



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 01:08 AM UTC

PDB ID : 4FD3 / pdb_00004fd3
Title : Crystal structure of apo-formed ymtOAR1
Authors : Zhang, Y.; Gao, Y.; Ning, F.; Niu, L.; Teng, M.
Deposited on : 2012-05-26
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 ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Xtrriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

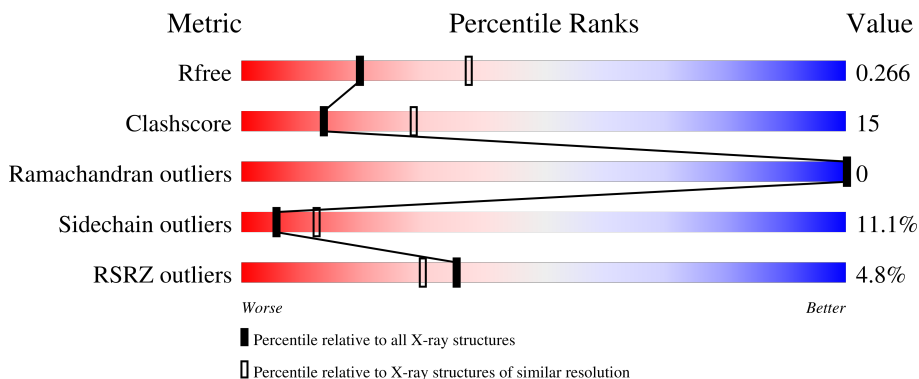
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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.



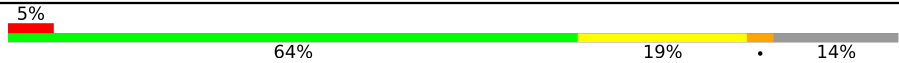
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	286	 5% 63% 21% 5% 11%
1	B	286	 3% 61% 22% 5% 12%
1	C	286	 2% 63% 21% 5% 13%
1	D	286	 6% 66% 21% 6% 6%
1	E	286	 5% 59% 25% 5% 13%

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Mol	Chain	Length	Quality of chain
1	F	286	 <p>A horizontal bar chart representing the quality of the chain. The bar is divided into four segments: a small red segment (5%), a large green segment (64%), a yellow segment (19%), and a small grey segment (14%).</p>

2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 11328 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 3-oxoacyl-[acyl-carrier-protein] reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	254	Total 1865	C 1189	N 320	O 342	S 14	0	0	0
1	B	253	Total 1903	C 1212	N 323	O 353	S 15	0	0	0
1	C	250	Total 1878	C 1198	N 319	O 346	S 15	0	0	0
1	D	268	Total 1961	C 1248	N 338	O 360	S 15	0	0	0
1	E	250	Total 1837	C 1174	N 311	O 337	S 15	0	0	0
1	F	246	Total 1820	C 1161	N 312	O 333	S 14	0	0	0

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	279	LEU	-	expression tag	UNP P35731
A	280	GLU	-	expression tag	UNP P35731
A	281	HIS	-	expression tag	UNP P35731
A	282	HIS	-	expression tag	UNP P35731
A	283	HIS	-	expression tag	UNP P35731
A	284	HIS	-	expression tag	UNP P35731
A	285	HIS	-	expression tag	UNP P35731
A	286	HIS	-	expression tag	UNP P35731
B	279	LEU	-	expression tag	UNP P35731
B	280	GLU	-	expression tag	UNP P35731
B	281	HIS	-	expression tag	UNP P35731
B	282	HIS	-	expression tag	UNP P35731
B	283	HIS	-	expression tag	UNP P35731
B	284	HIS	-	expression tag	UNP P35731
B	285	HIS	-	expression tag	UNP P35731
B	286	HIS	-	expression tag	UNP P35731
C	279	LEU	-	expression tag	UNP P35731

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Chain	Residue	Modelled	Actual	Comment	Reference
C	280	GLU	-	expression tag	UNP P35731
C	281	HIS	-	expression tag	UNP P35731
C	282	HIS	-	expression tag	UNP P35731
C	283	HIS	-	expression tag	UNP P35731
C	284	HIS	-	expression tag	UNP P35731
C	285	HIS	-	expression tag	UNP P35731
C	286	HIS	-	expression tag	UNP P35731
D	279	LEU	-	expression tag	UNP P35731
D	280	GLU	-	expression tag	UNP P35731
D	281	HIS	-	expression tag	UNP P35731
D	282	HIS	-	expression tag	UNP P35731
D	283	HIS	-	expression tag	UNP P35731
D	284	HIS	-	expression tag	UNP P35731
D	285	HIS	-	expression tag	UNP P35731
D	286	HIS	-	expression tag	UNP P35731
E	279	LEU	-	expression tag	UNP P35731
E	280	GLU	-	expression tag	UNP P35731
E	281	HIS	-	expression tag	UNP P35731
E	282	HIS	-	expression tag	UNP P35731
E	283	HIS	-	expression tag	UNP P35731
E	284	HIS	-	expression tag	UNP P35731
E	285	HIS	-	expression tag	UNP P35731
E	286	HIS	-	expression tag	UNP P35731
F	279	LEU	-	expression tag	UNP P35731
F	280	GLU	-	expression tag	UNP P35731
F	281	HIS	-	expression tag	UNP P35731
F	282	HIS	-	expression tag	UNP P35731
F	283	HIS	-	expression tag	UNP P35731
F	284	HIS	-	expression tag	UNP P35731
F	285	HIS	-	expression tag	UNP P35731
F	286	HIS	-	expression tag	UNP P35731

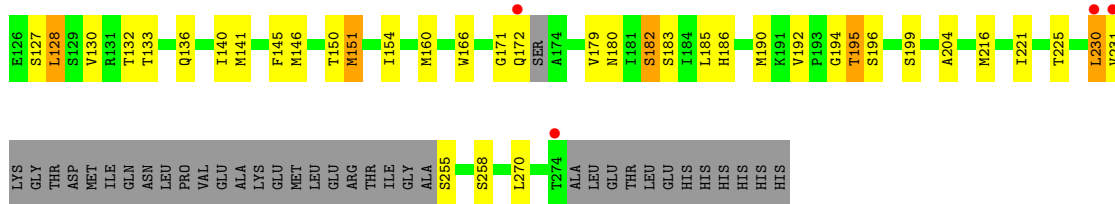
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	6	Total O 6 6	0	0
2	B	15	Total O 15 15	0	0
2	C	16	Total O 16 16	0	0
2	D	9	Total O 9 9	0	0

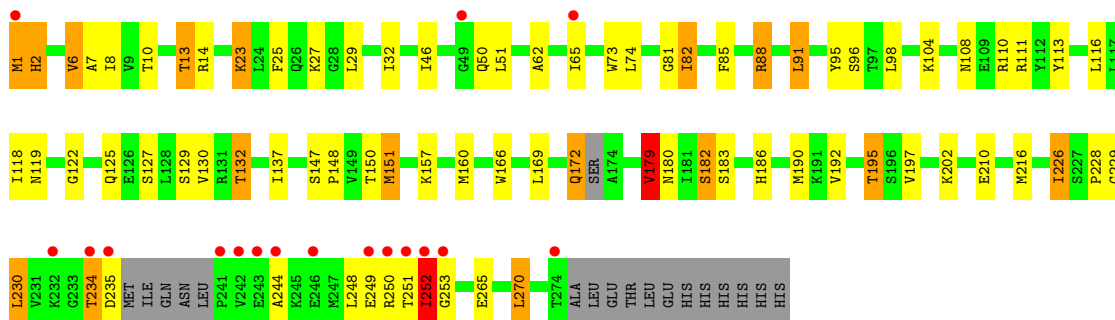
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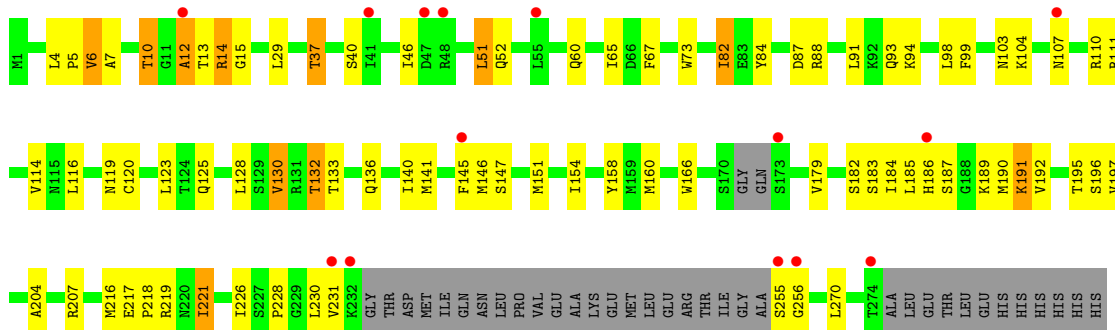
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	E	9	Total O 9 9	0	0
2	F	9	Total O 9 9	0	0



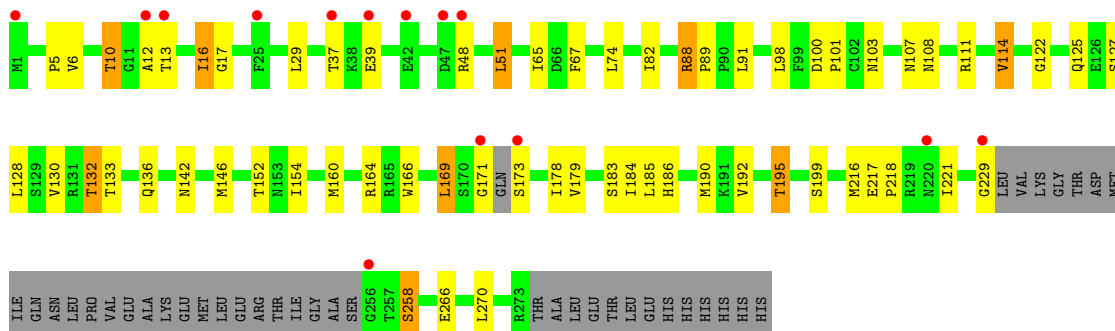
• Molecule 1: 3-oxoacyl-[acyl-carrier-protein] reductase



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4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	170.83Å 170.83Å 316.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.20 – 2.60 48.20 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.4 (48.20-2.60) 98.4 (48.20-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.73 (at 2.61Å)	Xtrriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.229 , 0.274 0.225 , 0.266	Depositor DCC
R_{free} test set	3577 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	50.2	Xtrriage
Anisotropy	0.097	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11328	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.20% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality 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	0.78	2/1905 (0.1%)	1.00	8/2598 (0.3%)
1	B	0.86	5/1944 (0.3%)	0.97	3/2644 (0.1%)
1	C	0.80	3/1918 (0.2%)	0.99	3/2609 (0.1%)
1	D	0.76	2/2002 (0.1%)	0.99	7/2727 (0.3%)
1	E	0.83	2/1876 (0.1%)	1.07	8/2559 (0.3%)
1	F	0.75	2/1859 (0.1%)	0.96	3/2532 (0.1%)
All	All	0.80	16/11504 (0.1%)	1.00	32/15669 (0.2%)

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	192	VAL	CA-C	-6.79	1.46	1.52
1	B	122	GLY	C-O	-6.64	1.17	1.24
1	E	192	VAL	CA-CB	-6.41	1.48	1.54
1	F	114	VAL	CA-CB	6.03	1.61	1.54
1	B	179	VAL	C-O	-5.86	1.18	1.24
1	D	179	VAL	C-O	-5.57	1.18	1.24
1	F	16	ILE	CA-CB	-5.57	1.47	1.54
1	B	180	ASN	C-O	-5.46	1.17	1.24
1	D	229	GLY	C-O	-5.28	1.18	1.23
1	B	119	ASN	C-O	-5.23	1.18	1.24
1	C	122	GLY	C-O	-5.16	1.19	1.24
1	A	181	ILE	CA-CB	-5.12	1.48	1.54
1	C	182	SER	C-O	-5.09	1.18	1.23
1	C	107	ASN	C-O	-5.02	1.18	1.23
1	B	182	SER	C-O	-5.01	1.18	1.23
1	A	180	ASN	C-O	-5.00	1.18	1.24

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	14	ARG	N-CA-C	11.06	127.10	113.17
1	A	232	LYS	N-CA-C	-10.36	99.75	112.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	233	GLY	N-CA-C	10.29	125.08	112.73
1	F	13	THR	N-CA-C	8.59	120.42	111.14
1	C	106	SER	N-CA-C	7.72	120.72	110.53
1	E	189	LYS	CA-C-N	-7.17	110.91	120.95
1	E	189	LYS	C-N-CA	-7.17	110.91	120.95
1	A	122	GLY	N-CA-C	7.08	121.05	112.49
1	A	234	THR	N-CA-C	6.83	119.55	109.24
1	B	122	GLY	N-CA-C	6.81	126.56	115.46
1	B	232	LYS	N-CA-C	6.75	118.63	111.28
1	E	13	THR	N-CA-C	6.71	122.35	114.04
1	E	184	ILE	CA-C-N	6.59	129.77	120.28
1	E	184	ILE	C-N-CA	6.59	129.77	120.28
1	D	234	THR	CB-CA-C	-6.50	109.09	116.63
1	E	12	ALA	N-CA-C	6.41	118.97	108.52
1	F	122	GLY	N-CA-C	6.25	123.51	115.32
1	E	191	LYS	N-CA-C	-6.21	98.27	108.76
1	B	118	ILE	CB-CA-C	-6.11	102.10	110.77
1	D	249	GLU	N-CA-C	-6.03	104.78	111.71
1	D	252	ILE	CA-C-N	5.96	126.20	121.61
1	D	252	ILE	C-N-CA	5.96	126.20	121.61
1	C	120	CYS	N-CA-C	5.82	118.57	111.82
1	A	229	GLY	N-CA-C	-5.75	103.72	112.51
1	A	197	VAL	N-CA-C	5.48	116.26	110.72
1	D	244	ALA	N-CA-C	-5.42	105.38	111.28
1	A	88	ARG	N-CA-C	5.38	112.86	108.07
1	C	122	GLY	N-CA-C	5.31	124.12	115.46
1	A	221	ILE	N-CA-C	5.14	116.15	108.54
1	D	110	ARG	N-CA-C	5.13	117.76	108.69
1	D	252	ILE	N-CA-C	5.10	115.00	107.75
1	F	114	VAL	N-CA-CB	5.01	116.18	110.72

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1865	0	1778	64	0
1	B	1903	0	1847	53	0
1	C	1878	0	1815	49	0
1	D	1961	0	1859	66	0
1	E	1837	0	1751	53	0
1	F	1820	0	1744	50	0
2	A	6	0	0	0	0
2	B	15	0	0	0	0
2	C	16	0	0	0	0
2	D	9	0	0	0	0
2	E	9	0	0	2	0
2	F	9	0	0	0	0
All	All	11328	0	10794	321	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 15.

All (321) 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:A:230:LEU:H	1:A:230:LEU:CD1	1.56	1.18
1:C:171:GLY:HA3	1:C:172:GLN:CB	1.71	1.18
1:A:230:LEU:HD12	1:A:230:LEU:N	1.52	1.08
1:D:1:MET:N	1:D:2:HIS:HA	1.58	1.08
1:D:88:ARG:HG2	1:D:88:ARG:HH11	0.95	1.06
1:D:1:MET:H2	1:D:2:HIS:HA	0.90	1.05
1:A:231:VAL:HG23	1:A:233:GLY:H	1.16	1.05
1:D:1:MET:N	1:D:2:HIS:CA	2.22	1.03
1:D:1:MET:O	1:D:1:MET:HG2	1.61	0.99
1:D:88:ARG:HG2	1:D:88:ARG:NH1	1.73	0.98
1:E:185:LEU:HD23	1:E:195:THR:HG21	1.45	0.96
1:A:230:LEU:H	1:A:230:LEU:HD12	0.79	0.95
1:F:12:ALA:CB	1:F:17:GLY:HA3	1.97	0.94
1:D:1:MET:H2	1:D:2:HIS:CA	1.80	0.93
1:E:12:ALA:HB2	1:E:120:CYS:SG	2.08	0.93
1:D:1:MET:H3	1:D:2:HIS:C	1.77	0.92
1:F:133:THR:H	1:F:136:GLN:HE21	1.17	0.92
1:D:230:LEU:N	1:D:230:LEU:HD12	1.82	0.92
1:E:98:LEU:HD13	1:E:151:MET:HE1	1.52	0.91
1:C:171:GLY:CA	1:C:172:GLN:CB	2.48	0.91
1:B:10:THR:HG21	1:B:119:ASN:OD1	1.70	0.91
1:D:192:VAL:O	1:D:195:THR:HB	1.69	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:230:LEU:HD12	1:D:230:LEU:H	1.36	0.90
1:C:160:MET:HE3	1:C:216:MET:HE1	1.51	0.88
1:D:160:MET:HE3	1:D:216:MET:HE1	1.57	0.87
1:F:12:ALA:HB1	1:F:17:GLY:HA3	1.57	0.86
1:F:74:LEU:HD21	1:F:154:ILE:HD11	1.57	0.86
1:A:231:VAL:HG23	1:A:233:GLY:N	1.91	0.84
1:C:230:LEU:O	1:C:231:VAL:HG23	1.79	0.82
1:A:133:THR:H	1:A:136:GLN:NE2	1.77	0.82
1:D:13:THR:CG2	1:D:14:ARG:HE	1.93	0.80
1:C:111:ARG:HD3	1:C:166:TRP:CZ2	2.17	0.80
1:A:118:ILE:HD13	1:A:179:VAL:HG13	1.63	0.80
1:A:185:LEU:HD23	1:A:195:THR:HG21	1.63	0.80
1:D:230:LEU:H	1:D:230:LEU:CD1	1.95	0.79
1:B:133:THR:H	1:B:136:GLN:NE2	1.80	0.79
1:D:23:LYS:HE3	1:D:265:GLU:OE2	1.82	0.79
1:C:133:THR:H	1:C:136:GLN:HE21	1.28	0.78
1:C:185:LEU:HD23	1:C:195:THR:HG21	1.63	0.78
1:B:192:VAL:O	1:B:195:THR:HB	1.83	0.78
1:E:14:ARG:N	1:E:15:GLY:HA2	2.00	0.77
1:A:46:ILE:H	1:A:60:GLN:HE22	1.32	0.77
1:B:151:MET:HA	1:B:151:MET:HE3	1.66	0.77
1:B:133:THR:H	1:B:136:GLN:HE21	1.32	0.77
1:F:266:GLU:O	1:F:270:LEU:HD12	1.84	0.76
1:A:111:ARG:HD3	1:A:166:TRP:CZ2	2.21	0.76
1:F:133:THR:H	1:F:136:GLN:NE2	1.83	0.75
1:D:98:LEU:HD13	1:D:151:MET:HE1	1.67	0.74
1:E:46:ILE:H	1:E:60:GLN:HE22	1.35	0.74
1:E:216:MET:HE2	1:E:221:ILE:HG13	1.69	0.73
1:D:65:ILE:HD11	1:D:73:TRP:CE3	2.24	0.73
1:C:133:THR:H	1:C:136:GLN:NE2	1.87	0.72
1:B:46:ILE:H	1:B:60:GLN:HE22	1.36	0.72
1:F:48:ARG:HA	1:F:51:LEU:HB2	1.71	0.71
1:A:107:ASN:C	1:A:107:ASN:OD1	2.33	0.71
1:D:234:THR:O	1:D:235:ASP:C	2.30	0.71
1:F:216:MET:HE3	1:F:221:ILE:HB	1.73	0.71
1:C:127:SER:OG	1:C:132:THR:HG22	1.90	0.71
1:A:195:THR:HG22	1:A:195:THR:O	1.90	0.70
1:F:160:MET:HE3	1:F:216:MET:HE1	1.74	0.69
1:F:183:SER:O	1:F:186:HIS:HD2	1.75	0.69
1:E:104:LYS:HE3	2:E:305:HOH:O	1.90	0.69
1:E:195:THR:O	1:E:195:THR:HG22	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:127:SER:OG	1:F:132:THR:HG22	1.92	0.69
1:A:12:ALA:HB1	1:A:17:GLY:HA3	1.75	0.69
1:C:36:SER:HA	1:C:65:ILE:O	1.92	0.69
1:F:133:THR:N	1:F:136:GLN:HE21	1.90	0.68
1:F:192:VAL:O	1:F:195:THR:HB	1.93	0.68
1:E:111:ARG:HD3	1:E:166:TRP:CZ2	2.29	0.68
1:F:229:GLY:H	1:F:258:SER:HB2	1.59	0.68
1:D:190:MET:HE1	1:D:248:LEU:CB	2.23	0.67
1:B:98:LEU:HD23	1:B:151:MET:HE1	1.75	0.67
1:B:10:THR:CG2	1:B:119:ASN:OD1	2.43	0.67
1:A:128:LEU:O	1:A:132:THR:HG23	1.95	0.66
1:B:157:LYS:HD2	1:E:130:VAL:HG22	1.78	0.66
1:E:104:LYS:CE	2:E:305:HOH:O	2.42	0.65
1:C:48:ARG:O	1:C:52:GLN:HG3	1.96	0.64
1:D:169:LEU:O	1:D:172:GLN:HB3	1.98	0.63
1:E:107:ASN:OD1	1:E:107:ASN:C	2.41	0.63
1:F:88:ARG:HB2	1:F:89:PRO:CD	2.28	0.63
1:B:111:ARG:HD3	1:B:166:TRP:CE2	2.34	0.63
1:C:183:SER:O	1:C:186:HIS:HD2	1.82	0.63
1:D:13:THR:HG22	1:D:14:ARG:HE	1.62	0.63
1:F:12:ALA:HB1	1:F:17:GLY:CA	2.28	0.63
1:D:88:ARG:HH11	1:D:88:ARG:CG	1.89	0.62
1:E:226:ILE:HG22	1:E:228:PRO:HD3	1.80	0.62
1:A:133:THR:H	1:A:136:GLN:HE21	1.43	0.62
1:A:231:VAL:C	1:A:233:GLY:N	2.51	0.62
1:A:12:ALA:CB	1:A:17:GLY:HA3	2.29	0.62
1:D:1:MET:N	1:D:2:HIS:C	2.47	0.62
1:B:111:ARG:HD3	1:B:166:TRP:CZ2	2.34	0.62
1:A:18:LYS:HG3	1:A:43:ARG:O	1.99	0.62
1:C:128:LEU:O	1:C:132:THR:HG23	1.99	0.61
1:B:160:MET:HE3	1:B:216:MET:HE1	1.82	0.61
1:C:185:LEU:HD23	1:C:195:THR:CG2	2.30	0.61
1:D:1:MET:H3	1:D:2:HIS:CA	2.04	0.61
1:F:65:ILE:HD11	1:F:98:LEU:HD21	1.83	0.61
1:E:141:MET:HB3	1:E:146:MET:HE1	1.81	0.61
1:B:151:MET:HA	1:B:151:MET:CE	2.28	0.60
1:C:108:ASN:HD22	1:C:108:ASN:H	1.48	0.60
1:D:111:ARG:HD3	1:D:166:TRP:CZ2	2.37	0.60
1:E:65:ILE:HD11	1:E:73:TRP:CE3	2.35	0.59
1:A:160:MET:HE3	1:A:216:MET:HE1	1.82	0.59
1:E:98:LEU:CD1	1:E:151:MET:HE1	2.28	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:VAL:HG22	1:A:30:SER:HB2	1.85	0.59
1:B:46:ILE:HD11	1:B:51:LEU:HD13	1.84	0.59
1:F:111:ARG:HD3	1:F:166:TRP:CZ2	2.38	0.59
1:C:8:ILE:HG12	1:C:32:ILE:HB	1.85	0.58
1:A:182:SER:OG	1:A:183:SER:N	2.36	0.58
1:B:26:GLN:HE21	1:B:27:LYS:NZ	2.01	0.58
1:F:74:LEU:CD2	1:F:154:ILE:HD11	2.31	0.58
1:D:250:ARG:O	1:D:252:ILE:N	2.37	0.58
1:A:127:SER:OG	1:A:132:THR:HG22	2.03	0.58
1:D:118:ILE:HG23	1:D:179:VAL:HG22	1.86	0.57
1:B:190:MET:HE2	1:B:190:MET:HA	1.85	0.57
1:C:65:ILE:HD11	1:C:98:LEU:HD21	1.86	0.57
1:A:231:VAL:C	1:A:233:GLY:H	2.05	0.57
1:F:10:THR:HG23	1:F:67:PHE:HE2	1.69	0.57
1:B:141:MET:HE2	1:E:145:PHE:CD2	2.39	0.57
1:C:230:LEU:O	1:C:231:VAL:CG2	2.52	0.56
1:B:157:LYS:CD	1:E:130:VAL:HG22	2.34	0.56
1:F:171:GLY:O	1:F:173:SER:N	2.38	0.56
1:C:125:GLN:HE21	1:C:132:THR:HG21	1.71	0.56
1:C:192:VAL:O	1:C:195:THR:HB	2.06	0.56
1:E:216:MET:CE	1:E:221:ILE:HG13	2.34	0.56
1:F:125:GLN:HE21	1:F:132:THR:HG21	1.71	0.56
1:D:252:ILE:HG23	1:D:253:GLY:N	2.21	0.56
1:B:26:GLN:HE21	1:B:27:LYS:HZ3	1.52	0.55
1:F:37:THR:HG22	1:F:39:GLU:H	1.72	0.55
1:F:185:LEU:HD23	1:F:195:THR:HG21	1.88	0.55
1:A:118:ILE:CD1	1:A:179:VAL:HG13	2.36	0.55
1:B:5:PRO:HG2	1:B:29:LEU:HD22	1.89	0.55
1:E:107:ASN:OD1	1:E:110:ARG:N	2.40	0.55
1:F:133:THR:OG1	1:F:136:GLN:HG3	2.07	0.55
1:C:216:MET:CE	1:C:221:ILE:HD12	2.37	0.55
1:D:183:SER:O	1:D:186:HIS:HD2	1.89	0.54
1:B:204:ALA:HB2	1:E:204:ALA:HB2	1.89	0.54
1:A:157:LYS:HG2	1:C:91:LEU:HD13	1.88	0.54
1:B:216:MET:HE3	1:B:221:ILE:HB	1.90	0.54
1:C:195:THR:CG2	1:C:199:SER:OG	2.56	0.54
1:A:226:ILE:HG22	1:A:228:PRO:HD3	1.90	0.54
1:D:157:LYS:HG2	1:F:91:LEU:HD13	1.89	0.54
1:E:160:MET:HE3	1:E:216:MET:HE1	1.89	0.54
1:D:23:LYS:CE	1:D:265:GLU:OE2	2.56	0.54
1:B:183:SER:O	1:B:186:HIS:HD2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:65:ILE:HD11	1:D:73:TRP:CD2	2.43	0.54
1:E:93:GLN:HG3	1:E:94:LYS:N	2.23	0.53
1:F:88:ARG:HB2	1:F:89:PRO:HD3	1.91	0.53
1:A:10:THR:O	1:A:120:CYS:HB2	2.09	0.53
1:A:128:LEU:O	1:A:132:THR:CG2	2.56	0.53
1:C:231:VAL:HA	1:C:258:SER:O	2.09	0.53
1:D:250:ARG:O	1:D:251:THR:C	2.51	0.53
1:A:93:GLN:HE21	1:A:94:LYS:N	2.06	0.53
1:A:234:THR:CG2	1:A:235:ASP:N	2.70	0.53
1:B:216:MET:CE	1:B:221:ILE:HD12	2.38	0.53
1:B:195:THR:HG22	1:B:199:SER:OG	2.07	0.53
1:F:103:ASN:HD22	1:F:111:ARG:NH2	2.07	0.53
1:D:195:THR:O	1:D:195:THR:CG2	2.57	0.53
1:E:216:MET:HE3	1:E:219:ARG:HB2	1.91	0.52
1:A:234:THR:HG22	1:A:235:ASP:N	2.22	0.52
1:E:103:ASN:OD1	1:E:111:ARG:NH2	2.42	0.52
1:A:65:ILE:HD11	1:A:98:LEU:HD21	1.92	0.52
1:E:6:VAL:HG22	1:E:114:VAL:HG22	1.91	0.52
1:C:123:LEU:HD23	1:C:140:ILE:HA	1.92	0.51
1:E:46:ILE:HD11	1:E:51:LEU:HD13	1.91	0.51
1:D:127:SER:OG	1:D:132:THR:HG23	2.11	0.51
1:E:160:MET:CE	1:E:216:MET:HE1	2.40	0.51
1:A:195:THR:HG22	1:A:199:SER:OG	2.10	0.51
1:C:35:GLY:HA3	1:C:41:ILE:HG23	1.93	0.51
1:F:111:ARG:HD3	1:F:166:TRP:CH2	2.46	0.51
1:B:130:VAL:HG12	1:B:131:ARG:HG3	1.92	0.51
1:A:153:ASN:OD1	1:D:129:SER:OG	2.29	0.50
1:A:216:MET:HE3	1:A:221:ILE:HB	1.93	0.50
1:C:99:PHE:CZ	1:C:154:ILE:HG22	2.46	0.50
1:A:190:MET:HA	1:A:190:MET:HE2	1.93	0.50
1:D:190:MET:CE	1:D:248:LEU:CB	2.89	0.50
1:A:226:ILE:HD11	1:A:270:LEU:HD22	1.93	0.50
1:C:195:THR:CG2	1:C:195:THR:O	2.59	0.50
1:A:142:ASN:HA	1:A:146:MET:HB2	1.93	0.50
1:D:125:GLN:HE21	1:D:132:THR:HG21	1.77	0.50
1:F:5:PRO:HG2	1:F:29:LEU:HD22	1.94	0.50
1:F:195:THR:CG2	1:F:199:SER:OG	2.60	0.50
1:D:62:ALA:HB3	1:D:82:ILE:HG23	1.93	0.50
1:D:127:SER:OG	1:D:132:THR:CG2	2.60	0.49
1:E:125:GLN:HE21	1:E:132:THR:HG21	1.77	0.49
1:A:195:THR:O	1:A:195:THR:CG2	2.57	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231:VAL:O	1:A:233:GLY:N	2.45	0.49
1:C:183:SER:O	1:C:186:HIS:CD2	2.65	0.49
1:E:93:GLN:NE2	1:E:104:LYS:HD3	2.27	0.49
1:F:12:ALA:HB2	1:F:17:GLY:HA3	1.87	0.49
1:F:142:ASN:HA	1:F:146:MET:HB2	1.93	0.49
1:A:130:VAL:CG2	1:D:157:LYS:HG3	2.43	0.49
1:D:190:MET:HA	1:D:190:MET:HE2	1.95	0.49
1:A:82:ILE:HD11	1:A:90:PRO:CB	2.43	0.49
1:C:141:MET:HB3	1:C:146:MET:HE1	1.95	0.49
1:F:103:ASN:HD22	1:F:111:ARG:HH22	1.61	0.49
1:F:183:SER:O	1:F:186:HIS:CD2	2.62	0.49
1:A:231:VAL:O	1:A:232:LYS:C	2.56	0.48
1:F:125:GLN:NE2	1:F:132:THR:HG21	2.28	0.48
1:C:35:GLY:HA3	1:C:41:ILE:CG2	2.44	0.48
1:D:182:SER:HA	1:D:202:LYS:HD2	1.94	0.48
1:E:185:LEU:N	1:E:185:LEU:HD12	2.28	0.48
1:F:195:THR:HG22	1:F:199:SER:OG	2.13	0.48
1:F:190:MET:HE3	1:F:190:MET:HA	1.96	0.48
1:B:195:THR:CG2	1:B:199:SER:OG	2.62	0.47
1:F:195:THR:CG2	1:F:195:THR:O	2.62	0.47
1:E:133:THR:OG1	1:E:136:GLN:HG3	2.14	0.47
1:F:88:ARG:CB	1:F:89:PRO:CD	2.92	0.47
1:F:100:ASP:N	1:F:101:PRO:CD	2.77	0.47
1:A:231:VAL:HG23	1:A:233:GLY:CA	2.44	0.47
1:B:18:LYS:O	1:B:22:GLN:HB2	2.14	0.47
1:B:118:ILE:HD12	1:B:267:VAL:HG11	1.96	0.47
1:B:207:ARG:NH2	1:E:190:MET:O	2.48	0.47
1:E:46:ILE:HG12	1:E:60:GLN:NE2	2.29	0.47
1:C:141:MET:HB3	1:C:146:MET:CE	2.45	0.47
1:A:182:SER:O	1:A:228:PRO:HD2	2.15	0.47
1:F:166:TRP:HB2	1:F:169:LEU:HD22	1.96	0.47
1:C:151:MET:HE3	1:C:151:MET:HA	1.96	0.46
1:E:133:THR:H	1:E:136:GLN:HE21	1.63	0.46
1:B:57:TYR:C	1:B:57:TYR:CD2	2.94	0.46
1:D:183:SER:O	1:D:186:HIS:CD2	2.69	0.46
1:D:62:ALA:HB3	1:D:82:ILE:CG2	2.45	0.46
1:A:160:MET:HG2	1:C:88:ARG:HH21	1.81	0.46
1:B:141:MET:HE2	1:E:145:PHE:CE2	2.50	0.46
1:A:118:ILE:HD13	1:A:179:VAL:CG1	2.39	0.46
1:E:99:PHE:CZ	1:E:154:ILE:HG22	2.51	0.46
1:A:130:VAL:HG22	1:D:157:LYS:HG3	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:185:LEU:HD23	1:B:195:THR:HG21	1.97	0.46
1:B:226:ILE:HG22	1:B:228:PRO:HD3	1.99	0.46
1:E:37:THR:HG23	1:E:40:SER:HB2	1.99	0.45
1:E:182:SER:OG	1:E:183:SER:N	2.45	0.45
1:B:120:CYS:HA	1:B:181:ILE:HD12	1.98	0.45
1:C:98:LEU:HD13	1:C:151:MET:HE1	1.96	0.45
1:D:182:SER:O	1:D:228:PRO:HD2	2.16	0.45
1:C:230:LEU:C	1:C:231:VAL:HG23	2.41	0.45
1:A:230:LEU:HD13	1:A:256:GLY:O	2.17	0.45
1:D:137:ILE:HG12	1:D:197:VAL:HG21	1.99	0.45
1:D:166:TRP:HZ3	1:F:108:ASN:HD22	1.65	0.45
1:B:46:ILE:HG12	1:B:60:GLN:NE2	2.32	0.44
1:B:195:THR:CG2	1:B:195:THR:O	2.65	0.44
1:C:185:LEU:CD2	1:C:195:THR:HG21	2.40	0.44
1:D:250:ARG:C	1:D:252:ILE:N	2.73	0.44
1:E:151:MET:HE3	1:E:151:MET:HA	1.98	0.44
1:E:217:GLU:N	1:E:218:PRO:HD2	2.33	0.44
1:D:166:TRP:O	1:D:169:LEU:HB2	2.18	0.44
1:F:152:THR:HG23	1:F:178:ILE:HG21	1.99	0.44
1:B:41:ILE:HD11	1:B:63:ILE:C	2.42	0.44
1:B:119:ASN:HB2	1:B:180:ASN:OD1	2.18	0.44
1:D:74:LEU:HD11	1:D:150:THR:HG21	2.00	0.44
1:B:54:GLY:O	1:B:86:LYS:NZ	2.43	0.44
1:B:108:ASN:H	1:B:108:ASN:HD22	1.65	0.44
1:C:99:PHE:CE2	1:C:154:ILE:HG22	2.53	0.44
1:B:26:GLN:NE2	1:B:27:LYS:NZ	2.65	0.44
1:C:119:ASN:HB2	1:C:180:ASN:OD1	2.17	0.44
1:A:184:ILE:HG23	1:A:185:LEU:HD12	2.00	0.44
1:A:183:SER:O	1:A:186:HIS:HD2	2.01	0.44
1:D:252:ILE:CG2	1:D:253:GLY:N	2.78	0.44
1:C:99:PHE:HE2	1:C:154:ILE:CG2	2.31	0.43
1:E:145:PHE:CE1	1:E:204:ALA:HB1	2.53	0.43
1:D:98:LEU:HD13	1:D:151:MET:CE	2.44	0.43
1:E:10:THR:HG21	1:E:119:ASN:OD1	2.18	0.43
1:A:147:SER:HB2	1:A:148:PRO:HD3	1.99	0.43
1:C:51:LEU:HD21	1:C:84:TYR:HB3	2.01	0.43
1:C:230:LEU:O	1:C:231:VAL:CB	2.65	0.43
1:D:85:PHE:HE2	1:D:91:LEU:HB2	1.82	0.43
1:F:217:GLU:N	1:F:218:PRO:HD2	2.34	0.43
1:A:192:VAL:O	1:A:195:THR:HB	2.18	0.43
1:B:74:LEU:HD11	1:B:150:THR:HG21	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:108:ASN:HD22	1:B:108:ASN:N	2.17	0.43
1:A:46:ILE:N	1:A:60:GLN:HE22	2.10	0.43
1:C:195:THR:O	1:C:195:THR:HG23	2.18	0.43
1:F:127:SER:OG	1:F:132:THR:CG2	2.64	0.43
1:A:156:ILE:HG13	1:A:178:ILE:CD1	2.49	0.43
1:A:182:SER:HA	1:A:202:LYS:HD2	2.01	0.43
1:B:147:SER:HB2	1:B:148:PRO:HD3	2.01	0.43
1:C:125:GLN:HG2	1:C:194:GLY:O	2.19	0.43
1:D:147:SER:HB2	1:D:148:PRO:HD3	2.01	0.42
1:B:118:ILE:HG23	1:B:179:VAL:HG22	2.01	0.42
1:D:8:ILE:HG12	1:D:32:ILE:HB	2.00	0.42
1:E:4:LEU:HA	1:E:5:PRO:HD3	1.92	0.42
1:E:255:SER:OG	1:E:256:GLY:N	2.52	0.42
1:B:137:ILE:HG12	1:B:197:VAL:HG21	2.02	0.42
1:D:166:TRP:CZ3	1:F:107:ASN:HB2	2.55	0.42
1:A:89:PRO:HA	1:A:90:PRO:HD3	1.88	0.42
1:C:190:MET:HE2	1:C:190:MET:HA	2.01	0.42
1:B:70:TRP:CZ3	1:B:146:MET:HG2	2.54	0.42
1:D:81:GLY:HA3	1:D:95:TYR:HE2	1.84	0.42
1:F:12:ALA:CB	1:F:17:GLY:CA	2.83	0.42
1:A:216:MET:CE	1:A:221:ILE:HB	2.50	0.42
1:C:74:LEU:HD11	1:C:150:THR:HG21	2.02	0.42
1:C:145:PHE:CE1	1:C:204:ALA:HB1	2.55	0.42
1:D:23:LYS:NZ	1:D:27:LYS:HE3	2.33	0.42
1:E:82:ILE:HD11	1:E:84:TYR:CZ	2.54	0.42
1:D:226:ILE:HD11	1:D:270:LEU:HD22	2.02	0.42
1:B:128:LEU:HD23	1:B:131:ARG:NH1	2.35	0.41
1:D:119:ASN:HB2	1:D:180:ASN:OD1	2.20	0.41
1:E:67:PHE:CD1	1:E:147:SER:HB3	2.56	0.41
1:A:74:LEU:HD11	1:A:150:THR:HG21	2.03	0.41
1:E:158:TYR:CD2	1:E:158:TYR:C	2.98	0.41
1:B:97:THR:H	1:B:100:ASP:HB2	1.85	0.41
1:D:13:THR:HG21	1:D:122:GLY:O	2.20	0.41
1:E:125:GLN:NE2	1:E:132:THR:HG21	2.35	0.41
1:D:6:VAL:HG13	1:D:113:TYR:O	2.21	0.41
1:D:25:PHE:HB3	1:D:50:GLN:NE2	2.35	0.41
1:A:65:ILE:HD11	1:A:98:LEU:CD2	2.50	0.41
1:C:12:ALA:HB1	1:C:17:GLY:HA3	2.02	0.41
1:F:185:LEU:HD23	1:F:195:THR:CG2	2.51	0.41
1:B:16:ILE:HD11	1:B:231:VAL:HG21	2.02	0.41
1:B:18:LYS:NZ	1:B:47:ASP:OD1	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:7:ALA:HA	1:E:116:LEU:O	2.21	0.41
1:E:123:LEU:HD23	1:E:140:ILE:HA	2.03	0.41
1:E:186:HIS:CD2	1:E:187:SER:HB3	2.55	0.41
1:A:125:GLN:HE21	1:A:132:THR:HG21	1.86	0.41
1:A:230:LEU:CD1	1:A:230:LEU:N	2.30	0.40
1:D:7:ALA:HA	1:D:116:LEU:O	2.21	0.40
1:B:263:ILE:O	1:B:267:VAL:HG23	2.22	0.40
1:C:99:PHE:CE2	1:C:154:ILE:CG2	3.05	0.40
1:A:230:LEU:CD1	1:A:256:GLY:O	2.70	0.40
1:E:87:ASP:OD1	1:E:87:ASP:N	2.52	0.40
1:F:169:LEU:HD12	1:F:169:LEU:HA	1.67	0.40
1:A:93:GLN:HE21	1:A:94:LYS:H	1.68	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	250/286 (87%)	245 (98%)	5 (2%)	0	100	100
1	B	249/286 (87%)	242 (97%)	7 (3%)	0	100	100
1	C	244/286 (85%)	240 (98%)	4 (2%)	0	100	100
1	D	262/286 (92%)	253 (97%)	9 (3%)	0	100	100
1	E	244/286 (85%)	239 (98%)	5 (2%)	0	100	100
1	F	240/286 (84%)	230 (96%)	10 (4%)	0	100	100
All	All	1489/1716 (87%)	1449 (97%)	40 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	187/250 (75%)	171 (91%)	16 (9%)	10	22
1	B	199/250 (80%)	173 (87%)	26 (13%)	4	8
1	C	194/250 (78%)	173 (89%)	21 (11%)	6	13
1	D	194/250 (78%)	167 (86%)	27 (14%)	3	7
1	E	185/250 (74%)	164 (89%)	21 (11%)	5	11
1	F	185/250 (74%)	169 (91%)	16 (9%)	10	22
All	All	1144/1500 (76%)	1017 (89%)	127 (11%)	6	12

All (127) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	THR
1	A	22	GLN
1	A	61	CYS
1	A	82	ILE
1	A	91	LEU
1	A	93	GLN
1	A	108	ASN
1	A	118	ILE
1	A	128	LEU
1	A	130	VAL
1	A	132	THR
1	A	154	ILE
1	A	179	VAL
1	A	230	LEU
1	A	257	THR
1	A	270	LEU
1	B	6	VAL
1	B	10	THR
1	B	13	THR
1	B	22	GLN
1	B	37	THR
1	B	41	ILE

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Mol	Chain	Res	Type
1	B	51	LEU
1	B	55	LEU
1	B	59	ARG
1	B	74	LEU
1	B	78	SER
1	B	91	LEU
1	B	98	LEU
1	B	108	ASN
1	B	118	ILE
1	B	128	LEU
1	B	130	VAL
1	B	135	SER
1	B	151	MET
1	B	172	GLN
1	B	179	VAL
1	B	195	THR
1	B	255	SER
1	B	258	SER
1	B	269	SER
1	B	270	LEU
1	C	6	VAL
1	C	14	ARG
1	C	29	LEU
1	C	55	LEU
1	C	59	ARG
1	C	65	ILE
1	C	88	ARG
1	C	96	SER
1	C	106	SER
1	C	108	ASN
1	C	128	LEU
1	C	130	VAL
1	C	151	MET
1	C	179	VAL
1	C	182	SER
1	C	195	THR
1	C	196	SER
1	C	225	THR
1	C	230	LEU
1	C	255	SER
1	C	270	LEU
1	D	1	MET

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Mol	Chain	Res	Type
1	D	2	HIS
1	D	6	VAL
1	D	10	THR
1	D	13	THR
1	D	23	LYS
1	D	29	LEU
1	D	46	ILE
1	D	51	LEU
1	D	82	ILE
1	D	88	ARG
1	D	91	LEU
1	D	96	SER
1	D	104	LYS
1	D	108	ASN
1	D	130	VAL
1	D	132	THR
1	D	151	MET
1	D	172	GLN
1	D	179	VAL
1	D	182	SER
1	D	195	THR
1	D	210	GLU
1	D	226	ILE
1	D	230	LEU
1	D	252	ILE
1	D	270	LEU
1	E	6	VAL
1	E	10	THR
1	E	29	LEU
1	E	37	THR
1	E	51	LEU
1	E	52	GLN
1	E	82	ILE
1	E	88	ARG
1	E	91	LEU
1	E	128	LEU
1	E	130	VAL
1	E	132	THR
1	E	179	VAL
1	E	191	LYS
1	E	196	SER
1	E	197	VAL

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Mol	Chain	Res	Type
1	E	207	ARG
1	E	221	ILE
1	E	230	LEU
1	E	231	VAL
1	E	270	LEU
1	F	6	VAL
1	F	10	THR
1	F	16	ILE
1	F	51	LEU
1	F	82	ILE
1	F	88	ARG
1	F	114	VAL
1	F	128	LEU
1	F	130	VAL
1	F	132	THR
1	F	164	ARG
1	F	169	LEU
1	F	179	VAL
1	F	184	ILE
1	F	195	THR
1	F	258	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
1	A	93	GLN
1	A	108	ASN
1	A	136	GLN
1	A	186	HIS
1	B	26	GLN
1	B	60	GLN
1	B	93	GLN
1	B	103	ASN
1	B	108	ASN
1	B	115	ASN
1	B	136	GLN
1	B	153	ASN
1	B	186	HIS
1	C	50	GLN
1	C	93	GLN
1	C	108	ASN

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Mol	Chain	Res	Type
1	C	115	ASN
1	C	136	GLN
1	C	153	ASN
1	C	186	HIS
1	D	2	HIS
1	D	115	ASN
1	D	172	GLN
1	D	186	HIS
1	E	60	GLN
1	E	136	GLN
1	E	153	ASN
1	F	103	ASN
1	F	108	ASN
1	F	136	GLN
1	F	153	ASN
1	F	186	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	254/286 (88%)	0.33	14 (5%) 30 25	28, 45, 68, 83	0
1	B	253/286 (88%)	0.07	9 (3%) 46 40	21, 39, 54, 63	0
1	C	250/286 (87%)	0.00	5 (2%) 65 60	20, 37, 54, 60	0
1	D	268/286 (93%)	0.37	17 (6%) 26 20	20, 44, 66, 90	0
1	E	250/286 (87%)	0.35	14 (5%) 30 24	23, 45, 66, 69	0
1	F	246/286 (86%)	0.27	14 (5%) 29 24	32, 45, 71, 78	0
All	All	1521/1716 (88%)	0.23	73 (4%) 35 30	20, 42, 64, 90	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	235	ASP	4.7
1	D	235	ASP	4.5
1	D	242	VAL	4.5
1	A	234	THR	4.5
1	A	122	GLY	4.4
1	D	244	ALA	4.3
1	B	233	GLY	4.3
1	E	274	THR	4.1
1	C	230	LEU	3.9
1	D	241	PRO	3.9
1	E	107	ASN	3.8
1	E	12	ALA	3.7
1	A	232	LYS	3.7
1	F	48	ARG	3.7
1	F	229	GLY	3.6
1	A	47	ASP	3.6
1	D	243	GLU	3.6
1	D	1	MET	3.5
1	F	47	ASP	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	233	GLY	3.4
1	E	231	VAL	3.4
1	D	250	ARG	3.3
1	F	13	THR	3.2
1	C	172	GLN	3.2
1	F	256	GLY	3.1
1	D	252	ILE	3.1
1	B	1	MET	3.1
1	D	65	ILE	3.0
1	D	253	GLY	3.0
1	B	2	HIS	2.9
1	C	274	THR	2.9
1	F	220	ASN	2.9
1	E	232	LYS	2.9
1	D	234	THR	2.9
1	D	251	THR	2.9
1	F	1	MET	2.8
1	C	231	VAL	2.8
1	A	1	MET	2.8
1	A	42	GLU	2.8
1	F	173	SER	2.8
1	E	256	GLY	2.7
1	F	12	ALA	2.6
1	A	173	SER	2.5
1	B	232	LYS	2.5
1	F	37	THR	2.5
1	D	274	THR	2.5
1	A	255	SER	2.5
1	E	186	HIS	2.5
1	F	39	GLU	2.5
1	B	274	THR	2.4
1	D	246	GLU	2.4
1	F	171	GLY	2.4
1	E	255	SER	2.4
1	B	170	SER	2.4
1	A	48	ARG	2.3
1	D	232	LYS	2.3
1	E	47	ASP	2.3
1	F	42	GLU	2.2
1	F	25	PHE	2.2
1	A	220	ASN	2.2
1	E	41	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	E	48	ARG	2.2
1	E	55	LEU	2.1
1	E	145	PHE	2.1
1	C	80	ASP	2.1
1	E	173	SER	2.1
1	B	14	ARG	2.1
1	D	249	GLU	2.1
1	D	49	GLY	2.1
1	B	220	ASN	2.0
1	A	226	ILE	2.0
1	B	13	THR	2.0
1	A	50	GLN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.