Full wwPDB X-ray Structure Validation Report (i)

Oct 15, 2023 – 10:05 AM EDT

PDB ID	:	1IOI
Title	:	x-ray crystalline structures of pyrrolidone carboxyl peptidase from a hyperthermophile, pyrococcus furiosus, and its cys-free mutant
Authors	:	Tanaka, H.; Chinami, M.; Ota, M.; Tsukihara, T.; Yutani, K.
Deposited on	:	2001-03-09
Resolution	:	2.70 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 (i)) were used in the production of this report:

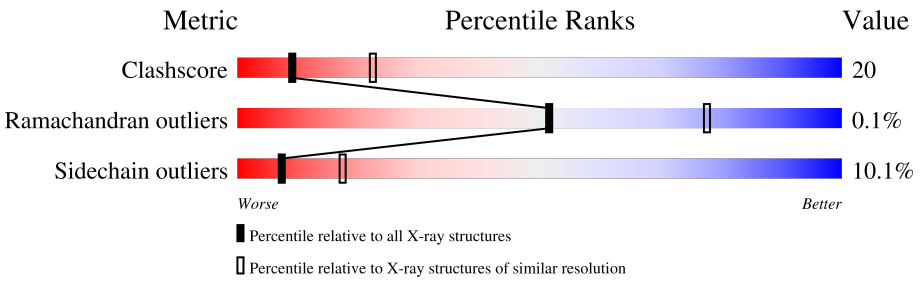
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.70 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))
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)

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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	208	55% 36% 8% •
1	B	208	58% 34% 7% •
1	C	208	55% 37% 7% •
1	D	208	49% 41% 8% •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 6544 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 PYRROLIDONE CARBOXYL PEPTIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	208	1605	1040	263	296	6	0	0	0
1	B	208	1605	1040	263	296	6	0	0	0
1	C	208	1605	1040	263	296	6	0	0	0
1	D	208	1605	1040	263	296	6	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	142	SER	CYS	engineered mutation	UNP O73944
A	188	SER	CYS	engineered mutation	UNP O73944
B	142	SER	CYS	engineered mutation	UNP O73944
B	188	SER	CYS	engineered mutation	UNP O73944
C	142	SER	CYS	engineered mutation	UNP O73944
C	188	SER	CYS	engineered mutation	UNP O73944
D	142	SER	CYS	engineered mutation	UNP O73944
D	188	SER	CYS	engineered mutation	UNP O73944

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	30	30	30	0	0
2	B	34	34	34	0	0
2	C	28	28	28	0	0
2	D	32	32	32	0	0

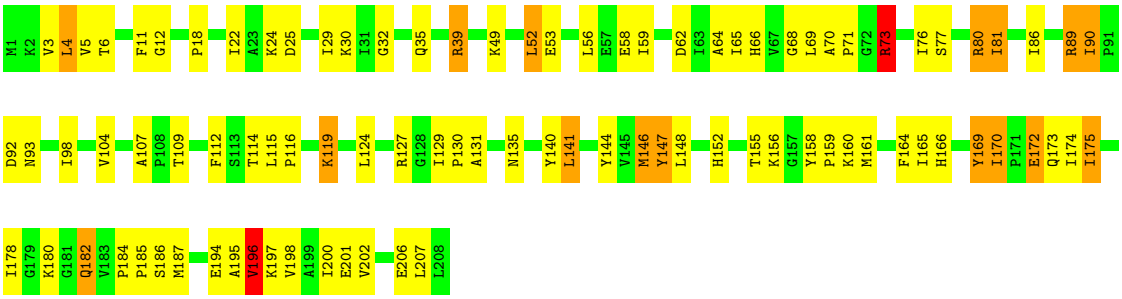
3 Residue-property plots (i)

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. 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. 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.

Note EDS was not executed.

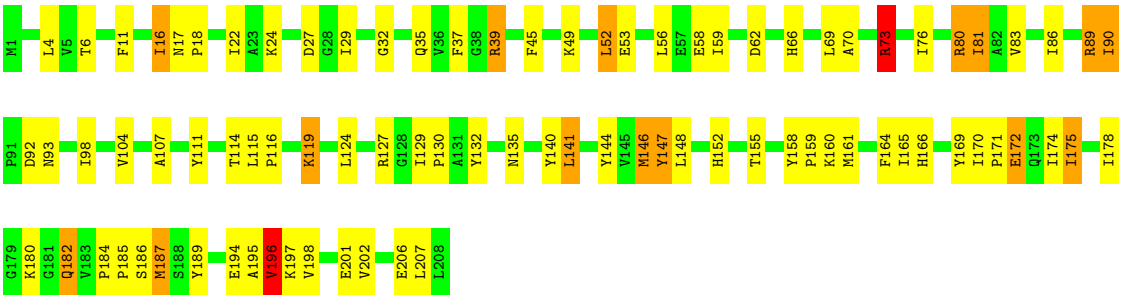
• Molecule 1: PYRROLIDONE CARBOXYL PEPTIDASE

Chain A:



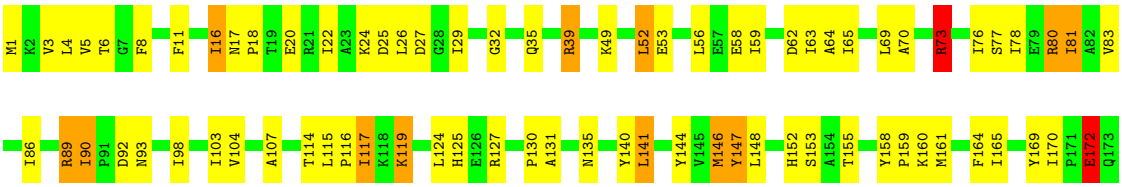
• Molecule 1: PYRROLIDONE CARBOXYL PEPTIDASE

Chain B:



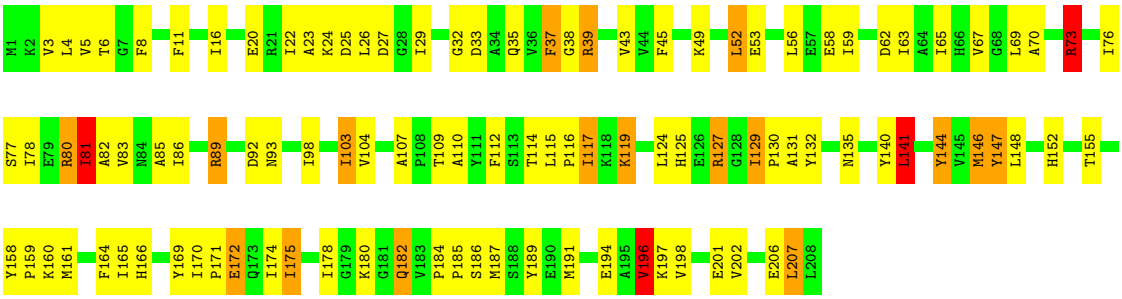
• Molecule 1: PYRROLIDONE CARBOXYL PEPTIDASE

Chain C:



• Molecule 1: PYRROLIDONE CARBOXYL PEPTIDASE

Chain D:



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	49.00Å 105.80Å 105.80Å 90.00° 96.20° 90.00°	Depositor
Resolution (Å)	10.00 – 2.70	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-2.70)	Depositor
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.210 , 0.243	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6544	wwPDB-VP
Average B, all atoms (Å^2)	26.0	wwPDB-VP

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	1.34	5/1639 (0.3%)	1.87	34/2216 (1.5%)
1	B	1.27	5/1639 (0.3%)	1.85	36/2216 (1.6%)
1	C	1.27	5/1639 (0.3%)	1.85	41/2216 (1.9%)
1	D	1.29	4/1639 (0.2%)	1.89	43/2216 (1.9%)
All	All	1.30	19/6556 (0.3%)	1.87	154/8864 (1.7%)

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	1
1	B	0	1
1	D	0	1
All	All	0	3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	196	VAL	CA-CB	7.78	1.71	1.54
1	B	202	VAL	CA-CB	6.80	1.69	1.54
1	A	196	VAL	CA-CB	6.77	1.69	1.54
1	D	202	VAL	CA-CB	6.75	1.69	1.54
1	C	196	VAL	CA-CB	6.74	1.69	1.54
1	C	202	VAL	CA-CB	6.60	1.68	1.54
1	A	202	VAL	CA-CB	6.46	1.68	1.54
1	D	172	GLU	CB-CG	6.33	1.64	1.52
1	D	175	ILE	CA-CB	6.15	1.69	1.54
1	C	172	GLU	CB-CG	5.98	1.63	1.52
1	A	172	GLU	CG-CD	5.93	1.60	1.51
1	A	172	GLU	CB-CG	5.84	1.63	1.52
1	B	196	VAL	CA-CB	5.81	1.67	1.54

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	175	ILE	CA-CB	5.55	1.67	1.54
1	B	175	ILE	CA-CB	5.54	1.67	1.54
1	B	111	TYR	CE2-CZ	5.51	1.45	1.38
1	C	175	ILE	CA-CB	5.31	1.67	1.54
1	C	172	GLU	CG-CD	5.16	1.59	1.51
1	B	66	HIS	CD2-NE2	-5.12	1.26	1.38

All (154) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	73	ARG	NE-CZ-NH1	15.33	127.97	120.30
1	A	73	ARG	NE-CZ-NH1	14.70	127.65	120.30
1	D	73	ARG	NE-CZ-NH1	13.97	127.28	120.30
1	B	73	ARG	NE-CZ-NH1	13.89	127.25	120.30
1	B	80	ARG	NE-CZ-NH2	-13.21	113.69	120.30
1	D	80	ARG	NE-CZ-NH1	12.80	126.70	120.30
1	B	80	ARG	NE-CZ-NH1	12.25	126.42	120.30
1	C	80	ARG	NE-CZ-NH2	-12.18	114.21	120.30
1	C	80	ARG	NE-CZ-NH1	11.93	126.27	120.30
1	A	80	ARG	NE-CZ-NH1	11.76	126.18	120.30
1	D	80	ARG	NE-CZ-NH2	-11.72	114.44	120.30
1	D	127	ARG	NE-CZ-NH2	-11.42	114.59	120.30
1	B	127	ARG	NE-CZ-NH2	-10.85	114.88	120.30
1	A	80	ARG	NE-CZ-NH2	-10.64	114.98	120.30
1	A	127	ARG	NE-CZ-NH2	-10.50	115.05	120.30
1	C	127	ARG	NE-CZ-NH2	-10.40	115.10	120.30
1	A	39	ARG	NE-CZ-NH1	9.79	125.19	120.30
1	C	140	TYR	CB-CG-CD2	-9.76	115.15	121.00
1	A	140	TYR	CB-CG-CD2	-9.47	115.32	121.00
1	B	146	MET	CG-SD-CE	9.26	115.02	100.20
1	D	140	TYR	CB-CG-CD2	-9.24	115.46	121.00
1	C	140	TYR	CB-CG-CD1	8.92	126.35	121.00
1	D	185	PRO	CA-C-N	-8.83	97.78	117.20
1	D	114	THR	CA-CB-CG2	-8.63	100.32	112.40
1	A	114	THR	CA-CB-CG2	-8.59	100.38	112.40
1	B	39	ARG	NE-CZ-NH1	8.52	124.56	120.30
1	B	185	PRO	CA-C-N	-8.49	98.52	117.20
1	A	92	ASP	CB-CG-OD2	8.42	125.88	118.30
1	B	114	THR	CA-CB-CG2	-8.31	100.76	112.40
1	A	185	PRO	CA-C-N	-8.24	99.07	117.20
1	A	73	ARG	NE-CZ-NH2	-8.21	116.20	120.30
1	C	114	THR	CA-CB-CG2	-8.12	101.03	112.40

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	140	TYR	CB-CG-CD2	-8.08	116.15	121.00
1	A	80	ARG	CB-CG-CD	-7.94	90.95	111.60
1	D	39	ARG	NE-CZ-NH1	7.89	124.25	120.30
1	B	52	LEU	CA-CB-CG	7.89	133.44	115.30
1	D	52	LEU	CA-CB-CG	7.83	133.30	115.30
1	C	52	LEU	CA-CB-CG	7.81	133.26	115.30
1	D	146	MET	CG-SD-CE	7.76	112.62	100.20
1	C	185	PRO	CA-C-N	-7.75	100.16	117.20
1	C	146	MET	CG-SD-CE	7.70	112.52	100.20
1	C	39	ARG	NE-CZ-NH1	7.62	124.11	120.30
1	A	53	GLU	CA-CB-CG	-7.38	97.17	113.40
1	D	80	ARG	CB-CG-CD	-7.33	92.53	111.60
1	A	197	LYS	CB-CG-CD	-7.27	92.69	111.60
1	D	132	TYR	CB-CG-CD1	-7.24	116.66	121.00
1	D	140	TYR	CB-CG-CD1	7.20	125.32	121.00
1	C	147	TYR	CB-CG-CD2	-7.19	116.69	121.00
1	B	140	TYR	CB-CG-CD1	7.19	125.31	121.00
1	D	197	LYS	CB-CG-CD	-7.04	93.30	111.60
1	B	53	GLU	CA-CB-CG	-6.99	98.01	113.40
1	B	197	LYS	CB-CG-CD	-6.97	93.47	111.60
1	C	53	GLU	CA-CB-CG	-6.96	98.09	113.40
1	D	53	GLU	CA-CB-CG	-6.86	98.31	113.40
1	D	24	LYS	CA-CB-CG	-6.82	98.41	113.40
1	A	52	LEU	CA-CB-CG	6.78	130.89	115.30
1	C	197	LYS	CB-CG-CD	-6.77	94.00	111.60
1	B	155	THR	CA-CB-CG2	6.76	121.86	112.40
1	B	58	GLU	CA-CB-CG	6.73	128.21	113.40
1	C	155	THR	CA-CB-CG2	6.69	121.77	112.40
1	A	147	TYR	CB-CG-CD2	-6.63	117.02	121.00
1	D	155	THR	CA-CB-CG2	6.63	121.68	112.40
1	B	169	TYR	CB-CG-CD1	-6.62	117.03	121.00
1	A	169	TYR	CB-CG-CD1	-6.59	117.05	121.00
1	C	80	ARG	CB-CG-CD	-6.51	94.67	111.60
1	D	147	TYR	CB-CG-CD2	-6.48	117.11	121.00
1	B	80	ARG	CB-CG-CD	-6.48	94.76	111.60
1	D	185	PRO	O-C-N	6.45	133.01	122.70
1	A	24	LYS	CA-CB-CG	-6.37	99.38	113.40
1	D	73	ARG	NE-CZ-NH2	-6.37	117.11	120.30
1	D	127	ARG	NE-CZ-NH1	6.37	123.48	120.30
1	D	92	ASP	CB-CG-OD2	6.32	123.98	118.30
1	C	92	ASP	CB-CG-OD2	6.27	123.94	118.30
1	A	140	TYR	CB-CG-CD1	6.26	124.76	121.00

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	155	THR	CA-CB-CG2	6.22	121.12	112.40
1	C	58	GLU	CA-CB-CG	6.18	126.99	113.40
1	D	25	ASP	CB-CG-OD2	6.10	123.79	118.30
1	B	172	GLU	CA-CB-CG	6.08	126.76	113.40
1	D	58	GLU	CA-CB-CG	6.07	126.75	113.40
1	B	89	ARG	CB-CG-CD	6.06	127.36	111.60
1	A	58	GLU	CA-CB-CG	6.04	126.69	113.40
1	B	147	TYR	CB-CG-CD2	-6.02	117.39	121.00
1	B	92	ASP	CB-CG-OD2	6.01	123.71	118.30
1	A	146	MET	CG-SD-CE	5.97	109.76	100.20
1	D	32	GLY	CA-C-N	-5.95	104.12	117.20
1	A	185	PRO	O-C-N	5.92	132.18	122.70
1	B	73	ARG	NE-CZ-NH2	-5.89	117.36	120.30
1	B	24	LYS	CA-CB-CG	-5.88	100.46	113.40
1	C	24	LYS	CA-CB-CG	-5.87	100.49	113.40
1	C	172	GLU	CA-CB-CG	5.86	126.30	113.40
1	D	169	TYR	CB-CG-CD1	-5.85	117.49	121.00
1	C	89	ARG	CB-CG-CD	5.84	126.78	111.60
1	D	114	THR	CA-CB-OG1	5.80	121.18	109.00
1	B	73	ARG	CD-NE-CZ	5.80	131.72	123.60
1	C	73	ARG	NE-CZ-NH2	-5.79	117.40	120.30
1	D	89	ARG	CB-CG-CD	5.75	126.55	111.60
1	B	185	PRO	O-C-N	5.71	131.84	122.70
1	A	89	ARG	CB-CG-CD	5.69	126.41	111.60
1	C	114	THR	CA-CB-OG1	5.69	120.95	109.00
1	D	172	GLU	CA-CB-CG	5.67	125.86	113.40
1	A	32	GLY	CA-C-N	-5.66	104.76	117.20
1	C	155	THR	N-CA-CB	-5.62	99.61	110.30
1	A	141	LEU	CA-CB-CG	-5.60	102.42	115.30
1	A	25	ASP	CB-CG-OD1	-5.59	113.27	118.30
1	B	114	THR	CA-CB-OG1	5.58	120.71	109.00
1	B	141	LEU	CA-CB-CG	-5.57	102.49	115.30
1	C	169	TYR	CB-CG-CD1	-5.54	117.68	121.00
1	A	155	THR	N-CA-CB	-5.52	99.82	110.30
1	C	127	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	B	201	GLU	CA-C-N	5.50	129.29	117.20
1	C	32	GLY	CA-C-N	-5.48	105.14	117.20
1	B	127	ARG	NE-CZ-NH1	5.47	123.03	120.30
1	D	27	ASP	CB-CG-OD1	-5.47	113.38	118.30
1	A	114	THR	CA-CB-OG1	5.46	120.47	109.00
1	D	89	ARG	NE-CZ-NH1	5.45	123.03	120.30
1	D	155	THR	N-CA-CB	-5.45	99.94	110.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	169	TYR	CG-CD2-CE2	-5.45	116.94	121.30
1	C	89	ARG	NE-CZ-NH1	5.43	123.02	120.30
1	A	172	GLU	CA-CB-CG	5.43	125.35	113.40
1	B	27	ASP	CB-CG-OD2	5.41	123.17	118.30
1	C	73	ARG	NH1-CZ-NH2	-5.39	113.47	119.40
1	B	187	MET	CG-SD-CE	5.39	108.82	100.20
1	B	132	TYR	CB-CG-CD1	-5.39	117.77	121.00
1	B	155	THR	N-CA-CB	-5.38	100.08	110.30
1	C	1	MET	CG-SD-CE	-5.37	91.61	100.20
1	C	141	LEU	CA-CB-CG	-5.34	103.01	115.30
1	A	77	SER	CA-C-N	-5.34	105.46	117.20
1	A	201	GLU	CA-C-N	5.33	128.92	117.20
1	C	73	ARG	CD-NE-CZ	5.28	130.99	123.60
1	C	207	LEU	O-C-N	5.27	131.13	122.70
1	C	185	PRO	O-C-N	5.26	131.11	122.70
1	C	27	ASP	CB-CG-OD2	5.25	123.02	118.30
1	D	189	TYR	CB-CG-CD2	-5.24	117.85	121.00
1	B	201	GLU	CA-C-O	-5.20	109.17	120.10
1	D	141	LEU	CA-CB-CG	-5.12	103.51	115.30
1	D	110	ALA	CA-C-N	-5.12	105.94	117.20
1	A	73	ARG	CD-NE-CZ	5.11	130.76	123.60
1	C	201	GLU	CA-C-N	5.11	128.44	117.20
1	D	89	ARG	CA-CB-CG	5.10	124.62	113.40
1	C	153	SER	CA-CB-OG	5.10	124.96	111.20
1	C	77	SER	CA-C-N	-5.08	106.02	117.20
1	D	201	GLU	CA-C-N	5.08	128.37	117.20
1	C	189	TYR	CB-CG-CD2	-5.06	117.96	121.00
1	D	81	ILE	CB-CA-C	-5.06	101.48	111.60
1	D	207	LEU	O-C-N	5.06	130.80	122.70
1	B	32	GLY	CA-C-N	-5.05	106.09	117.20
1	D	37	PHE	CB-CG-CD1	-5.04	117.27	120.80
1	B	89	ARG	CA-CB-CG	5.04	124.49	113.40
1	C	25	ASP	CB-CG-OD1	-5.03	113.77	118.30
1	A	4	LEU	CB-CG-CD1	-5.02	102.46	111.00
1	D	77	SER	CA-C-N	-5.02	106.15	117.20
1	A	127	ARG	NE-CZ-NH1	5.01	122.81	120.30
1	D	25	ASP	CB-CG-OD1	-5.01	113.79	118.30
1	C	25	ASP	CB-CG-OD2	5.01	122.81	118.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	144	TYR	Sidechain
1	B	189	TYR	Sidechain
1	D	144	TYR	Sidechain

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	1605	0	1671	73	0
1	B	1605	0	1671	67	0
1	C	1605	0	1671	66	0
1	D	1605	0	1671	72	0
2	A	30	0	0	3	0
2	B	34	0	0	0	0
2	C	28	0	0	0	0
2	D	32	0	0	3	0
All	All	6544	0	6684	263	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 (263) 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:65:ILE:HD12	1:C:115:LEU:HD11	1.27	1.16
1:A:65:ILE:HD12	1:A:115:LEU:HD11	1.29	1.14
1:A:64:ALA:O	1:A:65:ILE:HD13	1.62	0.99
1:C:64:ALA:O	1:C:65:ILE:HD13	1.62	0.99
1:A:65:ILE:HD12	1:A:115:LEU:CD1	2.03	0.88
1:D:175:ILE:HA	1:D:178:ILE:HD12	1.56	0.88
1:C:175:ILE:HA	1:C:178:ILE:HD12	1.57	0.87
1:C:116:PRO:HB2	1:C:119:LYS:HD2	1.56	0.87
1:B:175:ILE:HA	1:B:178:ILE:HD12	1.57	0.87
1:C:90:ILE:HD13	1:C:90:ILE:H	1.39	0.87
1:A:90:ILE:H	1:A:90:ILE:HD13	1.40	0.87
1:C:65:ILE:HD12	1:C:115:LEU:CD1	2.05	0.86
1:B:90:ILE:HD13	1:B:90:ILE:H	1.41	0.86

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:116:PRO:HB2	1:D:119:LYS:HD2	1.56	0.85
1:B:76:ILE:HD12	1:B:129:ILE:CG2	2.06	0.85
1:A:116:PRO:HB2	1:A:119:LYS:HD2	1.57	0.84
1:C:103:ILE:HD12	1:C:144:TYR:HD1	1.44	0.83
1:A:76:ILE:HD12	1:A:129:ILE:CG2	2.09	0.83
1:B:171:PRO:HA	1:B:174:ILE:HD13	1.58	0.83
1:D:70:ALA:HB1	1:D:73:ARG:HD3	1.59	0.83
1:B:116:PRO:HB2	1:B:119:LYS:HD2	1.60	0.82
1:C:70:ALA:HB1	1:C:73:ARG:HD3	1.62	0.81
1:A:175:ILE:HA	1:A:178:ILE:HD12	1.61	0.81
1:A:70:ALA:HB1	1:A:73:ARG:HD3	1.62	0.81
1:A:65:ILE:CD1	1:A:115:LEU:HD11	2.09	0.79
1:B:76:ILE:HD12	1:B:129:ILE:HG22	1.64	0.78
1:D:129:ILE:HD12	1:D:191:MET:CE	2.13	0.77
1:B:76:ILE:HD13	1:B:195:ALA:CB	2.16	0.76
1:B:70:ALA:HB1	1:B:73:ARG:HD3	1.67	0.76
1:A:76:ILE:HD13	1:A:195:ALA:CB	2.15	0.75
1:A:65:ILE:HD11	1:A:115:LEU:CD2	2.17	0.75
1:C:65:ILE:CD1	1:C:115:LEU:HD11	2.14	0.74
1:C:65:ILE:HD11	1:C:115:LEU:HD21	1.68	0.74
1:A:65:ILE:HD11	1:A:115:LEU:HD21	1.67	0.74
1:C:65:ILE:HD11	1:C:115:LEU:CD2	2.20	0.71
1:B:174:ILE:HD12	1:B:174:ILE:N	2.06	0.71
1:A:76:ILE:HD12	1:A:129:ILE:HG22	1.74	0.69
1:B:81:ILE:HD13	1:C:83:VAL:HG11	1.73	0.69
1:A:170:ILE:HD13	1:A:173:GLN:OE1	1.94	0.68
1:A:56:LEU:HD22	1:A:159:PRO:HD3	1.76	0.68
1:D:129:ILE:HD12	1:D:191:MET:HE1	1.75	0.67
1:C:103:ILE:HD12	1:C:144:TYR:CD1	2.30	0.67
1:D:6:THR:HA	1:D:39:ARG:O	1.95	0.67
1:B:83:VAL:HG11	1:C:81:ILE:HD13	1.78	0.66
1:B:76:ILE:HD13	1:B:195:ALA:HB1	1.76	0.65
1:C:64:ALA:C	1:C:65:ILE:HD13	2.17	0.65
1:A:76:ILE:HD13	1:A:195:ALA:HB1	1.77	0.64
1:B:104:VAL:HB	1:B:107:ALA:CB	2.28	0.64
1:A:104:VAL:HB	1:A:107:ALA:HB3	1.80	0.64
1:B:174:ILE:HD11	1:B:186:SER:HB3	1.79	0.64
1:B:104:VAL:HB	1:B:107:ALA:HB3	1.79	0.64
1:B:174:ILE:CD1	1:B:186:SER:HB3	2.27	0.64
1:D:78:ILE:HD13	1:D:165:ILE:HG12	1.78	0.63
1:D:117:ILE:H	1:D:117:ILE:HD13	1.63	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:VAL:HB	1:A:107:ALA:CB	2.29	0.63
1:D:56:LEU:HD22	1:D:159:PRO:HD3	1.81	0.63
1:D:49:LYS:HG3	1:D:148:LEU:HD13	1.79	0.63
1:A:81:ILE:HD13	1:D:83:VAL:HG11	1.81	0.63
1:A:6:THR:HA	1:A:39:ARG:O	1.99	0.62
1:C:6:THR:HA	1:C:39:ARG:O	2.00	0.62
1:B:49:LYS:HG3	1:B:148:LEU:HD13	1.82	0.61
1:C:56:LEU:HD22	1:C:159:PRO:HD3	1.81	0.61
1:B:6:THR:HA	1:B:39:ARG:O	1.99	0.61
1:C:104:VAL:HB	1:C:107:ALA:HB3	1.83	0.61
1:A:49:LYS:HG3	1:A:148:LEU:HD13	1.82	0.61
1:B:56:LEU:HD22	1:B:159:PRO:HD3	1.82	0.60
1:C:104:VAL:HB	1:C:107:ALA:CB	2.32	0.60
1:D:104:VAL:HB	1:D:107:ALA:HB3	1.82	0.60
1:C:117:ILE:HD13	1:C:117:ILE:H	1.65	0.60
1:D:104:VAL:HB	1:D:107:ALA:CB	2.32	0.60
1:C:65:ILE:CD1	1:C:115:LEU:HD21	2.30	0.60
1:C:78:ILE:HD13	1:C:165:ILE:HG12	1.84	0.60
1:B:76:ILE:CD1	1:B:129:ILE:CG2	2.78	0.59
1:D:45:PHE:CE2	1:D:86:ILE:HD12	2.37	0.59
1:C:90:ILE:H	1:C:90:ILE:CD1	2.14	0.59
1:A:64:ALA:C	1:A:65:ILE:HD13	2.22	0.59
1:C:49:LYS:HG3	1:C:148:LEU:HD13	1.85	0.59
1:B:115:LEU:HD23	1:B:161:MET:HB3	1.85	0.58
1:D:146:MET:HG3	1:D:164:PHE:HB2	1.85	0.58
1:D:80:ARG:HB3	1:D:81:ILE:HD13	1.84	0.58
1:A:18:PRO:HB3	1:A:170:ILE:HD12	1.86	0.58
1:B:170:ILE:O	1:B:174:ILE:CD1	2.52	0.57
1:A:115:LEU:HD23	1:A:161:MET:HB3	1.86	0.57
1:D:16:ILE:HD12	2:D:211:HOH:O	2.04	0.57
1:D:45:PHE:HE2	1:D:86:ILE:HD12	1.68	0.57
1:C:16:ILE:HD13	1:C:17:ASN:N	2.20	0.56
1:A:65:ILE:CD1	1:A:115:LEU:HD21	2.33	0.56
1:D:115:LEU:HD23	1:D:161:MET:HB3	1.86	0.56
1:A:62:ASP:HB3	1:A:207:LEU:HD11	1.87	0.56
1:C:174:ILE:HG13	1:C:186:SER:HB3	1.87	0.56
1:D:78:ILE:CD1	1:D:165:ILE:HG12	2.34	0.56
1:D:174:ILE:HG13	1:D:186:SER:HB3	1.88	0.56
1:D:22:ILE:HG21	1:D:196:VAL:HG11	1.88	0.56
1:D:11:PHE:HA	1:D:93:ASN:HD21	1.70	0.56
1:D:129:ILE:HD12	1:D:191:MET:HE2	1.85	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:11:PHE:HA	1:C:93:ASN:HD21	1.70	0.55
1:C:78:ILE:CD1	1:C:165:ILE:HG12	2.35	0.55
1:A:76:ILE:HD12	1:A:129:ILE:HG21	1.88	0.55
1:C:115:LEU:HD23	1:C:161:MET:HB3	1.88	0.55
1:B:76:ILE:CD1	1:B:129:ILE:HG21	2.37	0.55
1:B:130:PRO:HA	1:D:175:ILE:HG22	1.88	0.55
1:A:11:PHE:HA	1:A:93:ASN:HD21	1.71	0.55
1:B:11:PHE:HA	1:B:93:ASN:HD21	1.70	0.55
1:D:129:ILE:N	1:D:129:ILE:HD13	2.23	0.54
1:B:22:ILE:HG21	1:B:196:VAL:HG11	1.90	0.54
1:A:3:VAL:HG11	1:A:200:ILE:HD12	1.90	0.54
1:B:194:GLU:O	1:B:198:VAL:HG23	2.08	0.53
1:B:174:ILE:HD12	1:B:174:ILE:H	1.73	0.53
1:B:174:ILE:N	1:B:174:ILE:CD1	2.71	0.53
1:B:171:PRO:CA	1:B:174:ILE:HD13	2.33	0.53
1:C:182:GLN:HA	1:C:182:GLN:NE2	2.24	0.53
1:D:86:ILE:HG21	1:D:98:ILE:HB	1.91	0.53
1:A:65:ILE:CD1	1:A:115:LEU:CD1	2.78	0.52
1:B:81:ILE:CD1	1:C:83:VAL:HG11	2.39	0.52
1:D:127:ARG:O	1:D:129:ILE:HD13	2.10	0.52
1:A:76:ILE:CD1	1:A:129:ILE:CG2	2.84	0.52
1:B:90:ILE:H	1:B:90:ILE:CD1	2.17	0.52
1:C:146:MET:HG3	1:C:164:PHE:HB2	1.91	0.52
1:A:76:ILE:CD1	1:A:129:ILE:HG21	2.41	0.51
1:A:22:ILE:HG21	1:A:196:VAL:HG11	1.92	0.51
1:A:194:GLU:O	1:A:198:VAL:HG23	2.10	0.51
1:C:86:ILE:HG13	1:C:103:ILE:CD1	2.39	0.51
1:B:76:ILE:CD1	1:B:195:ALA:CB	2.88	0.51
1:D:152:HIS:HD2	1:D:158:TYR:O	1.93	0.51
1:A:182:GLN:HA	1:A:182:GLN:NE2	2.25	0.51
1:C:39:ARG:NH2	1:C:59:ILE:HD11	2.25	0.51
1:D:33:ASP:HA	2:D:216:HOH:O	2.10	0.51
1:D:182:GLN:HA	1:D:182:GLN:NE2	2.26	0.51
1:A:152:HIS:HD2	1:A:158:TYR:O	1.92	0.51
1:B:16:ILE:CD1	1:B:18:PRO:HD3	2.40	0.51
1:A:90:ILE:H	1:A:90:ILE:CD1	2.16	0.51
1:C:22:ILE:HG21	1:C:196:VAL:HG11	1.93	0.50
1:B:146:MET:HG3	1:B:164:PHE:HB2	1.93	0.50
1:C:152:HIS:HD2	1:C:158:TYR:O	1.94	0.50
1:D:194:GLU:O	1:D:198:VAL:HG23	2.11	0.50
1:B:37:PHE:CE2	1:B:59:ILE:HD12	2.47	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:ARG:NH2	1:A:59:ILE:HD11	2.27	0.50
1:B:86:ILE:HD11	1:B:144:TYR:CD1	2.48	0.49
1:A:66:HIS:HD2	2:A:227:HOH:O	1.94	0.49
1:A:174:ILE:HG13	1:A:186:SER:HB3	1.94	0.49
1:B:76:ILE:CD1	1:B:195:ALA:HB2	2.42	0.49
1:B:152:HIS:HD2	1:B:158:TYR:O	1.95	0.49
1:B:174:ILE:CD1	1:B:174:ILE:H	2.25	0.49
1:D:70:ALA:O	1:D:73:ARG:HB2	2.12	0.49
1:D:62:ASP:HB3	1:D:207:LEU:HD11	1.94	0.49
1:A:146:MET:HG3	1:A:164:PHE:HB2	1.95	0.49
1:B:182:GLN:HA	1:B:182:GLN:NE2	2.27	0.49
1:C:16:ILE:CD1	1:C:18:PRO:HD3	2.43	0.49
1:D:37:PHE:CE2	1:D:59:ILE:HD12	2.48	0.49
1:A:86:ILE:HG21	1:A:98:ILE:HB	1.94	0.49
1:B:45:PHE:HE2	1:B:86:ILE:HD12	1.78	0.49
1:B:174:ILE:HG12	1:B:186:SER:HB3	1.95	0.49
1:D:80:ARG:CB	1:D:81:ILE:HD13	2.43	0.49
1:D:117:ILE:HD13	1:D:117:ILE:N	2.28	0.48
1:A:66:HIS:CD2	2:A:227:HOH:O	2.65	0.48
1:A:76:ILE:CD1	1:A:195:ALA:CB	2.89	0.48
1:B:86:ILE:HG21	1:B:98:ILE:HB	1.94	0.48
1:C:117:ILE:HD13	1:C:117:ILE:N	2.27	0.48
1:D:104:VAL:HG23	1:D:147:TYR:HE2	1.79	0.48
1:B:62:ASP:HB3	1:B:207:LEU:HD11	1.95	0.48
1:B:175:ILE:HD13	1:D:125:HIS:O	2.13	0.48
1:D:86:ILE:HD13	1:D:103:ILE:HD11	1.95	0.48
1:A:70:ALA:O	1:A:73:ARG:HB2	2.13	0.48
1:B:175:ILE:HG22	1:D:130:PRO:HA	1.95	0.48
1:C:16:ILE:HD13	1:C:16:ILE:C	2.34	0.47
1:A:90:ILE:HD13	1:A:90:ILE:N	2.20	0.47
1:C:62:ASP:HB3	1:C:207:LEU:HD11	1.96	0.47
1:C:90:ILE:HD13	1:C:90:ILE:N	2.19	0.47
1:D:86:ILE:HD11	1:D:144:TYR:CD1	2.49	0.47
1:A:156:LYS:HD3	2:A:221:HOH:O	2.15	0.47
1:B:83:VAL:HG11	1:C:81:ILE:CD1	2.44	0.47
1:C:65:ILE:CD1	1:C:115:LEU:CD1	2.84	0.47
1:D:81:ILE:HD13	1:D:81:ILE:N	2.29	0.47
1:A:174:ILE:HG23	1:A:184:PRO:HD2	1.97	0.47
1:B:80:ARG:CD	1:B:135:ASN:ND2	2.78	0.46
1:D:23:ALA:O	1:D:38:GLY:HA3	2.15	0.46
1:A:104:VAL:HG23	1:A:147:TYR:HE2	1.80	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:PRO:CB	1:A:119:LYS:HD2	2.38	0.46
1:B:45:PHE:CE2	1:B:86:ILE:HD12	2.50	0.46
1:B:16:ILE:HD13	1:B:17:ASN:N	2.30	0.46
1:B:164:PHE:C	1:B:165:ILE:HD12	2.35	0.46
1:D:80:ARG:CD	1:D:135:ASN:ND2	2.78	0.46
1:B:16:ILE:HD13	1:B:18:PRO:HD3	1.97	0.46
1:C:65:ILE:CD1	1:C:115:LEU:CD2	2.93	0.45
1:A:130:PRO:HA	1:C:175:ILE:HG22	1.99	0.45
1:B:16:ILE:HD13	1:B:16:ILE:C	2.37	0.45
1:D:76:ILE:O	1:D:131:ALA:HA	2.16	0.45
1:D:170:ILE:HA	1:D:187:MET:O	2.16	0.45
1:A:175:ILE:HG22	1:C:130:PRO:HA	1.98	0.45
1:B:80:ARG:HD2	1:B:135:ASN:ND2	2.31	0.45
1:B:174:ILE:HG23	1:B:184:PRO:HD2	1.98	0.45
1:A:76:ILE:CD1	1:A:195:ALA:HB2	2.46	0.45
1:B:90:ILE:HD13	1:B:90:ILE:N	2.20	0.45
1:C:117:ILE:H	1:C:117:ILE:CD1	2.30	0.45
1:A:164:PHE:C	1:A:165:ILE:HD12	2.37	0.45
1:A:109:THR:HG23	1:D:112:PHE:CE1	2.51	0.45
1:B:124:LEU:HD23	1:B:124:LEU:HA	1.79	0.45
1:C:174:ILE:HG23	1:C:184:PRO:HD2	1.99	0.45
1:C:70:ALA:O	1:C:73:ARG:HB2	2.17	0.45
1:A:65:ILE:CD1	1:A:115:LEU:CD2	2.93	0.44
1:C:194:GLU:O	1:C:198:VAL:HG23	2.17	0.44
1:B:130:PRO:CA	1:D:175:ILE:HG22	2.47	0.44
1:C:80:ARG:CD	1:C:135:ASN:ND2	2.80	0.44
1:D:80:ARG:HD2	1:D:135:ASN:ND2	2.32	0.44
1:A:5:VAL:HA	1:A:65:ILE:O	2.17	0.44
1:D:80:ARG:C	1:D:81:ILE:HD13	2.38	0.44
1:A:70:ALA:HB2	1:A:166:HIS:CD2	2.52	0.44
1:B:174:ILE:CG1	1:B:186:SER:HB3	2.47	0.44
1:D:16:ILE:HA	2:D:211:HOH:O	2.18	0.44
1:A:124:LEU:HD23	1:A:124:LEU:HA	1.78	0.44
1:C:104:VAL:HG23	1:C:147:TYR:HE2	1.82	0.44
1:C:86:ILE:HG13	1:C:103:ILE:HD13	1.98	0.44
1:C:175:ILE:C	1:C:175:ILE:HD12	2.37	0.44
1:C:174:ILE:CG1	1:C:186:SER:HB3	2.48	0.44
1:A:76:ILE:HD13	1:A:195:ALA:HB2	1.99	0.44
1:D:117:ILE:H	1:D:117:ILE:CD1	2.29	0.44
1:A:71:PRO:HB3	1:A:169:TYR:CZ	2.53	0.43
1:A:112:PHE:CE1	1:D:109:THR:HG23	2.52	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:ILE:CD1	1:D:83:VAL:HG11	2.47	0.43
1:D:174:ILE:HG23	1:D:184:PRO:HD2	2.00	0.43
1:D:175:ILE:HD12	1:D:175:ILE:C	2.38	0.43
1:C:3:VAL:HG22	1:C:63:ILE:HB	2.00	0.43
1:B:170:ILE:HA	1:B:187:MET:O	2.19	0.43
1:D:5:VAL:HA	1:D:65:ILE:O	2.19	0.43
1:D:116:PRO:CB	1:D:119:LYS:HD2	2.38	0.43
1:C:86:ILE:HG21	1:C:98:ILE:HB	2.00	0.43
1:A:80:ARG:CD	1:A:135:ASN:ND2	2.81	0.43
1:B:29:ILE:C	1:B:29:ILE:HD12	2.39	0.42
1:B:104:VAL:HG23	1:B:147:TYR:HE2	1.83	0.42
1:C:26:LEU:O	1:C:29:ILE:HG12	2.18	0.42
1:C:90:ILE:CD1	1:C:90:ILE:N	2.81	0.42
1:D:26:LEU:O	1:D:29:ILE:HG13	2.19	0.42
1:D:67:VAL:HA	1:D:165:ILE:O	2.19	0.42
1:A:175:ILE:HD13	1:C:125:HIS:O	2.19	0.42
1:C:80:ARG:HD3	1:C:135:ASN:ND2	2.35	0.42
1:D:29:ILE:HD12	1:D:29:ILE:C	2.39	0.42
1:D:70:ALA:HB2	1:D:166:HIS:CD2	2.54	0.42
1:D:171:PRO:HD3	1:D:187:MET:O	2.19	0.42
1:A:68:GLY:O	1:A:166:HIS:HA	2.20	0.42
1:A:170:ILE:HA	1:A:187:MET:O	2.20	0.42
1:A:174:ILE:CG1	1:A:186:SER:HB3	2.49	0.42
1:B:70:ALA:HB2	1:B:166:HIS:CD2	2.55	0.42
1:A:29:ILE:HD12	1:A:29:ILE:C	2.40	0.42
1:D:124:LEU:HD23	1:D:124:LEU:HA	1.81	0.41
1:C:8:PHE:O	1:C:20:GLU:HB2	2.20	0.41
1:D:3:VAL:HG22	1:D:63:ILE:HB	2.03	0.41
1:A:80:ARG:HD3	1:A:135:ASN:ND2	2.35	0.41
1:C:170:ILE:HD12	1:C:172:GLU:HG2	2.02	0.41
1:D:103:ILE:HD13	1:D:103:ILE:N	2.35	0.41
1:D:129:ILE:N	1:D:129:ILE:CD1	2.84	0.41
1:D:82:ALA:HB1	1:D:147:TYR:HB2	2.02	0.41
1:A:146:MET:O	1:A:146:MET:HG2	2.21	0.41
1:A:30:LYS:HB3	1:A:30:LYS:HE2	1.90	0.41
1:D:8:PHE:O	1:D:20:GLU:HB2	2.21	0.41
1:B:171:PRO:HD3	1:B:187:MET:O	2.21	0.41
1:A:76:ILE:O	1:A:131:ALA:HA	2.21	0.40
1:C:76:ILE:O	1:C:131:ALA:HA	2.21	0.40
1:C:124:LEU:HA	1:C:124:LEU:HD23	1.81	0.40
1:B:119:LYS:HE3	1:B:119:LYS:N	2.37	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:85:ALA:C	1:D:103:ILE:HD11	2.41	0.40
1:B:37:PHE:HD2	1:B:39:ARG:HH22	1.69	0.40
1:C:5:VAL:HA	1:C:65:ILE:O	2.22	0.40
1:D:43:VAL:CG1	1:D:141:LEU:HD13	2.52	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	206/208 (99%)	189 (92%)	16 (8%)	1 (0%)	29 54
1	B	206/208 (99%)	190 (92%)	16 (8%)	0	100 100
1	C	206/208 (99%)	188 (91%)	18 (9%)	0	100 100
1	D	206/208 (99%)	189 (92%)	17 (8%)	0	100 100
All	All	824/832 (99%)	756 (92%)	67 (8%)	1 (0%)	51 78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	12	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	174/174 (100%)	157 (90%)	17 (10%)	8	18
1	B	174/174 (100%)	157 (90%)	17 (10%)	8	18
1	C	174/174 (100%)	156 (90%)	18 (10%)	7	16
1	D	174/174 (100%)	156 (90%)	18 (10%)	7	16
All	All	696/696 (100%)	626 (90%)	70 (10%)	7	17

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	35	GLN
1	A	52	LEU
1	A	69	LEU
1	A	73	ARG
1	A	81	ILE
1	A	89	ARG
1	A	90	ILE
1	A	119	LYS
1	A	141	LEU
1	A	160	LYS
1	A	170	ILE
1	A	172	GLU
1	A	180	LYS
1	A	182	GLN
1	A	196	VAL
1	A	206	GLU
1	B	4	LEU
1	B	16	ILE
1	B	35	GLN
1	B	52	LEU
1	B	69	LEU
1	B	73	ARG
1	B	81	ILE
1	B	89	ARG
1	B	90	ILE
1	B	119	LYS
1	B	141	LEU
1	B	160	LYS
1	B	172	GLU
1	B	180	LYS
1	B	182	GLN
1	B	196	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	206	GLU
1	C	4	LEU
1	C	16	ILE
1	C	35	GLN
1	C	52	LEU
1	C	69	LEU
1	C	73	ARG
1	C	81	ILE
1	C	89	ARG
1	C	90	ILE
1	C	117	ILE
1	C	119	LYS
1	C	141	LEU
1	C	160	LYS
1	C	172	GLU
1	C	180	LYS
1	C	182	GLN
1	C	196	VAL
1	C	206	GLU
1	D	4	LEU
1	D	35	GLN
1	D	52	LEU
1	D	69	LEU
1	D	73	ARG
1	D	81	ILE
1	D	89	ARG
1	D	103	ILE
1	D	117	ILE
1	D	119	LYS
1	D	129	ILE
1	D	141	LEU
1	D	160	LYS
1	D	172	GLU
1	D	180	LYS
1	D	182	GLN
1	D	196	VAL
1	D	206	GLU

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	93	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	135	ASN
1	A	152	HIS
1	B	93	ASN
1	B	135	ASN
1	B	152	HIS
1	C	93	ASN
1	C	135	ASN
1	C	152	HIS
1	D	93	ASN
1	D	135	ASN
1	D	152	HIS

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)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.