Full wwPDB X-ray Structure Validation Report (i)

Nov 1, 2023 – 12:26 PM JST

PDB ID	:	5HAC
Title	:	Crystal structure of Proliferating Cell Nuclear Antigen from Leishmania donovani at 2.95 A resolution
Authors	:	Singh, P.K.; Yadav, S.P.; Sharma, P.; Kaur, P.; Sharma, S.; Singh, T.P.
Deposited on	:	2015-12-30
Resolution	:	2.95 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:

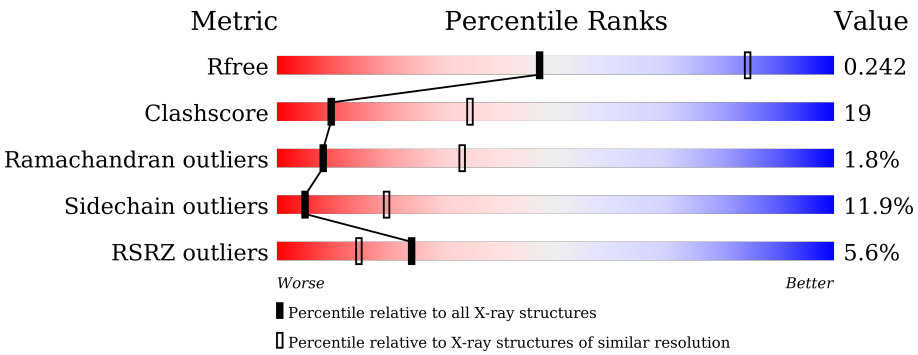
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 2.95 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	3104 (3.00-2.92)
Clashscore	141614	3462 (3.00-2.92)
Ramachandran outliers	138981	3340 (3.00-2.92)
Sidechain outliers	138945	3343 (3.00-2.92)
RSRZ outliers	127900	2986 (3.00-2.92)

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	300	8% 57% 21% 8% • 13%
1	B	300	4% 63% 19% • 13%
1	C	300	5% 62% 18% 5% • 13%
1	D	300	4% 62% 19% 6% 13%
1	E	300	5% 58% 24% • • 14%
1	F	300	4% 62% 18% 5% • 13%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 12076 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 Proliferating cell nuclear antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	260	2008	1263	337	393	15	0	0	0
1	B	261	2012	1265	338	394	15	0	0	0
1	C	260	2008	1263	337	393	15	0	0	0
1	D	261	2012	1265	338	394	15	0	0	0
1	E	259	2000	1257	336	392	15	0	0	0
1	F	261	2012	1265	338	394	15	0	0	0

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLY	-	expression tag	UNP B5TV91
A	-6	LEU	-	expression tag	UNP B5TV91
A	-5	VAL	-	expression tag	UNP B5TV91
A	-4	PRO	-	expression tag	UNP B5TV91
A	-3	ARG	-	expression tag	UNP B5TV91
A	-2	GLY	-	expression tag	UNP B5TV91
A	-1	SER	-	expression tag	UNP B5TV91
A	0	HIS	-	expression tag	UNP B5TV91
B	-7	GLY	-	expression tag	UNP B5TV91
B	-6	LEU	-	expression tag	UNP B5TV91
B	-5	VAL	-	expression tag	UNP B5TV91
B	-4	PRO	-	expression tag	UNP B5TV91
B	-3	ARG	-	expression tag	UNP B5TV91
B	-2	GLY	-	expression tag	UNP B5TV91
B	-1	SER	-	expression tag	UNP B5TV91
B	0	HIS	-	expression tag	UNP B5TV91
C	-7	GLY	-	expression tag	UNP B5TV91

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	-6	LEU	-	expression tag	UNP B5TV91
C	-5	VAL	-	expression tag	UNP B5TV91
C	-4	PRO	-	expression tag	UNP B5TV91
C	-3	ARG	-	expression tag	UNP B5TV91
C	-2	GLY	-	expression tag	UNP B5TV91
C	-1	SER	-	expression tag	UNP B5TV91
C	0	HIS	-	expression tag	UNP B5TV91
D	-7	GLY	-	expression tag	UNP B5TV91
D	-6	LEU	-	expression tag	UNP B5TV91
D	-5	VAL	-	expression tag	UNP B5TV91
D	-4	PRO	-	expression tag	UNP B5TV91
D	-3	ARG	-	expression tag	UNP B5TV91
D	-2	GLY	-	expression tag	UNP B5TV91
D	-1	SER	-	expression tag	UNP B5TV91
D	0	HIS	-	expression tag	UNP B5TV91
E	-7	GLY	-	expression tag	UNP B5TV91
E	-6	LEU	-	expression tag	UNP B5TV91
E	-5	VAL	-	expression tag	UNP B5TV91
E	-4	PRO	-	expression tag	UNP B5TV91
E	-3	ARG	-	expression tag	UNP B5TV91
E	-2	GLY	-	expression tag	UNP B5TV91
E	-1	SER	-	expression tag	UNP B5TV91
E	0	HIS	-	expression tag	UNP B5TV91
F	-7	GLY	-	expression tag	UNP B5TV91
F	-6	LEU	-	expression tag	UNP B5TV91
F	-5	VAL	-	expression tag	UNP B5TV91
F	-4	PRO	-	expression tag	UNP B5TV91
F	-3	ARG	-	expression tag	UNP B5TV91
F	-2	GLY	-	expression tag	UNP B5TV91
F	-1	SER	-	expression tag	UNP B5TV91
F	0	HIS	-	expression tag	UNP B5TV91

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	B	4	4	4	0	0
			Total	O		
2	C	7	7	7	0	0
			Total	O		
2	D	1	1	1	0	0
			Total	O		
2	E	3	3	3	0	0

Continued on next page...

Continued from previous page...

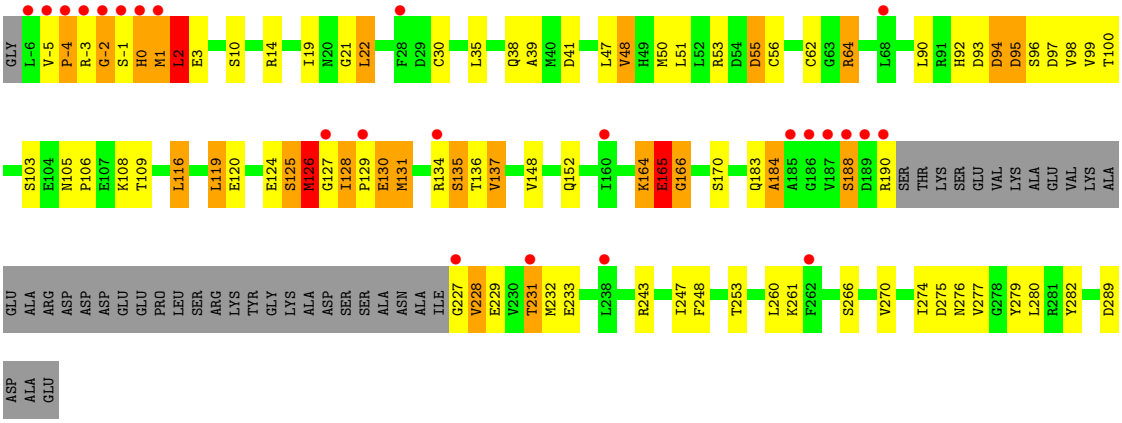
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	F	9	9	9	0	0

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.

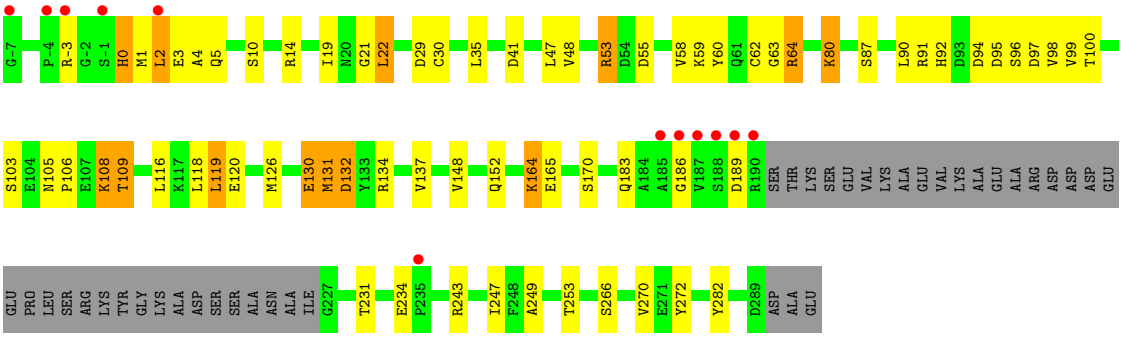
• Molecule 1: Proliferating cell nuclear antigen

Chain A: 8% 57% 21% 8% • 13%



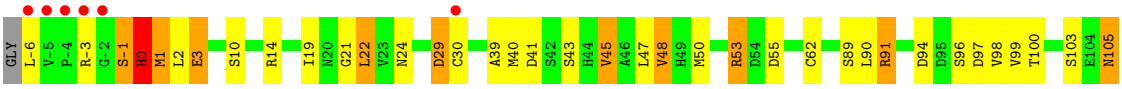
• Molecule 1: Proliferating cell nuclear antigen

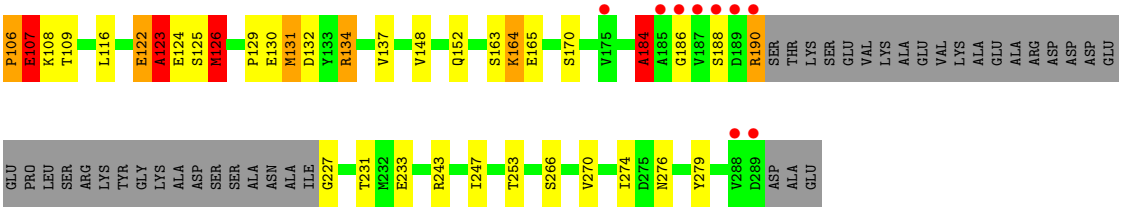
Chain B: 4% 63% 19% • 13%



• Molecule 1: Proliferating cell nuclear antigen

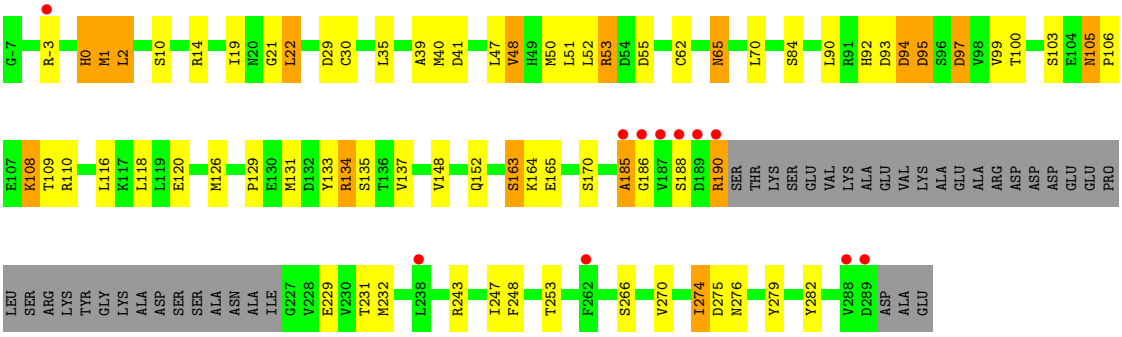
Chain C: 5% 62% 18% 5% • 13%





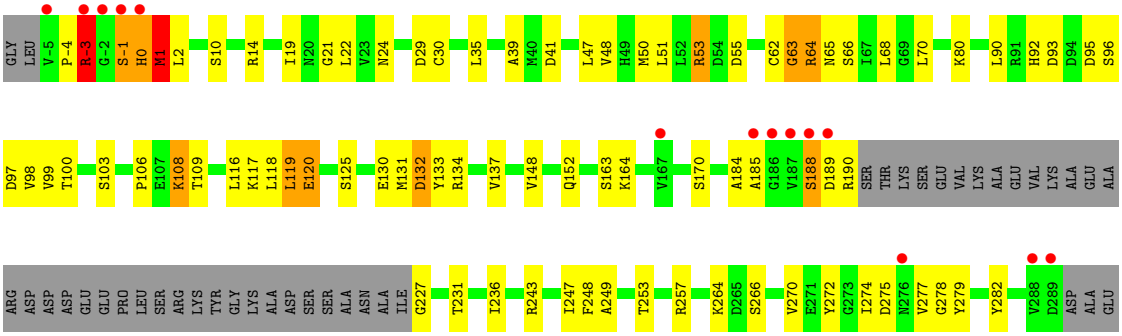
• Molecule 1: Proliferating cell nuclear antigen

Chain D:



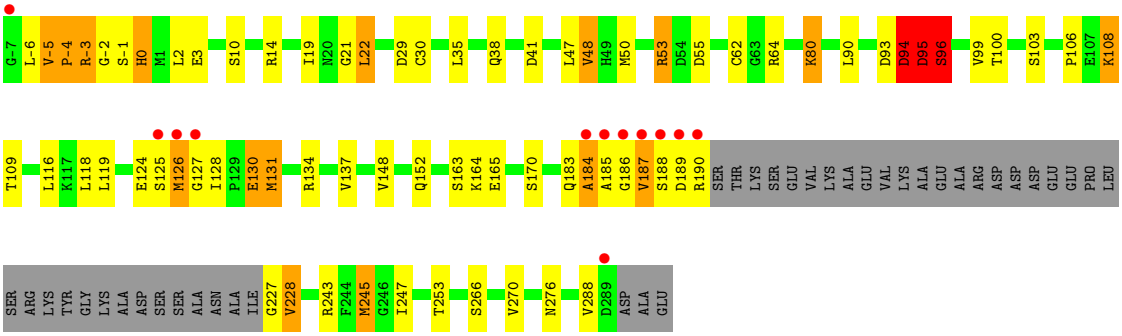
• Molecule 1: Proliferating cell nuclear antigen

Chain E:



• Molecule 1: Proliferating cell nuclear antigen

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	133.99Å 150.19Å 170.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.95 42.87 – 2.95	Depositor EDS
% Data completeness (in resolution range)	88.4 (50.00-2.95) 88.4 (42.87-2.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.19 (at 2.95Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.205 , 0.232 0.217 , 0.242	Depositor DCC
R_{free} test set	966 reflections (1.50%)	wwPDB-VP
Wilson B-factor (Å^2)	91.4	Xtriage
Anisotropy	0.267	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 64.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	12076	wwPDB-VP
Average B, all atoms (Å^2)	107.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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	0.73	0/2038	1.06	16/2750 (0.6%)
1	B	0.78	1/2042 (0.0%)	1.01	8/2755 (0.3%)
1	C	0.77	1/2038 (0.0%)	1.08	10/2750 (0.4%)
1	D	0.76	1/2042 (0.0%)	1.05	12/2755 (0.4%)
1	E	0.75	0/2030	1.02	9/2739 (0.3%)
1	F	0.78	1/2042 (0.0%)	1.05	10/2755 (0.4%)
All	All	0.76	4/12232 (0.0%)	1.05	65/16504 (0.4%)

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	3
1	B	0	1
1	C	0	5
1	D	0	3
1	E	0	1
1	F	0	4
All	All	0	17

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	119	LEU	C-O	5.72	1.34	1.23
1	D	275	ASP	CB-CG	5.63	1.63	1.51
1	B	186	GLY	N-CA	5.52	1.54	1.46
1	C	0	HIS	CA-C	5.23	1.66	1.52

All (65) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	64	ARG	NE-CZ-NH1	9.48	125.04	120.30
1	E	0	HIS	CB-CA-C	9.41	129.22	110.40
1	F	96	SER	CB-CA-C	9.00	127.21	110.10
1	F	95	ASP	CB-CA-C	-7.75	94.90	110.40
1	A	2	LEU	N-CA-C	7.60	131.53	111.00
1	F	-6	LEU	CB-CA-C	7.59	124.62	110.20
1	D	1	MET	N-CA-CB	-7.47	97.15	110.60
1	A	55	ASP	CB-CG-OD2	7.21	124.78	118.30
1	B	105	ASN	CB-CA-C	-7.12	96.17	110.40
1	E	63	GLY	N-CA-C	-6.93	95.78	113.10
1	A	48	VAL	CB-CA-C	-6.91	98.27	111.40
1	C	107	GLU	N-CA-CB	6.88	122.99	110.60
1	A	1	MET	N-CA-C	6.83	129.46	111.00
1	C	48	VAL	CB-CA-C	-6.81	98.47	111.40
1	F	48	VAL	CB-CA-C	-6.74	98.59	111.40
1	C	91	ARG	NE-CZ-NH1	6.73	123.67	120.30
1	E	164	LYS	N-CA-CB	6.67	122.61	110.60
1	D	48	VAL	CB-CA-C	-6.55	98.96	111.40
1	A	2	LEU	CB-CG-CD2	6.36	121.81	111.00
1	C	126	MET	CB-CG-SD	6.32	131.35	112.40
1	C	106	PRO	N-CA-C	-6.26	95.82	112.10
1	C	186	GLY	N-CA-C	-6.26	97.46	113.10
1	F	134	ARG	NE-CZ-NH1	6.24	123.42	120.30
1	D	135	SER	N-CA-CB	-6.21	101.18	110.50
1	D	275	ASP	N-CA-C	-6.21	94.24	111.00
1	C	105	ASN	CB-CA-C	-6.16	98.08	110.40
1	A	119	LEU	O-C-N	-6.10	112.94	122.70
1	A	131	MET	CG-SD-CE	6.06	109.90	100.20
1	B	22	LEU	CB-CG-CD2	5.94	121.09	111.00
1	E	119	LEU	O-C-N	-5.93	113.22	122.70
1	A	95	ASP	N-CA-C	-5.86	95.19	111.00
1	C	1	MET	CG-SD-CE	5.82	109.50	100.20
1	E	66	SER	N-CA-CB	5.75	119.13	110.50
1	E	-3	ARG	NE-CZ-NH2	5.70	123.15	120.30
1	D	0	HIS	CB-CA-C	-5.66	99.07	110.40
1	D	-3	ARG	NE-CZ-NH1	5.66	123.13	120.30
1	F	-5	VAL	N-CA-C	-5.65	95.75	111.00
1	F	288	VAL	CB-CA-C	5.58	122.01	111.40
1	F	94	ASP	CB-CG-OD1	5.58	123.32	118.30
1	A	119	LEU	CA-C-N	5.55	129.42	117.20
1	A	116	LEU	CA-CB-CG	5.54	128.04	115.30
1	A	166	GLY	N-CA-C	5.47	126.78	113.10
1	C	1	MET	N-CA-C	5.46	125.73	111.00

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	95	ASP	CB-CA-C	5.44	121.28	110.40
1	F	134	ARG	NE-CZ-NH2	-5.39	117.61	120.30
1	D	95	ASP	N-CA-C	5.32	125.36	111.00
1	B	119	LEU	O-C-N	-5.30	114.21	122.70
1	B	118	LEU	CA-CB-CG	5.29	127.45	115.30
1	D	0	HIS	C-N-CA	5.27	134.88	121.70
1	D	0	HIS	N-CA-C	-5.26	96.81	111.00
1	D	-3	ARG	NE-CZ-NH2	-5.25	117.67	120.30
1	E	119	LEU	CA-C-N	5.22	128.69	117.20
1	A	128	ILE	CB-CA-C	5.22	122.04	111.60
1	D	19	ILE	CB-CA-C	-5.21	101.17	111.60
1	A	3	GLU	N-CA-CB	-5.21	101.22	110.60
1	A	165	GLU	N-CA-C	-5.21	96.94	111.00
1	A	-2	GLY	N-CA-C	5.20	126.10	113.10
1	D	2	LEU	CA-CB-CG	5.18	127.23	115.30
1	F	118	LEU	CA-CB-CG	5.18	127.21	115.30
1	C	184	ALA	CB-CA-C	-5.13	102.41	110.10
1	E	1	MET	N-CA-C	5.12	124.84	111.00
1	A	128	ILE	N-CA-C	-5.12	97.18	111.00
1	B	119	LEU	CA-C-N	5.09	128.40	117.20
1	E	117	LYS	CA-CB-CG	5.08	124.58	113.40
1	B	131	MET	CA-CB-CG	5.04	121.87	113.30

There are no chirality outliers.

All (17) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	130	GLU	Peptide
1	A	94	ASP	Peptide
1	A	95	ASP	Peptide
1	B	130	GLU	Peptide
1	C	0	HIS	Peptide
1	C	122	GLU	Peptide
1	C	123	ALA	Peptide
1	C	184	ALA	Peptide
1	C	29	ASP	Mainchain
1	D	129	PRO	Peptide
1	D	134	ARG	Peptide
1	D	274	ILE	Peptide
1	E	130	GLU	Peptide
1	F	130	GLU	Peptide
1	F	227	GLY	Peptide

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Group
1	F	94	ASP	Peptide
1	F	95	ASP	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	2008	0	2003	125	1
1	B	2012	0	2005	70	1
1	C	2008	0	2002	42	0
1	D	2012	0	2006	46	0
1	E	2000	0	1992	79	0
1	F	2012	0	2006	103	0
2	B	4	0	0	0	0
2	C	7	0	0	0	0
2	D	1	0	0	0	0
2	E	3	0	0	1	0
2	F	9	0	0	0	0
All	All	12076	0	12014	464	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 19.

All (464) 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:F:47:LEU:HD11	1:F:126:MET:CE	1.08	1.54
1:B:30:CYS:SG	1:B:35:LEU:CD2	2.04	1.46
1:D:30:CYS:SG	1:D:35:LEU:CD2	2.06	1.43
1:F:47:LEU:CD1	1:F:126:MET:HE2	1.49	1.43
1:E:30:CYS:SG	1:E:35:LEU:CD2	2.11	1.37
1:D:30:CYS:SG	1:D:35:LEU:HD21	1.64	1.37
1:A:30:CYS:SG	1:A:35:LEU:CD2	2.12	1.36
1:E:30:CYS:SG	1:E:35:LEU:HD21	1.67	1.33
1:B:30:CYS:SG	1:B:35:LEU:HD21	1.62	1.33
1:F:30:CYS:SG	1:F:35:LEU:CD2	2.16	1.33

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:SER:HB3	1:A:232:MET:CA	1.60	1.30
1:F:47:LEU:CD1	1:F:126:MET:CE	2.01	1.30
1:F:30:CYS:SG	1:F:35:LEU:HD21	1.73	1.29
1:A:2:LEU:HD22	1:A:2:LEU:O	1.12	1.27
1:A:1:MET:O	1:A:92:HIS:O	1.52	1.25
1:A:135:SER:CB	1:A:232:MET:HA	1.65	1.24
1:A:30:CYS:SG	1:A:35:LEU:HD21	1.73	1.24
1:A:-4:PRO:CD	1:A:-3:ARG:HA	1.71	1.20
1:A:136:THR:HG23	1:A:260:LEU:O	1.43	1.19
1:A:2:LEU:O	1:A:2:LEU:CD2	1.89	1.19
1:B:48:VAL:HG22	1:B:282:TYR:CD1	1.78	1.18
1:F:-1:SER:HB3	1:F:0:HIS:CB	1.75	1.16
1:F:47:LEU:HD11	1:F:126:MET:HE3	1.24	1.12
1:F:-1:SER:CB	1:F:0:HIS:HB2	1.80	1.11
1:A:0:HIS:HB2	1:A:94:ASP:HB2	1.16	1.11
1:A:274:ILE:HD13	1:A:279:TYR:HA	1.33	1.11
1:F:185:ALA:HA	1:F:187:VAL:HG12	1.24	1.10
1:E:134:ARG:NH2	1:E:264:LYS:H	1.47	1.10
1:A:-1:SER:HB2	1:A:0:HIS:HB3	1.11	1.10
1:C:131:MET:HB2	1:C:132:ASP:HA	1.27	1.09
1:A:136:THR:CG2	1:A:260:LEU:O	1.99	1.09
1:B:30:CYS:O	1:B:62:CYS:SG	2.11	1.09
1:F:186:GLY:HA2	1:F:187:VAL:HG12	1.14	1.08
1:B:48:VAL:HG22	1:B:282:TYR:HD1	1.11	1.08
1:A:-4:PRO:HD2	1:A:-3:ARG:HA	1.22	1.08
1:A:274:ILE:HD13	1:A:279:TYR:CA	1.83	1.07
1:F:-1:SER:H	1:F:0:HIS:HB3	1.18	1.05
1:F:127:GLY:O	1:F:128:ILE:HG13	1.58	1.02
1:F:186:GLY:HA2	1:F:187:VAL:CG1	1.90	1.02
1:A:30:CYS:O	1:A:62:CYS:SG	2.17	1.02
1:A:0:HIS:CB	1:A:94:ASP:HB2	1.90	1.01
1:E:274:ILE:HD12	1:E:274:ILE:H	1.25	1.00
1:A:1:MET:HG2	1:A:93:ASP:OD1	1.62	1.00
1:E:134:ARG:HH21	1:E:264:LYS:N	1.58	0.99
1:C:30:CYS:O	1:C:62:CYS:SG	2.20	0.99
1:D:30:CYS:SG	1:D:35:LEU:HD23	2.02	0.98
1:B:30:CYS:SG	1:B:35:LEU:HD23	2.01	0.97
1:F:185:ALA:HA	1:F:187:VAL:CG1	1.96	0.95
1:F:184:ALA:HA	1:F:187:VAL:HG11	1.48	0.95
1:B:164:LYS:HD2	1:B:165:GLU:H	1.31	0.95
1:A:-1:SER:CB	1:A:0:HIS:HB3	1.97	0.94

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:30:CYS:O	1:D:62:CYS:SG	2.26	0.94
1:A:136:THR:HG22	1:A:137:VAL:H	1.32	0.94
1:F:126:MET:SD	1:F:127:GLY:HA2	2.08	0.94
1:E:30:CYS:O	1:E:62:CYS:SG	2.25	0.94
1:E:134:ARG:HH22	1:E:236:ILE:HD12	1.31	0.93
1:A:274:ILE:CD1	1:A:279:TYR:HA	1.97	0.93
1:F:30:CYS:O	1:F:62:CYS:SG	2.26	0.93
1:A:275:ASP:OD2	1:A:276:ASN:N	2.02	0.93
1:C:130:GLU:N	1:C:130:GLU:OE1	2.01	0.93
1:E:30:CYS:SG	1:E:35:LEU:HD23	2.07	0.93
1:A:64:ARG:HH11	1:A:64:ARG:HG2	1.33	0.91
1:A:30:CYS:SG	1:A:35:LEU:HD23	2.06	0.91
1:A:135:SER:HB3	1:A:232:MET:HA	0.90	0.90
1:A:135:SER:HB3	1:A:232:MET:CB	2.01	0.90
1:F:-1:SER:N	1:F:0:HIS:HB3	1.87	0.89
1:A:135:SER:OG	1:A:233:GLU:N	2.05	0.89
1:F:30:CYS:SG	1:F:35:LEU:HD23	2.12	0.88
1:A:-1:SER:HB2	1:A:0:HIS:CB	2.01	0.88
1:F:184:ALA:CB	1:F:187:VAL:HG11	2.05	0.87
1:D:0:HIS:CB	1:D:1:MET:HB2	2.04	0.87
1:F:184:ALA:CA	1:F:187:VAL:HG11	2.06	0.86
1:B:22:LEU:CD2	1:B:48:VAL:HG23	2.05	0.86
1:E:190:ARG:HB2	1:E:227:GLY:O	1.75	0.86
1:F:47:LEU:CD1	1:F:126:MET:HE3	1.87	0.86
1:F:47:LEU:CG	1:F:126:MET:HE3	2.05	0.86
1:F:184:ALA:O	1:F:228:VAL:HG22	1.75	0.86
1:C:131:MET:CB	1:C:132:ASP:HA	2.04	0.86
1:F:124:GLU:OE2	1:F:124:GLU:N	2.09	0.86
1:F:124:GLU:HA	1:F:125:SER:HB2	1.58	0.85
1:A:0:HIS:HB2	1:A:94:ASP:CB	2.04	0.85
1:E:53:ARG:HB2	1:E:277:VAL:O	1.76	0.84
1:F:184:ALA:HA	1:F:187:VAL:CG1	2.07	0.84
1:D:65:ASN:HD22	1:D:65:ASN:C	1.80	0.83
1:E:70:LEU:HD21	1:E:118:LEU:CD2	2.09	0.83
1:D:0:HIS:HB2	1:D:1:MET:HB2	1.57	0.83
1:F:-1:SER:HB3	1:F:0:HIS:HB2	0.89	0.83
1:E:274:ILE:HD13	1:E:278:GLY:C	1.98	0.82
1:B:3:GLU:HB2	1:B:91:ARG:HG3	1.59	0.82
1:A:-4:PRO:HD2	1:A:-3:ARG:CA	2.10	0.81
1:A:2:LEU:CD2	1:A:2:LEU:C	2.49	0.81
1:E:70:LEU:CD2	1:E:118:LEU:CD2	2.58	0.81

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:22:LEU:HD23	1:B:48:VAL:HG23	1.63	0.80
1:E:189:ASP:O	1:E:190:ARG:O	1.98	0.80
1:A:136:THR:HG22	1:A:137:VAL:N	1.97	0.80
1:A:1:MET:C	1:A:92:HIS:O	2.19	0.80
1:F:185:ALA:CA	1:F:187:VAL:HG12	2.11	0.80
1:C:105:ASN:HB3	1:C:106:PRO:O	1.81	0.79
1:F:126:MET:CG	1:F:127:GLY:HA2	2.13	0.79
1:E:134:ARG:NH2	1:E:264:LYS:N	2.21	0.79
1:B:5:GLN:HE21	1:B:87:SER:HB2	1.46	0.79
1:A:1:MET:HA	1:A:92:HIS:O	1.83	0.78
1:A:2:LEU:HD22	1:A:2:LEU:C	2.03	0.78
1:E:134:ARG:HH22	1:E:236:ILE:CD1	1.96	0.77
1:B:164:LYS:HD2	1:B:165:GLU:N	1.99	0.76
1:F:186:GLY:CA	1:F:187:VAL:HG12	2.08	0.75
1:A:-4:PRO:CG	1:A:-3:ARG:HA	2.16	0.74
1:F:47:LEU:HD21	1:F:126:MET:CE	2.17	0.74
1:B:48:VAL:CG2	1:B:282:TYR:CD1	2.67	0.74
1:A:2:LEU:O	1:A:2:LEU:CG	2.35	0.74
1:F:124:GLU:CA	1:F:125:SER:HB2	2.17	0.74
1:B:164:LYS:NZ	1:B:165:GLU:HG3	2.02	0.74
1:E:274:ILE:HD12	1:E:274:ILE:N	2.02	0.73
1:F:-1:SER:CB	1:F:0:HIS:CB	2.54	0.73
1:F:95:ASP:O	1:F:96:SER:HB2	1.87	0.73
1:E:133:TYR:O	1:E:133:TYR:HD1	1.72	0.73
1:F:124:GLU:HB2	1:F:125:SER:C	2.10	0.73
1:A:136:THR:HG22	1:A:260:LEU:O	1.89	0.73
1:F:188:SER:HB3	1:F:189:ASP:OD2	1.88	0.73
1:D:30:CYS:SG	1:D:35:LEU:HD22	2.24	0.72
1:E:24:ASN:HB3	2:E:303:HOH:O	1.89	0.72
1:F:47:LEU:HD11	1:F:126:MET:HE2	0.73	0.72
1:A:30:CYS:SG	1:A:35:LEU:HD22	2.29	0.72
1:A:135:SER:CA	1:A:232:MET:HA	2.20	0.72
1:F:47:LEU:HD11	1:F:126:MET:SD	2.29	0.72
1:E:133:TYR:O	1:E:133:TYR:CD1	2.43	0.71
1:A:136:THR:HG21	1:A:261:LYS:HG2	1.71	0.71
1:F:183:GLN:O	1:F:184:ALA:HB3	1.88	0.71
1:A:64:ARG:N	1:A:64:ARG:HD2	2.05	0.71
1:F:126:MET:HG2	1:F:127:GLY:CA	2.20	0.71
1:D:65:ASN:O	1:D:65:ASN:ND2	2.23	0.71
1:E:274:ILE:H	1:E:274:ILE:CD1	2.01	0.70
1:A:136:THR:HG23	1:A:261:LYS:HA	1.72	0.70

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:30:CYS:SG	1:E:35:LEU:HD22	2.29	0.70
1:B:30:CYS:SG	1:B:35:LEU:HD22	2.23	0.70
1:F:126:MET:CE	1:F:127:GLY:O	2.38	0.70
1:F:127:GLY:O	1:F:128:ILE:CG1	2.38	0.70
1:B:0:HIS:CE1	1:B:2:LEU:HD23	2.27	0.70
1:E:99:VAL:CG1	1:E:118:LEU:HD21	2.22	0.70
1:E:134:ARG:HH21	1:E:264:LYS:H	0.75	0.70
1:A:274:ILE:HD12	1:A:274:ILE:N	2.06	0.69
1:B:48:VAL:HG22	1:B:282:TYR:CE1	2.28	0.69
1:A:0:HIS:CE1	1:A:1:MET:HG3	2.27	0.69
1:E:64:ARG:HG3	1:E:65:ASN:N	2.06	0.69
1:A:-3:ARG:HD2	1:A:-3:ARG:C	2.15	0.68
1:D:0:HIS:HB3	1:D:1:MET:HB2	1.75	0.67
1:F:0:HIS:H	1:F:94:ASP:HB2	1.58	0.67
1:B:22:LEU:HD22	1:B:48:VAL:CG2	2.25	0.67
1:E:99:VAL:HG12	1:E:118:LEU:HD21	1.76	0.67
1:E:274:ILE:CD1	1:E:279:TYR:HA	2.24	0.67
1:E:70:LEU:HD23	1:E:118:LEU:CD2	2.25	0.67
1:A:275:ASP:OD2	1:A:277:VAL:N	2.27	0.66
1:C:43:SER:OG	1:C:45:VAL:HG13	1.95	0.66
1:B:22:LEU:HD22	1:B:48:VAL:HG23	1.76	0.66
1:D:190:ARG:O	1:D:229:GLU:HG3	1.96	0.65
1:F:47:LEU:HD13	1:F:126:MET:HE2	1.70	0.65
1:F:-3:ARG:NH1	1:F:-2:GLY:HA3	2.12	0.65
1:F:-1:SER:CA	1:F:0:HIS:CB	2.75	0.65
1:D:134:ARG:NH2	1:D:232:MET:O	2.30	0.64
1:F:126:MET:HE1	1:F:127:GLY:O	1.96	0.64
1:A:64:ARG:HG2	1:A:64:ARG:NH1	2.09	0.64
1:A:135:SER:CB	1:A:232:MET:CA	2.45	0.64
1:E:70:LEU:HD21	1:E:118:LEU:HD21	1.78	0.63
1:F:47:LEU:HD21	1:F:126:MET:HE3	1.80	0.63
1:A:274:ILE:CD1	1:A:279:TYR:CA	2.63	0.63
1:B:48:VAL:CG2	1:B:282:TYR:CE1	2.81	0.63
1:E:63:GLY:O	1:E:64:ARG:HG2	1.98	0.63
1:A:136:THR:CG2	1:A:137:VAL:H	2.08	0.63
1:B:5:GLN:HE21	1:B:87:SER:CB	2.10	0.63
1:F:-3:ARG:HH12	1:F:-1:SER:C	2.02	0.63
1:A:-3:ARG:HB2	1:A:-2:GLY:O	1.98	0.63
1:A:-4:PRO:CD	1:A:-3:ARG:CA	2.65	0.63
1:C:0:HIS:HB3	1:C:94:ASP:HB2	1.80	0.63
1:C:131:MET:HB2	1:C:132:ASP:CA	2.18	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:70:LEU:HD21	1:D:118:LEU:CD2	2.29	0.62
1:F:47:LEU:CD2	1:F:126:MET:HE3	2.30	0.62
1:B:22:LEU:CD2	1:B:48:VAL:CG2	2.77	0.62
1:B:3:GLU:CB	1:B:91:ARG:HG3	2.30	0.62
1:B:80:LYS:HE2	1:B:80:LYS:N	2.15	0.62
1:B:119:LEU:HB3	1:B:120:GLU:HB2	1.82	0.62
1:A:64:ARG:HH11	1:A:64:ARG:CG	2.11	0.62
1:A:135:SER:HA	1:A:231:THR:O	2.00	0.62
1:C:-1:SER:HB3	1:C:0:HIS:CG	2.35	0.62
1:F:187:VAL:O	1:F:187:VAL:HG13	1.99	0.61
1:F:124:GLU:HA	1:F:125:SER:CB	2.30	0.61
1:E:99:VAL:HB	1:E:118:LEU:HD11	1.82	0.61
1:F:124:GLU:CB	1:F:125:SER:HB2	2.31	0.61
1:A:190:ARG:HB3	1:A:227:GLY:HA2	1.82	0.60
1:D:99:VAL:CG1	1:D:118:LEU:HD21	2.31	0.60
1:B:3:GLU:CD	1:B:91:ARG:NE	2.55	0.60
1:B:130:GLU:N	1:B:131:MET:HB3	2.17	0.60
1:A:275:ASP:OD2	1:A:277:VAL:HG22	2.01	0.60
1:F:47:LEU:CD2	1:F:126:MET:CE	2.80	0.60
1:F:30:CYS:SG	1:F:35:LEU:HD22	2.33	0.60
1:A:1:MET:CA	1:A:92:HIS:O	2.50	0.60
1:E:119:LEU:HA	1:E:120:GLU:HG2	1.84	0.60
1:A:274:ILE:HD11	1:A:280:LEU:N	2.16	0.59
1:A:184:ALA:HB3	1:A:228:VAL:HG23	1.85	0.59
1:F:-3:ARG:NH1	1:F:-2:GLY:CA	2.66	0.59
1:A:274:ILE:HD13	1:A:279:TYR:C	2.22	0.59
1:B:30:CYS:C	1:B:62:CYS:SG	2.81	0.59
1:A:30:CYS:C	1:A:62:CYS:SG	2.82	0.59
1:F:245:MET:HE3	1:F:245:MET:HA	1.84	0.58
1:F:80:LYS:N	1:F:80:LYS:HE2	2.18	0.58
1:F:126:MET:SD	1:F:126:MET:N	2.75	0.58
1:A:-3:ARG:HD2	1:A:-3:ARG:O	2.03	0.58
1:F:183:GLN:O	1:F:184:ALA:CB	2.51	0.58
1:E:119:LEU:HB3	1:E:120:GLU:HB2	1.86	0.58
1:E:274:ILE:HD13	1:E:279:TYR:N	2.18	0.58
1:E:274:ILE:HG22	1:E:277:VAL:CG2	2.34	0.57
1:A:130:GLU:HB3	1:A:131:MET:HB2	1.87	0.57
1:A:136:THR:HG23	1:A:260:LEU:C	2.22	0.57
1:A:274:ILE:CD1	1:A:279:TYR:C	2.72	0.57
1:F:-1:SER:CA	1:F:0:HIS:HB3	2.33	0.57
1:E:70:LEU:CD2	1:E:118:LEU:HD22	2.35	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:51:LEU:HD23	1:E:279:TYR:CE1	2.40	0.57
1:A:0:HIS:CE1	1:A:1:MET:CG	2.87	0.56
1:C:43:SER:OG	1:C:45:VAL:CG1	2.53	0.56
1:A:188:SER:H	1:A:229:GLU:HG2	1.69	0.56
1:B:4:ALA:HB2	1:B:60:TYR:HD1	1.71	0.56
1:A:274:ILE:HD13	1:A:279:TYR:N	2.20	0.56
1:E:70:LEU:HD23	1:E:118:LEU:HD23	1.86	0.56
1:A:136:THR:CG2	1:A:261:LYS:HG2	2.36	0.56
1:F:124:GLU:CA	1:F:125:SER:CB	2.85	0.55
1:B:4:ALA:CB	1:B:60:TYR:HD1	2.20	0.55
1:B:5:GLN:NE2	1:B:87:SER:HB2	2.20	0.55
1:D:65:ASN:C	1:D:65:ASN:ND2	2.55	0.55
1:F:126:MET:HG2	1:F:127:GLY:C	2.27	0.55
1:B:164:LYS:CD	1:B:165:GLU:HG3	2.37	0.55
1:E:48:VAL:O	1:E:48:VAL:HG13	2.06	0.54
1:A:90:LEU:HD11	1:A:99:VAL:HG21	1.89	0.54
1:B:4:ALA:HB2	1:B:60:TYR:CD1	2.42	0.54
1:D:92:HIS:CE1	1:D:94:ASP:O	2.60	0.54
1:F:184:ALA:HB2	1:F:187:VAL:HG11	1.85	0.54
1:D:70:LEU:CD2	1:D:118:LEU:CD2	2.85	0.54
1:F:47:LEU:CG	1:F:126:MET:CE	2.71	0.54
1:F:245:MET:HA	1:F:245:MET:CE	2.38	0.54
1:E:90:LEU:HD11	1:E:99:VAL:HG21	1.89	0.54
1:F:124:GLU:HB2	1:F:125:SER:HB2	1.90	0.54
1:B:90:LEU:HD11	1:B:99:VAL:HG21	1.91	0.53
1:F:126:MET:CG	1:F:127:GLY:CA	2.81	0.53
1:C:134:ARG:CZ	1:C:233:GLU:OE1	2.56	0.53
1:D:51:LEU:HD23	1:D:279:TYR:CE1	2.43	0.53
1:E:274:ILE:CG2	1:E:277:VAL:CG2	2.85	0.53
1:D:40:MET:SD	1:D:126:MET:HG3	2.48	0.53
1:D:30:CYS:C	1:D:62:CYS:SG	2.87	0.53
1:B:3:GLU:OE2	1:B:91:ARG:CZ	2.57	0.53
1:E:53:ARG:HB3	1:E:55:ASP:OD1	2.09	0.53
1:F:-3:ARG:NH1	1:F:-1:SER:O	2.42	0.53
1:C:90:LEU:HD11	1:C:99:VAL:HG21	1.91	0.52
1:A:51:LEU:HD23	1:A:279:TYR:CE1	2.43	0.52
1:A:-3:ARG:HH21	1:A:64:ARG:HH12	1.57	0.52
1:B:3:GLU:OE2	1:B:91:ARG:NH2	2.43	0.52
1:A:119:LEU:HB3	1:A:120:GLU:HB2	1.91	0.52
1:D:90:LEU:HD11	1:D:99:VAL:HG21	1.91	0.52
1:D:190:ARG:C	1:D:229:GLU:HG3	2.29	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:64:ARG:HG3	1:E:65:ASN:H	1.72	0.52
1:F:30:CYS:C	1:F:62:CYS:SG	2.87	0.52
1:C:131:MET:CB	1:C:132:ASP:CA	2.85	0.52
1:A:-4:PRO:N	1:A:-3:ARG:HA	2.20	0.52
1:D:53:ARG:HB3	1:D:55:ASP:OD1	2.10	0.52
1:A:1:MET:HA	1:A:93:ASP:HA	1.92	0.51
1:A:274:ILE:CD1	1:A:274:ILE:N	2.73	0.51
1:B:164:LYS:HZ2	1:B:165:GLU:HG3	1.73	0.51
1:A:-3:ARG:NH2	1:A:64:ARG:HH22	2.09	0.51
1:B:53:ARG:HB3	1:B:55:ASP:OD1	2.11	0.51
1:F:124:GLU:HB2	1:F:125:SER:CA	2.40	0.51
1:F:90:LEU:HD11	1:F:99:VAL:HG21	1.93	0.51
1:A:130:GLU:HB3	1:A:131:MET:CB	2.41	0.51
1:C:122:GLU:HA	1:C:123:ALA:HB3	1.93	0.50
1:D:39:ALA:O	1:D:47:LEU:HD22	2.12	0.50
1:E:70:LEU:CD2	1:E:118:LEU:HD23	2.40	0.50
1:A:125:SER:O	1:A:126:MET:HB2	2.11	0.50
1:B:164:LYS:NZ	1:B:165:GLU:CG	2.73	0.50
1:C:164:LYS:O	1:C:165:GLU:HG2	2.11	0.50
1:A:275:ASP:CG	1:A:277:VAL:HG22	2.32	0.50
1:D:93:ASP:O	1:D:94:ASP:HB3	2.10	0.50
1:D:92:HIS:NE2	1:D:94:ASP:O	2.44	0.49
1:C:165:GLU:HA	1:C:184:ALA:HB3	1.93	0.49
1:A:-5:VAL:HA	1:A:-3:ARG:HB3	1.94	0.49
1:A:125:SER:O	1:A:126:MET:HE3	2.12	0.49
1:C:129:PRO:C	1:C:130:GLU:OE1	2.49	0.49
1:D:70:LEU:CD2	1:D:118:LEU:HD23	2.42	0.49
1:A:64:ARG:NH1	1:A:64:ARG:CG	2.72	0.49
1:A:164:LYS:O	1:A:165:GLU:CG	2.60	0.49
1:B:130:GLU:CD	1:B:130:GLU:H	2.16	0.49
1:C:30:CYS:C	1:C:62:CYS:SG	2.89	0.49
1:F:53:ARG:HB3	1:F:55:ASP:OD1	2.13	0.49
1:A:125:SER:O	1:A:126:MET:CE	2.60	0.49
1:E:96:SER:O	1:E:98:VAL:N	2.46	0.49
1:D:163:SER:OG	1:D:165:GLU:OE2	2.30	0.49
1:F:21:GLY:O	1:F:247:ILE:HD13	2.13	0.49
1:A:39:ALA:O	1:A:47:LEU:HD22	2.13	0.48
1:A:135:SER:HB3	1:A:232:MET:HB2	1.89	0.48
1:A:-3:ARG:NH2	1:A:64:ARG:NH2	2.61	0.48
1:F:185:ALA:N	1:F:228:VAL:O	2.46	0.48
1:A:64:ARG:N	1:A:64:ARG:CD	2.73	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:95:ASP:O	1:F:95:ASP:OD1	2.31	0.48
1:A:-3:ARG:HH21	1:A:64:ARG:NH1	2.12	0.48
1:B:164:LYS:HZ3	1:B:165:GLU:HG3	1.78	0.48
1:C:130:GLU:HB2	1:C:131:MET:SD	2.53	0.48
1:F:-3:ARG:HH11	1:F:-2:GLY:HA3	1.79	0.48
1:B:96:SER:O	1:B:98:VAL:N	2.47	0.48
1:C:53:ARG:HB3	1:C:55:ASP:OD1	2.13	0.48
1:C:188:SER:O	1:C:227:GLY:C	2.52	0.48
1:D:94:ASP:C	1:D:94:ASP:OD1	2.50	0.48
1:E:39:ALA:O	1:E:47:LEU:HD22	2.14	0.48
1:A:56:CYS:HB2	1:A:277:VAL:HB	1.96	0.48
1:D:99:VAL:HG12	1:D:118:LEU:HD21	1.95	0.48
1:D:105:ASN:N	1:D:105:ASN:ND2	2.61	0.48
1:A:-4:PRO:HG2	1:A:-3:ARG:HA	1.93	0.47
1:A:165:GLU:HG3	1:A:166:GLY:H	1.80	0.47
1:C:14:ARG:HD2	1:C:253:THR:HB	1.96	0.47
1:C:39:ALA:O	1:C:47:LEU:HD22	2.14	0.47
1:C:125:SER:OG	1:C:126:MET:N	2.48	0.47
1:F:126:MET:N	1:F:127:GLY:HA2	2.28	0.47
1:B:164:LYS:CD	1:B:165:GLU:N	2.73	0.47
1:E:190:ARG:HB2	1:E:227:GLY:C	2.33	0.47
1:A:14:ARG:HD2	1:A:253:THR:HB	1.97	0.47
1:A:135:SER:HA	1:A:232:MET:HA	1.93	0.47
1:E:274:ILE:CD1	1:E:279:TYR:CA	2.92	0.47
1:F:-5:VAL:HG23	1:F:-4:PRO:HD2	1.97	0.47
1:B:80:LYS:HE2	1:B:80:LYS:CA	2.44	0.47
1:D:70:LEU:HD23	1:D:118:LEU:HD23	1.96	0.47
1:C:134:ARG:NH1	1:C:233:GLU:OE1	2.47	0.47
1:F:126:MET:HE2	1:F:127:GLY:O	2.15	0.46
1:D:14:ARG:HD2	1:D:253:THR:HB	1.97	0.46
1:E:21:GLY:O	1:E:247:ILE:HD13	2.15	0.46
1:A:119:LEU:CA	1:A:120:GLU:HB2	2.45	0.46
1:E:184:ALA:O	1:E:185:ALA:C	2.54	0.46
1:A:135:SER:CB	1:A:232:MET:CB	2.86	0.46
1:A:274:ILE:CD1	1:A:280:LEU:N	2.78	0.46
1:F:80:LYS:HE2	1:F:80:LYS:CA	2.46	0.46
1:C:21:GLY:O	1:C:247:ILE:HD13	2.16	0.46
1:F:130:GLU:HB3	1:F:131:MET:CB	2.46	0.46
1:A:21:GLY:O	1:A:247:ILE:HD13	2.16	0.46
1:A:136:THR:HG23	1:A:261:LYS:CA	2.42	0.46
1:E:274:ILE:CG2	1:E:277:VAL:HG23	2.45	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:124:GLU:HB2	1:F:125:SER:CB	2.46	0.46
1:A:275:ASP:OD1	1:A:277:VAL:HG22	2.15	0.46
1:B:21:GLY:O	1:B:247:ILE:HD13	2.16	0.46
1:E:274:ILE:HD12	1:E:279:TYR:HA	1.95	0.46
1:F:14:ARG:HD2	1:F:253:THR:HB	1.97	0.46
1:A:105:ASN:O	1:A:106:PRO:C	2.54	0.45
1:F:184:ALA:HA	1:F:185:ALA:HA	1.64	0.45
1:B:14:ARG:HD2	1:B:253:THR:HB	1.98	0.45
1:B:131:MET:O	1:B:132:ASP:HB2	2.16	0.45
1:E:64:ARG:CG	1:E:65:ASN:N	2.75	0.45
1:E:68:LEU:HD22	1:E:92:HIS:CD2	2.52	0.45
1:A:96:SER:O	1:A:98:VAL:N	2.50	0.45
1:B:119:LEU:CA	1:B:120:GLU:HB2	2.46	0.45
1:E:30:CYS:C	1:E:62:CYS:SG	2.95	0.45
1:E:92:HIS:CE1	1:E:93:ASP:O	2.70	0.45
1:E:132:ASP:O	1:E:133:TYR:CB	2.65	0.45
1:D:10:SER:HB2	1:D:14:ARG:NH1	2.32	0.45
1:B:4:ALA:HA	1:B:59:LYS:O	2.17	0.45
1:D:106:PRO:O	1:D:108:LYS:HD3	2.17	0.45
1:D:21:GLY:O	1:D:247:ILE:HD13	2.17	0.45
1:D:131:MET:O	1:D:133:TYR:N	2.50	0.45
1:E:14:ARG:HD2	1:E:253:THR:HB	1.99	0.45
1:F:93:ASP:C	1:F:95:ASP:O	2.56	0.45
1:E:190:ARG:HB2	1:E:227:GLY:HA3	1.99	0.45
1:A:-4:PRO:HD2	1:A:-2:GLY:HA3	1.99	0.44
1:C:-1:SER:HB3	1:C:0:HIS:CD2	2.53	0.44
1:C:89:SER:OG	1:C:91:ARG:NH2	2.49	0.44
1:E:106:PRO:O	1:E:108:LYS:HD3	2.18	0.44
1:B:48:VAL:HG21	1:B:282:TYR:HE1	1.82	0.44
1:E:63:GLY:O	1:E:64:ARG:CG	2.66	0.44
1:F:38:GLN:OE1	1:F:127:GLY:O	2.35	0.44
1:A:119:LEU:HA	1:A:120:GLU:HB2	1.99	0.44
1:A:164:LYS:O	1:A:165:GLU:HG2	2.18	0.44
1:B:5:GLN:O	1:B:58:VAL:HG22	2.18	0.44
1:E:63:GLY:O	1:E:64:ARG:CB	2.65	0.44
1:F:188:SER:HA	1:F:189:ASP:HA	1.68	0.44
1:A:165:GLU:CG	1:A:166:GLY:N	2.81	0.44
1:B:1:MET:O	1:B:92:HIS:O	2.36	0.44
1:C:243:ARG:O	1:C:247:ILE:HG13	2.18	0.44
1:F:2:LEU:HD12	1:F:3:GLU:N	2.32	0.44
1:B:243:ARG:O	1:B:247:ILE:HG13	2.18	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:243:ARG:O	1:D:247:ILE:HG13	2.17	0.44
1:E:132:ASP:O	1:E:133:TYR:HB3	2.18	0.44
1:F:-2:GLY:HA3	1:F:-1:SER:HA	1.67	0.44
1:D:185:ALA:HB1	1:D:186:GLY:HA3	2.00	0.43
1:E:1:MET:SD	1:E:1:MET:N	2.92	0.43
1:E:51:LEU:HD23	1:E:279:TYR:HE1	1.81	0.43
1:E:243:ARG:O	1:E:247:ILE:HG13	2.18	0.43
1:F:22:LEU:HD12	1:F:48:VAL:CG2	2.48	0.43
1:A:164:LYS:O	1:A:165:GLU:CB	2.66	0.43
1:B:106:PRO:O	1:B:108:LYS:HD3	2.18	0.43
1:C:131:MET:SD	1:C:131:MET:N	2.76	0.43
1:D:105:ASN:OD1	1:D:110:ARG:HB3	2.19	0.43
1:A:165:GLU:CG	1:A:166:GLY:H	2.32	0.43
1:E:148:VAL:O	1:E:152:GLN:HG3	2.19	0.43
1:F:243:ARG:O	1:F:247:ILE:HG13	2.18	0.43
1:A:10:SER:HB2	1:A:14:ARG:NH1	2.33	0.43
1:C:96:SER:O	1:C:98:VAL:N	2.52	0.43
1:F:95:ASP:OD1	1:F:95:ASP:N	2.51	0.43
1:C:148:VAL:O	1:C:152:GLN:HG3	2.19	0.43
1:C:188:SER:HA	1:C:227:GLY:HA2	2.01	0.43
1:D:22:LEU:HD12	1:D:48:VAL:CG2	2.49	0.43
1:F:130:GLU:HB3	1:F:131:MET:HB3	2.00	0.43
1:D:97:ASP:O	1:D:118:LEU:HB2	2.18	0.43
1:F:106:PRO:O	1:F:108:LYS:HD3	2.19	0.43
1:B:48:VAL:HG21	1:B:282:TYR:CE1	2.54	0.43
1:C:22:LEU:HD12	1:C:48:VAL:CG2	2.49	0.43
1:E:10:SER:HB2	1:E:14:ARG:NH1	2.33	0.43
1:E:249:ALA:HA	1:E:272:TYR:OH	2.19	0.43
1:F:47:LEU:HG	1:F:126:MET:HE3	1.95	0.43
1:A:-4:PRO:N	1:A:-3:ARG:CA	2.81	0.42
1:B:130:GLU:N	1:B:130:GLU:CD	2.72	0.42
1:B:130:GLU:H	1:B:131:MET:HB3	1.83	0.42
1:C:2:LEU:HD12	1:C:3:GLU:N	2.34	0.42
1:E:-4:PRO:O	1:E:-3:ARG:HB2	2.19	0.42
1:A:0:HIS:CB	1:A:94:ASP:CB	2.80	0.42
1:A:135:SER:CB	1:A:232:MET:HB2	2.48	0.42
1:D:248:PHE:CD1	1:D:282:TYR:CG	3.07	0.42
1:E:134:ARG:NH2	1:E:236:ILE:HD12	2.14	0.42
1:E:188:SER:OG	1:E:189:ASP:HB3	2.20	0.42
1:E:120:GLU:HA	1:E:120:GLU:OE1	2.19	0.42
1:A:90:LEU:HD11	1:A:99:VAL:CG2	2.50	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:243:ARG:O	1:A:247:ILE:HG13	2.19	0.42
1:B:148:VAL:O	1:B:152:GLN:HG3	2.20	0.42
1:A:134:ARG:NH2	1:A:233:GLU:OE1	2.53	0.42
1:A:190:ARG:NH2	1:B:109:THR:OG1	2.52	0.42
1:C:10:SER:HB2	1:C:14:ARG:NH1	2.35	0.42
1:D:148:VAL:O	1:D:152:GLN:HG3	2.19	0.42
1:D:274:ILE:HD12	1:D:279:TYR:HA	2.02	0.42
1:E:190:ARG:HB3	1:E:257:ARG:HH12	1.84	0.42
1:F:189:ASP:O	1:F:190:ARG:HB2	2.19	0.42
1:B:164:LYS:HZ2	1:B:165:GLU:CG	2.32	0.42
1:C:190:ARG:HB3	1:C:190:ARG:HH11	1.85	0.42
1:A:-3:ARG:C	1:A:-3:ARG:CD	2.86	0.41
1:A:126:MET:HG3	1:A:127:GLY:C	2.41	0.41
1:B:10:SER:HB2	1:B:14:ARG:NH1	2.35	0.41
1:B:164:LYS:NZ	1:B:165:GLU:CD	2.73	0.41
1:C:40:MET:HB2	1:C:126:MET:HG3	2.01	0.41
1:E:90:LEU:HD11	1:E:99:VAL:CG2	2.50	0.41
1:C:39:ALA:HA	1:C:126:MET:CE	2.51	0.41
1:D:51:LEU:HD12	1:D:52:LEU:N	2.35	0.41
1:A:38:GLN:OE1	1:A:127:GLY:O	2.39	0.41
1:B:119:LEU:CB	1:B:120:GLU:HB2	2.49	0.41
1:B:134:ARG:HH12	1:B:234:GLU:HB2	1.86	0.41
1:C:90:LEU:HD11	1:C:99:VAL:CG2	2.51	0.41
1:F:148:VAL:O	1:F:152:GLN:HG3	2.20	0.41
1:A:148:VAL:O	1:A:152:GLN:HG3	2.20	0.41
1:B:249:ALA:HA	1:B:272:TYR:OH	2.21	0.41
1:C:122:GLU:O	1:C:124:GLU:HG2	2.20	0.41
1:C:274:ILE:HD12	1:C:279:TYR:HA	2.03	0.41
1:E:248:PHE:CD1	1:E:282:TYR:CG	3.09	0.41
1:F:184:ALA:C	1:F:228:VAL:O	2.58	0.41
1:E:133:TYR:CD1	1:E:133:TYR:C	2.93	0.41
1:F:10:SER:HB2	1:F:14:ARG:NH1	2.36	0.41
1:A:248:PHE:CD1	1:A:282:TYR:CG	3.09	0.41
1:E:-1:SER:CB	1:E:0:HIS:HB3	2.51	0.41
1:A:22:LEU:HD12	1:A:48:VAL:CG2	2.50	0.40
1:B:22:LEU:HD12	1:B:22:LEU:HA	1.91	0.40
1:B:30:CYS:HA	1:B:35:LEU:HD23	2.04	0.40
1:E:274:ILE:HB	1:E:277:VAL:HG23	2.03	0.40
1:A:-3:ARG:HH21	1:A:64:ARG:HH22	1.68	0.40
1:B:90:LEU:HD11	1:B:99:VAL:CG2	2.51	0.40
1:F:95:ASP:O	1:F:96:SER:CB	2.63	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:5:GLN:NE2	1:B:87:SER:CB	2.80	0.40
1:F:184:ALA:HA	1:F:187:VAL:HG12	1.96	0.40
1:B:164:LYS:HD3	1:B:165:GLU:HG3	2.02	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:A:-3:ARG:NH1	1:B:120:GLU:OE2[4_544]	2.01	0.19

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	256/300 (85%)	227 (89%)	20 (8%)	9 (4%)	3	17
1	B	257/300 (86%)	232 (90%)	22 (9%)	3 (1%)	13	43
1	C	256/300 (85%)	233 (91%)	20 (8%)	3 (1%)	13	43
1	D	257/300 (86%)	234 (91%)	20 (8%)	3 (1%)	13	43
1	E	255/300 (85%)	228 (89%)	21 (8%)	6 (2%)	6	26
1	F	257/300 (86%)	223 (87%)	30 (12%)	4 (2%)	9	36
All	All	1538/1800 (85%)	1377 (90%)	133 (9%)	28 (2%)	8	33

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	97	ASP
1	A	165	GLU
1	B	97	ASP
1	C	97	ASP
1	D	94	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	95	ASP
1	E	97	ASP
1	E	275	ASP
1	F	0	HIS
1	A	126	MET
1	A	128	ILE
1	C	123	ALA
1	F	-4	PRO
1	A	0	HIS
1	D	185	ALA
1	E	64	ARG
1	E	125	SER
1	F	96	SER
1	A	125	SER
1	E	95	ASP
1	F	184	ALA
1	B	132	ASP
1	E	1	MET
1	A	-4	PRO
1	A	129	PRO
1	A	184	ALA
1	C	107	GLU
1	B	63	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	226/257 (88%)	200 (88%)	26 (12%)	5 21
1	B	226/257 (88%)	201 (89%)	25 (11%)	6 22
1	C	226/257 (88%)	195 (86%)	31 (14%)	3 15
1	D	226/257 (88%)	200 (88%)	26 (12%)	5 21
1	E	225/257 (88%)	199 (88%)	26 (12%)	5 20
1	F	226/257 (88%)	199 (88%)	27 (12%)	5 20

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1355/1542 (88%)	1194 (88%)	161 (12%)	5 20

All (161) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LEU
1	A	19	ILE
1	A	22	LEU
1	A	41	ASP
1	A	50	MET
1	A	53	ARG
1	A	55	ASP
1	A	64	ARG
1	A	100	THR
1	A	103	SER
1	A	108	LYS
1	A	109	THR
1	A	116	LEU
1	A	124	GLU
1	A	126	MET
1	A	135	SER
1	A	137	VAL
1	A	164	LYS
1	A	170	SER
1	A	183	GLN
1	A	188	SER
1	A	228	VAL
1	A	231	THR
1	A	266	SER
1	A	270	VAL
1	A	289	ASP
1	B	-3	ARG
1	B	0	HIS
1	B	2	LEU
1	B	19	ILE
1	B	29	ASP
1	B	41	ASP
1	B	47	LEU
1	B	53	ARG
1	B	64	ARG
1	B	80	LYS
1	B	94	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	100	THR
1	B	103	SER
1	B	108	LYS
1	B	109	THR
1	B	116	LEU
1	B	126	MET
1	B	137	VAL
1	B	164	LYS
1	B	170	SER
1	B	183	GLN
1	B	189	ASP
1	B	231	THR
1	B	266	SER
1	B	270	VAL
1	C	-6	LEU
1	C	-3	ARG
1	C	-1	SER
1	C	1	MET
1	C	3	GLU
1	C	19	ILE
1	C	22	LEU
1	C	24	ASN
1	C	29	ASP
1	C	41	ASP
1	C	45	VAL
1	C	50	MET
1	C	53	ARG
1	C	100	THR
1	C	103	SER
1	C	107	GLU
1	C	108	LYS
1	C	109	THR
1	C	116	LEU
1	C	126	MET
1	C	131	MET
1	C	134	ARG
1	C	137	VAL
1	C	163	SER
1	C	164	LYS
1	C	170	SER
1	C	190	ARG
1	C	231	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	266	SER
1	C	270	VAL
1	C	276	ASN
1	D	2	LEU
1	D	22	LEU
1	D	29	ASP
1	D	41	ASP
1	D	50	MET
1	D	53	ARG
1	D	65	ASN
1	D	84	SER
1	D	97	ASP
1	D	100	THR
1	D	103	SER
1	D	105	ASN
1	D	108	LYS
1	D	109	THR
1	D	116	LEU
1	D	120	GLU
1	D	137	VAL
1	D	163	SER
1	D	164	LYS
1	D	170	SER
1	D	188	SER
1	D	190	ARG
1	D	231	THR
1	D	266	SER
1	D	270	VAL
1	D	276	ASN
1	E	-3	ARG
1	E	-1	SER
1	E	1	MET
1	E	2	LEU
1	E	19	ILE
1	E	22	LEU
1	E	29	ASP
1	E	41	ASP
1	E	50	MET
1	E	53	ARG
1	E	80	LYS
1	E	100	THR
1	E	103	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	E	108	LYS
1	E	109	THR
1	E	116	LEU
1	E	120	GLU
1	E	131	MET
1	E	132	ASP
1	E	137	VAL
1	E	163	SER
1	E	170	SER
1	E	188	SER
1	E	231	THR
1	E	266	SER
1	E	270	VAL
1	F	-3	ARG
1	F	19	ILE
1	F	22	LEU
1	F	29	ASP
1	F	41	ASP
1	F	50	MET
1	F	53	ARG
1	F	64	ARG
1	F	80	LYS
1	F	100	THR
1	F	103	SER
1	F	108	LYS
1	F	109	THR
1	F	116	LEU
1	F	126	MET
1	F	131	MET
1	F	137	VAL
1	F	163	SER
1	F	164	LYS
1	F	165	GLU
1	F	170	SER
1	F	187	VAL
1	F	228	VAL
1	F	245	MET
1	F	266	SER
1	F	270	VAL
1	F	276	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	HIS
1	A	276	ASN
1	B	5	GLN
1	B	7	GLN
1	C	7	GLN
1	C	49	HIS
1	D	7	GLN
1	D	65	ASN
1	E	7	GLN
1	E	44	HIS
1	E	276	ASN
1	F	7	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	260/300 (86%)	0.33	24 (9%)	9	5	70, 110, 188, 230	0
1	B	261/300 (87%)	0.19	12 (4%)	32	20	56, 92, 163, 247	0
1	C	260/300 (86%)	0.13	15 (5%)	23	14	60, 95, 167, 221	0
1	D	261/300 (87%)	0.05	11 (4%)	36	23	69, 104, 161, 224	0
1	E	259/300 (86%)	0.17	14 (5%)	25	16	69, 104, 181, 241	0
1	F	261/300 (87%)	0.11	12 (4%)	32	20	56, 88, 160, 229	0
All	All	1562/1800 (86%)	0.17	88 (5%)	24	15	56, 99, 173, 247	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	187	VAL	12.3
1	A	189	ASP	11.9
1	C	188	SER	11.8
1	B	188	SER	11.5
1	F	190	ARG	11.1
1	A	187	VAL	10.2
1	B	187	VAL	10.1
1	D	188	SER	9.5
1	D	186	GLY	8.9
1	F	189	ASP	8.7
1	E	188	SER	8.6
1	A	190	ARG	8.2
1	F	127	GLY	8.1
1	A	188	SER	8.0
1	C	189	ASP	7.6
1	D	187	VAL	7.6
1	F	185	ALA	7.5
1	A	-6	LEU	7.4
1	A	186	GLY	6.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	187	VAL	6.8
1	B	189	ASP	6.7
1	A	-1	SER	6.4
1	C	-6	LEU	6.2
1	F	188	SER	6.1
1	F	187	VAL	6.1
1	C	186	GLY	5.9
1	E	186	GLY	5.9
1	B	190	ARG	5.9
1	F	186	GLY	5.7
1	B	186	GLY	5.6
1	E	-5	VAL	5.6
1	A	-4	PRO	5.4
1	A	-2	GLY	5.2
1	F	126	MET	5.1
1	E	189	ASP	5.0
1	D	190	ARG	4.9
1	A	0	HIS	4.9
1	B	-3	ARG	4.8
1	C	-2	GLY	4.7
1	C	185	ALA	4.7
1	A	-5	VAL	4.7
1	C	-3	ARG	4.6
1	D	185	ALA	4.5
1	D	189	ASP	4.3
1	E	288	VAL	4.3
1	A	185	ALA	4.2
1	E	0	HIS	4.2
1	E	-1	SER	4.2
1	B	-1	SER	4.1
1	B	-4	PRO	4.0
1	C	289	ASP	4.0
1	D	-3	ARG	3.9
1	E	-2	GLY	3.8
1	E	289	ASP	3.8
1	F	289	ASP	3.8
1	E	-3	ARG	3.8
1	D	289	ASP	3.6
1	B	185	ALA	3.6
1	C	-4	PRO	3.6
1	D	238	LEU	3.5
1	C	-5	VAL	3.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	2	LEU	3.3
1	E	185	ALA	3.2
1	A	68	LEU	3.0
1	C	288	VAL	3.0
1	F	-7	GLY	2.9
1	A	238	LEU	2.8
1	E	276	ASN	2.7
1	D	288	VAL	2.7
1	B	-7	GLY	2.7
1	A	129	PRO	2.6
1	A	-3	ARG	2.6
1	F	184	ALA	2.6
1	F	125	SER	2.6
1	A	28	PHE	2.5
1	D	262	PHE	2.5
1	C	190	ARG	2.4
1	A	231	THR	2.4
1	A	127	GLY	2.4
1	C	30	CYS	2.4
1	A	262	PHE	2.2
1	A	227	GLY	2.2
1	A	1	MET	2.2
1	E	167	VAL	2.2
1	C	175	VAL	2.2
1	B	235	PRO	2.1
1	A	134	ARG	2.1
1	A	160	ILE	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.