARG	NE-CZ-NH1	-7.26	116.67	120.30
1	D	79	LYS	CA-CB-CG	7.22	129.28	113.40
1	A	224	ASP	CB-CG-OD1	-7.11	111.90	118.30
1	A	79	LYS	CA-CB-CG	7.11	129.04	113.40
1	B	12	GLU	N-CA-CB	7.05	123.28	110.60
1	B	79	LYS	CA-CB-CG	7.04	128.90	113.40
1	C	79	LYS	CA-CB-CG	6.97	128.75	113.40
1	D	12	GLU	N-CA-CB	6.96	123.14	110.60
1	A	137	GLU	OE1-CD-OE2	6.91	131.59	123.30
1	B	101	ASP	CB-CG-OD1	6.89	124.50	118.30
1	C	12	GLU	N-CA-CB	6.86	122.95	110.60
1	A	12	GLU	N-CA-CB	6.86	122.95	110.60
1	B	224	ASP	CB-CG-OD1	-6.80	112.18	118.30
1	C	137	GLU	OE1-CD-OE2	6.52	131.13	123.30
1	C	142	SER	N-CA-CB	6.49	120.23	110.50
1	A	70	ARG	NE-CZ-NH1	-6.45	117.08	120.30
1	B	182	ASP	CB-CG-OD2	-6.43	112.51	118.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	182	ASP	CB-CG-OD2	-6.39	112.55	118.30
1	B	165	ASP	CB-CG-OD1	6.33	124.00	118.30
1	B	137	GLU	OE1-CD-OE2	6.32	130.88	123.30
1	B	142	SER	N-CA-CB	6.31	119.96	110.50
1	D	182	ASP	N-CA-CB	-6.30	99.25	110.60
1	D	70	ARG	NE-CZ-NH1	-6.28	117.16	120.30
1	A	182	ASP	CB-CG-OD2	-6.24	112.69	118.30
1	D	49	ARG	NH1-CZ-NH2	-6.24	112.54	119.40
1	C	224	ASP	CB-CG-OD1	-6.18	112.74	118.30
1	C	165	ASP	CA-C-N	6.17	128.55	116.20
1	C	92	VAL	CA-CB-CG2	-6.16	101.66	110.90
1	A	89	VAL	O-C-N	-6.15	112.86	122.70
1	C	165	ASP	CB-CG-OD1	6.13	123.82	118.30
1	D	142	SER	N-CA-CB	6.10	119.65	110.50
1	D	182	ASP	CB-CG-OD2	-6.10	112.81	118.30
1	A	101	ASP	CB-CG-OD1	6.09	123.78	118.30
1	A	165	ASP	CA-C-N	6.09	128.38	116.20
1	A	142	SER	N-CA-CB	6.06	119.59	110.50
1	C	182	ASP	N-CA-CB	-6.05	99.70	110.60
1	D	219	ASP	CB-CG-OD1	6.05	123.75	118.30
1	B	92	VAL	CA-CB-CG2	-6.04	101.84	110.90
1	A	165	ASP	CB-CG-OD1	6.03	123.72	118.30
1	B	165	ASP	CA-C-N	5.99	128.19	116.20
1	B	182	ASP	N-CA-CB	-5.94	99.91	110.60
1	C	219	ASP	OD1-CG-OD2	5.90	134.51	123.30
1	A	182	ASP	N-CA-CB	-5.89	100.00	110.60
1	D	165	ASP	CA-C-N	5.84	127.88	116.20
1	A	49	ARG	NH1-CZ-NH2	-5.83	112.99	119.40
1	B	49	ARG	NH1-CZ-NH2	-5.83	112.99	119.40
1	C	49	ARG	NH1-CZ-NH2	-5.82	113.00	119.40
1	B	101	ASP	CB-CG-OD2	-5.70	113.17	118.30
1	C	105	LYS	CA-CB-CG	5.70	125.93	113.40
1	B	67	GLY	CA-C-N	-5.68	104.70	117.20
1	D	67	GLY	CA-C-N	-5.68	104.71	117.20
1	C	121	LYS	CD-CE-NZ	5.66	124.71	111.70
1	B	219	ASP	CB-CG-OD1	5.63	123.37	118.30
1	D	165	ASP	CB-CG-OD1	5.62	123.36	118.30
1	D	92	VAL	CA-CB-CG2	-5.62	102.47	110.90
1	B	219	ASP	OD1-CG-OD2	5.61	133.96	123.30
1	A	67	GLY	CA-C-N	-5.60	104.87	117.20
1	A	219	ASP	OD1-CG-OD2	5.60	133.95	123.30
1	D	121	LYS	CD-CE-NZ	5.55	124.47	111.70

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	204	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	C	67	GLY	CA-C-N	-5.54	105.01	117.20
1	D	224	ASP	CB-CG-OD1	-5.53	113.33	118.30
1	B	121	LYS	CD-CE-NZ	5.51	124.37	111.70
1	A	121	LYS	CD-CE-NZ	5.50	124.35	111.70
1	D	177	ARG	NE-CZ-NH2	-5.49	117.56	120.30
1	D	219	ASP	OD1-CG-OD2	5.48	133.71	123.30
1	A	177	ARG	NE-CZ-NH2	-5.38	117.61	120.30
1	B	61	LYS	N-CA-CB	5.37	120.27	110.60
1	D	13	ARG	NE-CZ-NH2	5.36	122.98	120.30
1	B	104	HIS	CA-CB-CG	-5.33	104.54	113.60
1	A	224	ASP	OD1-CG-OD2	5.33	133.42	123.30
1	D	204	ARG	NE-CZ-NH1	5.32	122.96	120.30
1	B	82	ARG	NE-CZ-NH2	-5.32	117.64	120.30
1	A	82	ARG	NE-CZ-NH2	-5.28	117.66	120.30
1	C	101	ASP	CB-CG-OD2	-5.28	113.55	118.30
1	A	61	LYS	N-CA-CB	5.27	120.09	110.60
1	C	15	ASP	CB-CG-OD1	5.26	123.04	118.30
1	D	61	LYS	N-CA-CB	5.26	120.07	110.60
1	B	15	ASP	CB-CG-OD1	5.18	122.96	118.30
1	C	70	ARG	NE-CZ-NH2	5.17	122.89	120.30
1	C	61	LYS	N-CA-CB	5.14	119.86	110.60
1	B	89	VAL	O-C-N	-5.14	114.47	122.70
1	A	70	ARG	CB-CA-C	5.13	120.65	110.40
1	A	92	VAL	CA-CB-CG2	-5.12	103.21	110.90
1	C	89	VAL	O-C-N	-5.11	114.53	122.70
1	D	104	HIS	CA-CB-CG	-5.08	104.97	113.60
1	C	70	ARG	CB-CA-C	5.04	120.48	110.40

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	1887	0	1861	84	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1885	0	1865	78	0
1	C	1891	0	1867	78	0
1	D	1885	0	1860	85	0
2	A	26	0	0	1	0
2	B	19	0	0	0	0
2	C	14	0	0	0	0
2	D	15	0	0	2	0
All	All	7622	0	7453	267	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 18.

All (267) 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:C:245:ILE:HD12	1:D:245:ILE:HD12	1.29	1.12
1:A:245:ILE:HD12	1:B:245:ILE:HD12	1.29	1.11
1:A:69:THR:HB	1:A:72:HIS:HD2	1.34	0.91
1:C:69:THR:HB	1:C:72:HIS:HD2	1.36	0.91
1:D:69:THR:HB	1:D:72:HIS:HD2	1.35	0.90
1:B:69:THR:HB	1:B:72:HIS:HD2	1.35	0.89
1:C:69:THR:HG22	1:C:71:VAL:H	1.44	0.81
1:C:237:GLN:NE2	1:D:122:ASN:HD22	1.79	0.81
1:B:69:THR:HG22	1:B:71:VAL:H	1.46	0.80
1:A:69:THR:HG22	1:A:71:VAL:H	1.45	0.80
1:D:69:THR:HG22	1:D:71:VAL:H	1.46	0.79
1:A:128:MET:HG2	1:B:131:LEU:CD1	2.13	0.79
1:C:122:ASN:HD22	1:D:237:GLN:NE2	1.82	0.78
1:A:146:GLN:HE22	1:A:244:SER:HB2	1.51	0.76
1:A:131:LEU:CD1	1:B:128:MET:HG2	2.16	0.75
1:D:146:GLN:HE22	1:D:244:SER:HB2	1.52	0.75
1:D:33:GLY:HA2	2:D:2002:HOH:O	1.87	0.75
1:A:122:ASN:HD22	1:B:237:GLN:NE2	1.85	0.75
1:B:146:GLN:HE22	1:B:244:SER:HB2	1.52	0.74
1:C:146:GLN:HE22	1:C:244:SER:HB2	1.52	0.74
1:C:237:GLN:HE22	1:D:122:ASN:HD22	1.35	0.73
1:C:131:LEU:CD1	1:D:128:MET:HG2	2.20	0.72
1:C:135:ASN:HD21	1:C:137:GLU:HB2	1.55	0.71
1:D:69:THR:HG22	1:D:71:VAL:N	2.07	0.70
1:A:135:ASN:HD21	1:A:137:GLU:HB2	1.55	0.70
1:C:146:GLN:NE2	1:C:244:SER:HB2	2.07	0.70

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:146:GLN:NE2	1:B:244:SER:HB2	2.06	0.70
1:C:69:THR:HG22	1:C:71:VAL:N	2.06	0.70
1:D:146:GLN:NE2	1:D:244:SER:HB2	2.07	0.70
1:B:135:ASN:HD21	1:B:137:GLU:HB2	1.56	0.69
1:D:135:ASN:HD21	1:D:137:GLU:HB2	1.55	0.69
1:C:122:ASN:HD22	1:D:237:GLN:HE22	1.37	0.69
1:A:69:THR:HG22	1:A:71:VAL:N	2.07	0.69
1:A:146:GLN:NE2	1:A:244:SER:HB2	2.06	0.69
1:A:237:GLN:NE2	1:B:122:ASN:HD22	1.89	0.69
1:B:69:THR:HG22	1:B:71:VAL:N	2.07	0.68
1:A:69:THR:HB	1:A:72:HIS:CD2	2.25	0.68
1:C:128:MET:HG2	1:D:131:LEU:CD1	2.23	0.68
1:B:130:GLY:O	1:B:133:GLN:NE2	2.27	0.67
1:A:130:GLY:O	1:A:133:GLN:NE2	2.28	0.67
1:D:130:GLY:O	1:D:133:GLN:NE2	2.28	0.66
1:D:63:PRO:HD2	1:D:79:LYS:NZ	2.11	0.66
1:D:69:THR:HB	1:D:72:HIS:CD2	2.26	0.66
1:A:124:MET:CE	1:B:131:LEU:HD22	2.25	0.66
1:A:124:MET:HE2	1:B:131:LEU:HD13	1.77	0.65
1:A:131:LEU:HD13	1:B:124:MET:HE2	1.77	0.65
1:B:69:THR:HB	1:B:72:HIS:CD2	2.26	0.65
1:C:130:GLY:O	1:C:133:GLN:NE2	2.29	0.65
1:A:63:PRO:HD2	1:A:79:LYS:NZ	2.12	0.64
1:B:63:PRO:HD2	1:B:79:LYS:NZ	2.10	0.64
1:C:63:PRO:HD2	1:C:79:LYS:NZ	2.12	0.64
1:A:122:ASN:HD22	1:B:237:GLN:HE22	1.45	0.64
1:A:131:LEU:HD22	1:B:124:MET:CE	2.28	0.63
1:B:87:ASN:O	1:B:88:ASN:HB2	1.97	0.63
1:C:144:VAL:HG11	1:C:156:VAL:HG11	1.82	0.62
1:A:128:MET:HG2	1:B:131:LEU:HD12	1.81	0.62
1:D:144:VAL:HG11	1:D:156:VAL:HG11	1.83	0.61
1:A:87:ASN:O	1:A:88:ASN:HB2	2.00	0.61
1:B:144:VAL:HG11	1:B:156:VAL:HG11	1.82	0.61
1:C:87:ASN:O	1:C:88:ASN:HB2	2.00	0.61
1:A:127:ILE:HG21	1:B:245:ILE:HD11	1.83	0.60
1:A:237:GLN:HE22	1:B:122:ASN:HD22	1.48	0.60
1:C:131:LEU:HD13	1:D:124:MET:HE2	1.82	0.60
1:D:63:PRO:HD2	1:D:79:LYS:HZ2	1.66	0.60
1:A:131:LEU:HD12	1:B:128:MET:HG2	1.82	0.59
1:B:63:PRO:HD2	1:B:79:LYS:HZ2	1.66	0.59
1:A:144:VAL:HG11	1:A:156:VAL:HG11	1.83	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:ILE:HD11	1:B:127:ILE:HG21	1.85	0.59
1:A:49:ARG:HG2	2:A:2006:HOH:O	2.01	0.59
1:D:87:ASN:O	1:D:88:ASN:HB2	2.02	0.59
1:C:124:MET:HE2	1:D:131:LEU:HD13	1.83	0.58
1:C:131:LEU:HD22	1:D:124:MET:CE	2.33	0.58
1:D:35:GLN:O	1:D:49:ARG:NH2	2.37	0.58
1:C:67:GLY:HA3	1:C:72:HIS:HB2	1.86	0.58
1:D:67:GLY:HA3	1:D:72:HIS:HB2	1.87	0.56
1:C:69:THR:HB	1:C:72:HIS:CD2	2.27	0.56
1:C:124:MET:CE	1:D:131:LEU:HD22	2.35	0.56
1:D:143:TRP:HB2	1:D:243:VAL:HG12	1.87	0.56
1:B:67:GLY:HA3	1:B:72:HIS:HB2	1.86	0.56
1:C:130:GLY:HA3	1:C:246:GLU:HG2	1.88	0.56
1:B:130:GLY:HA3	1:B:246:GLU:HG2	1.88	0.56
1:A:67:GLY:HA3	1:A:72:HIS:HB2	1.87	0.56
1:D:130:GLY:HA3	1:D:246:GLU:HG2	1.88	0.55
1:C:116:LEU:HD12	1:C:120:VAL:HB	1.88	0.55
1:C:143:TRP:HB2	1:C:243:VAL:HG12	1.87	0.55
1:D:231:TYR:O	1:D:234:SER:HB2	2.06	0.55
1:A:130:GLY:HA3	1:A:246:GLU:HG2	1.89	0.55
1:A:143:TRP:HB2	1:A:243:VAL:HG12	1.88	0.55
1:C:35:GLN:O	1:C:49:ARG:NH2	2.39	0.55
1:B:35:GLN:O	1:B:49:ARG:NH2	2.40	0.55
1:A:53:LEU:HG	1:A:57:LEU:HD12	1.89	0.54
1:B:143:TRP:HB2	1:B:243:VAL:HG12	1.88	0.54
1:A:116:LEU:HD12	1:A:120:VAL:HB	1.90	0.54
1:A:35:GLN:O	1:A:49:ARG:NH2	2.41	0.54
1:C:53:LEU:HG	1:C:57:LEU:HD12	1.90	0.54
1:D:53:LEU:HG	1:D:57:LEU:CD1	2.38	0.54
1:B:231:TYR:O	1:B:234:SER:HB2	2.08	0.54
1:C:53:LEU:HG	1:C:57:LEU:CD1	2.37	0.54
1:D:53:LEU:HG	1:D:57:LEU:HD12	1.89	0.54
1:B:53:LEU:HG	1:B:57:LEU:HD12	1.90	0.54
1:B:53:LEU:HG	1:B:57:LEU:CD1	2.38	0.53
1:C:112:TRP:CZ2	1:D:226:LYS:HG2	2.44	0.53
1:D:116:LEU:HD12	1:D:120:VAL:HB	1.90	0.53
1:A:53:LEU:HG	1:A:57:LEU:CD1	2.39	0.53
1:B:67:GLY:HA3	1:B:72:HIS:CB	2.38	0.53
1:B:116:LEU:HD12	1:B:120:VAL:HB	1.91	0.53
1:C:231:TYR:O	1:C:234:SER:HB2	2.09	0.53
1:D:67:GLY:HA3	1:D:72:HIS:CB	2.39	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231:TYR:O	1:A:234:SER:HB2	2.09	0.52
1:D:133:GLN:HG2	1:D:133:GLN:O	2.10	0.52
1:B:133:GLN:O	1:B:133:GLN:HG2	2.09	0.52
1:C:67:GLY:HA3	1:C:72:HIS:CB	2.39	0.52
1:A:67:GLY:HA3	1:A:72:HIS:CB	2.39	0.52
1:B:66:LYS:HE2	1:B:66:LYS:N	2.26	0.51
1:A:133:GLN:O	1:A:133:GLN:HG2	2.10	0.51
1:A:123:VAL:O	1:A:127:ILE:HG12	2.11	0.51
1:C:133:GLN:O	1:C:133:GLN:HG2	2.10	0.50
1:C:241:GLN:HE22	1:D:126:ASN:HD22	1.58	0.50
1:D:123:VAL:O	1:D:127:ILE:HG12	2.12	0.50
1:A:226:LYS:HG2	1:B:112:TRP:CZ2	2.46	0.50
1:C:66:LYS:N	1:C:66:LYS:HE2	2.27	0.50
1:A:63:PRO:HD2	1:A:79:LYS:HZ1	1.75	0.50
1:C:126:ASN:HD22	1:D:241:GLN:HE22	1.60	0.50
1:C:131:LEU:HD12	1:D:128:MET:HG2	1.93	0.50
1:C:226:LYS:HG2	1:D:112:TRP:CZ2	2.46	0.50
1:C:135:ASN:HD21	1:C:137:GLU:CB	2.24	0.50
1:B:135:ASN:HD21	1:B:137:GLU:CB	2.24	0.49
1:A:112:TRP:CZ2	1:B:226:LYS:HG2	2.46	0.49
1:D:66:LYS:HE2	1:D:66:LYS:N	2.27	0.49
1:C:65:GLU:CD	1:C:75:ASN:HD22	2.16	0.49
1:B:123:VAL:O	1:B:127:ILE:HG12	2.12	0.49
1:A:66:LYS:N	1:A:66:LYS:HE2	2.27	0.49
1:B:47:GLY:HA3	1:B:76:ASN:HB3	1.95	0.49
1:C:123:VAL:O	1:C:127:ILE:HG12	2.12	0.48
1:C:128:MET:HG2	1:D:131:LEU:HD12	1.95	0.48
1:B:65:GLU:CD	1:B:75:ASN:HD22	2.16	0.48
1:D:47:GLY:HA3	1:D:76:ASN:HB3	1.96	0.48
1:D:42:SER:O	1:D:45:GLN:HB2	2.14	0.48
1:A:47:GLY:HA3	1:A:76:ASN:HB3	1.95	0.48
1:C:245:ILE:HD11	1:D:127:ILE:HG21	1.94	0.48
1:D:135:ASN:HD21	1:D:137:GLU:CB	2.24	0.48
1:C:47:GLY:HA3	1:C:76:ASN:HB3	1.95	0.48
1:D:65:GLU:CD	1:D:75:ASN:HD22	2.17	0.48
1:D:124:MET:HG3	1:D:124:MET:O	2.13	0.47
1:B:144:VAL:CG1	1:B:156:VAL:HG11	2.45	0.47
1:A:66:LYS:HB2	1:A:67:GLY:H	1.39	0.47
1:C:144:VAL:CG1	1:C:156:VAL:HG11	2.44	0.47
1:A:65:GLU:CD	1:A:75:ASN:HD22	2.17	0.47
1:C:63:PRO:HD2	1:C:79:LYS:HZ2	1.76	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:124:MET:O	1:C:124:MET:HG3	2.15	0.47
1:C:245:ILE:CD1	1:D:245:ILE:HD12	2.21	0.47
1:D:193:THR:O	1:D:197:GLU:HG3	2.15	0.47
1:A:135:ASN:HD21	1:A:137:GLU:CB	2.25	0.47
1:B:45:GLN:HG2	1:B:100:VAL:O	2.14	0.47
1:B:66:LYS:HE2	1:B:66:LYS:H	1.80	0.47
1:B:42:SER:O	1:B:45:GLN:HB2	2.15	0.47
1:A:131:LEU:HD21	1:A:245:ILE:CD1	2.44	0.47
1:C:127:ILE:HG21	1:D:245:ILE:HD11	1.96	0.47
1:A:124:MET:O	1:A:124:MET:HG3	2.15	0.46
1:C:131:LEU:HD21	1:C:245:ILE:CD1	2.45	0.46
1:D:131:LEU:HD21	1:D:245:ILE:CD1	2.45	0.46
1:A:32:PHE:CE2	1:A:56:GLY:HA3	2.50	0.46
1:D:183:TRP:O	1:D:187:VAL:HG23	2.15	0.46
1:C:69:THR:O	1:C:70:ARG:C	2.54	0.46
1:C:242:GLN:HA	1:D:245:ILE:O	2.15	0.46
1:D:144:VAL:CG1	1:D:156:VAL:HG11	2.45	0.46
1:A:69:THR:O	1:A:70:ARG:C	2.53	0.46
1:B:124:MET:O	1:B:124:MET:HG3	2.16	0.46
1:A:144:VAL:CG1	1:A:156:VAL:HG11	2.46	0.46
1:B:69:THR:O	1:B:70:ARG:C	2.53	0.46
1:B:131:LEU:HD21	1:B:245:ILE:CD1	2.46	0.46
1:C:183:TRP:O	1:C:187:VAL:HG23	2.15	0.46
1:D:66:LYS:HE2	1:D:66:LYS:H	1.80	0.46
1:A:193:THR:O	1:A:197:GLU:HG3	2.16	0.46
1:B:66:LYS:HB2	1:B:67:GLY:H	1.39	0.46
1:C:66:LYS:HE2	1:C:66:LYS:H	1.81	0.46
1:C:131:LEU:CD2	1:C:245:ILE:HD13	2.46	0.46
1:A:66:LYS:HE2	1:A:66:LYS:H	1.80	0.45
1:C:245:ILE:HD12	1:D:245:ILE:CD1	2.21	0.45
1:A:124:MET:HE1	1:B:131:LEU:HD22	1.96	0.45
1:A:245:ILE:HG22	1:A:246:GLU:N	2.31	0.45
1:B:193:THR:O	1:B:197:GLU:HG3	2.16	0.45
1:C:245:ILE:HG22	1:C:246:GLU:N	2.32	0.45
1:A:131:LEU:CD2	1:A:245:ILE:HD13	2.46	0.45
1:B:95:GLY:O	1:B:98:ASP:HB2	2.17	0.45
1:B:245:ILE:HG22	1:B:246:GLU:N	2.32	0.45
1:A:45:GLN:HG2	1:A:100:VAL:O	2.16	0.45
1:A:183:TRP:O	1:A:187:VAL:HG23	2.16	0.45
1:A:70:ARG:NH2	1:A:101:ASP:OD2	2.49	0.44
1:B:183:TRP:O	1:B:187:VAL:HG23	2.16	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:245:ILE:CG2	1:C:246:GLU:N	2.80	0.44
1:D:45:GLN:HG2	1:D:100:VAL:O	2.17	0.44
1:C:171:ALA:HB2	1:C:183:TRP:CH2	2.53	0.44
1:A:245:ILE:CG2	1:A:246:GLU:N	2.80	0.44
1:C:42:SER:O	1:C:45:GLN:HB2	2.17	0.44
1:A:42:SER:O	1:A:45:GLN:HB2	2.17	0.44
1:B:131:LEU:CD2	1:B:245:ILE:HD13	2.47	0.44
1:B:245:ILE:CG2	1:B:246:GLU:N	2.81	0.44
1:C:193:THR:O	1:C:197:GLU:HG3	2.17	0.44
1:D:131:LEU:CD2	1:D:245:ILE:HD13	2.47	0.44
1:C:66:LYS:HB2	1:C:67:GLY:H	1.39	0.44
1:D:69:THR:O	1:D:70:ARG:C	2.55	0.44
1:D:245:ILE:HG22	1:D:246:GLU:N	2.33	0.44
1:D:245:ILE:CG2	1:D:246:GLU:N	2.81	0.44
1:A:93:ASN:ND2	1:D:78:ASN:OD1	2.42	0.43
1:A:131:LEU:HD13	1:B:128:MET:HG2	1.99	0.43
1:C:45:GLN:HG2	1:C:100:VAL:O	2.18	0.43
1:A:127:ILE:CG2	1:B:245:ILE:HD11	2.46	0.43
1:B:134:THR:CG2	1:B:139:ILE:HD11	2.48	0.43
1:B:216:GLU:H	1:B:216:GLU:CD	2.22	0.43
1:A:78:ASN:OD1	1:D:93:ASN:ND2	2.44	0.43
1:A:126:ASN:HD22	1:B:241:GLN:HE22	1.67	0.43
1:A:171:ALA:HB2	1:A:183:TRP:CH2	2.54	0.43
1:B:153:GLN:NE2	1:B:183:TRP:HE1	2.17	0.43
1:C:236:PHE:CE2	1:C:241:GLN:HB3	2.53	0.43
1:C:245:ILE:O	1:D:242:GLN:HA	2.19	0.43
1:A:32:PHE:CD2	1:A:56:GLY:HA3	2.54	0.43
1:D:67:GLY:HA3	1:D:72:HIS:CG	2.54	0.43
1:D:70:ARG:NH2	1:D:101:ASP:OD2	2.50	0.43
1:D:58:THR:HA	2:D:2008:HOH:O	2.19	0.42
1:D:66:LYS:HB2	1:D:67:GLY:H	1.38	0.42
1:D:216:GLU:H	1:D:216:GLU:CD	2.22	0.42
1:A:67:GLY:HA3	1:A:72:HIS:CG	2.54	0.42
1:C:134:THR:CG2	1:C:139:ILE:HD11	2.49	0.42
1:A:245:ILE:HD11	1:B:127:ILE:CG2	2.48	0.42
1:A:128:MET:HG2	1:B:131:LEU:HD13	1.96	0.42
1:C:63:PRO:HD2	1:C:79:LYS:HZ1	1.81	0.42
1:A:112:TRP:CE2	1:B:226:LYS:HG2	2.54	0.42
1:A:236:PHE:CE2	1:A:241:GLN:HB3	2.54	0.42
1:D:117:HIS:HA	1:D:121:LYS:HE3	2.02	0.42
1:B:67:GLY:HA3	1:B:72:HIS:CG	2.54	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:192:ALA:O	1:B:193:THR:C	2.57	0.42
1:B:236:PHE:CE2	1:B:241:GLN:HB3	2.55	0.42
1:C:128:MET:HG2	1:D:131:LEU:HD13	2.02	0.42
1:D:153:GLN:NE2	1:D:183:TRP:HE1	2.18	0.42
1:D:171:ALA:HB2	1:D:183:TRP:CH2	2.54	0.42
1:D:192:ALA:O	1:D:193:THR:C	2.56	0.42
1:C:117:HIS:HA	1:C:121:LYS:HE3	2.02	0.42
1:C:153:GLN:NE2	1:C:183:TRP:HE1	2.18	0.42
1:C:241:GLN:OE1	1:D:126:ASN:HB3	2.19	0.42
1:D:134:THR:CG2	1:D:139:ILE:HD11	2.49	0.42
1:C:95:GLY:O	1:C:98:ASP:HB2	2.20	0.42
1:A:95:GLY:O	1:A:98:ASP:HB2	2.20	0.42
1:C:67:GLY:HA3	1:C:72:HIS:CG	2.55	0.42
1:D:236:PHE:CE2	1:D:241:GLN:HB3	2.55	0.42
1:A:192:ALA:O	1:A:193:THR:C	2.58	0.41
1:C:131:LEU:HD22	1:D:124:MET:HE1	2.01	0.41
1:A:117:HIS:HA	1:A:121:LYS:HE3	2.02	0.41
1:A:153:GLN:NE2	1:A:183:TRP:HE1	2.18	0.41
1:B:70:ARG:NH2	1:B:101:ASP:OD2	2.52	0.41
1:A:63:PRO:HD2	1:A:79:LYS:HZ2	1.82	0.41
1:B:117:HIS:HA	1:B:121:LYS:HE3	2.01	0.41
1:B:171:ALA:HB2	1:B:183:TRP:CH2	2.55	0.41
1:D:19:LYS:O	1:D:23:LYS:HG3	2.20	0.41
1:C:212:LEU:O	1:D:20:THR:HG22	2.21	0.41
1:A:19:LYS:NZ	1:B:217:ASP:OD1	2.52	0.41
1:D:50:LEU:HD23	1:D:50:LEU:HA	1.91	0.41
1:A:19:LYS:O	1:A:23:LYS:HG3	2.21	0.41
1:C:112:TRP:CE2	1:D:226:LYS:HG2	2.55	0.41
1:A:134:THR:CG2	1:A:139:ILE:HD11	2.51	0.41
1:A:217:ASP:OD1	1:B:19:LYS:NZ	2.52	0.41
1:A:230:MET:HE3	1:B:20:THR:HB	2.02	0.41
1:B:19:LYS:O	1:B:23:LYS:HG3	2.20	0.41
1:D:95:GLY:O	1:D:98:ASP:HB2	2.20	0.41
1:C:57:LEU:HD23	1:C:57:LEU:HA	1.96	0.41
1:C:226:LYS:HG2	1:D:112:TRP:CE2	2.56	0.41
1:A:169:LEU:HG	1:A:199:ALA:HB1	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	236/246 (96%)	219 (93%)	15 (6%)	2 (1%)	19	39
1	B	236/246 (96%)	221 (94%)	13 (6%)	2 (1%)	19	39
1	C	236/246 (96%)	220 (93%)	14 (6%)	2 (1%)	19	39
1	D	236/246 (96%)	220 (93%)	14 (6%)	2 (1%)	19	39
All	All	944/984 (96%)	880 (93%)	56 (6%)	8 (1%)	19	39

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	66	LYS
1	A	67	GLY
1	B	66	LYS
1	B	67	GLY
1	C	66	LYS
1	C	67	GLY
1	D	66	LYS
1	D	67	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	208/226 (92%)	193 (93%)	15 (7%)	14	29
1	B	208/226 (92%)	193 (93%)	15 (7%)	14	29

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	209/226 (92%)	194 (93%)	15 (7%)	14	29
1	D	208/226 (92%)	193 (93%)	15 (7%)	14	29
All	All	833/904 (92%)	773 (93%)	60 (7%)	14	29

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	SER
1	A	36	HIS
1	A	66	LYS
1	A	79	LYS
1	A	124	MET
1	A	133	GLN
1	A	134	THR
1	A	142	SER
1	A	179	ASP
1	A	193	THR
1	A	202	ILE
1	A	227	SER
1	A	241	GLN
1	A	244	SER
1	A	246	GLU
1	B	10	SER
1	B	36	HIS
1	B	66	LYS
1	B	79	LYS
1	B	124	MET
1	B	133	GLN
1	B	134	THR
1	B	142	SER
1	B	179	ASP
1	B	193	THR
1	B	202	ILE
1	B	227	SER
1	B	241	GLN
1	B	244	SER
1	B	246	GLU
1	C	10	SER
1	C	36	HIS
1	C	66	LYS
1	C	79	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	124	MET
1	C	133	GLN
1	C	134	THR
1	C	142	SER
1	C	179	ASP
1	C	193	THR
1	C	202	ILE
1	C	227	SER
1	C	241	GLN
1	C	244	SER
1	C	246	GLU
1	D	10	SER
1	D	36	HIS
1	D	66	LYS
1	D	79	LYS
1	D	124	MET
1	D	133	GLN
1	D	134	THR
1	D	142	SER
1	D	179	ASP
1	D	193	THR
1	D	202	ILE
1	D	227	SER
1	D	241	GLN
1	D	244	SER
1	D	246	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	HIS
1	A	72	HIS
1	A	117	HIS
1	A	126	ASN
1	A	133	GLN
1	A	146	GLN
1	A	237	GLN
1	B	36	HIS
1	B	72	HIS
1	B	117	HIS
1	B	126	ASN
1	B	133	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	146	GLN
1	B	237	GLN
1	C	36	HIS
1	C	72	HIS
1	C	117	HIS
1	C	126	ASN
1	C	133	GLN
1	C	146	GLN
1	C	237	GLN
1	D	36	HIS
1	D	72	HIS
1	D	117	HIS
1	D	126	ASN
1	D	133	GLN
1	D	146	GLN
1	D	237	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no 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)

There are no chain breaks in this entry.

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	238/246 (96%)	-0.18	5 (2%)	63	58	18, 35, 60, 74	0
1	B	238/246 (96%)	-0.05	8 (3%)	45	38	19, 35, 60, 75	0
1	C	238/246 (96%)	-0.13	7 (2%)	51	45	19, 35, 60, 75	0
1	D	238/246 (96%)	-0.15	4 (1%)	70	66	19, 35, 60, 75	0
All	All	952/984 (96%)	-0.13	24 (2%)	57	51	18, 35, 60, 75	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	9	ASP	8.2
1	B	9	ASP	5.5
1	C	9	ASP	4.3
1	A	9	ASP	4.1
1	B	67	GLY	3.6
1	D	190	GLN	3.6
1	B	68	SER	3.5
1	C	190	GLN	3.5
1	B	131	LEU	3.5
1	C	67	GLY	3.4
1	D	67	GLY	3.2
1	C	68	SER	3.0
1	A	67	GLY	2.7
1	A	190	GLN	2.7
1	C	132	GLN	2.6
1	D	68	SER	2.5
1	A	133	GLN	2.4
1	B	66	LYS	2.3
1	A	132	GLN	2.2
1	B	35	GLN	2.2
1	B	206	GLN	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	34	LYS	2.2
1	C	33	GLY	2.0
1	C	32	PHE	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.