



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:16 AM GMT

PDB ID : 3E1H
Title : Crystal structure of a type III polyketide synthase PKSIIINc from *Neurospora crassa*
Authors : Goyal, A.; Rahman, A.; Sankaranarayanan, R.
Deposited on : 2008-08-04
Resolution : 2.58 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

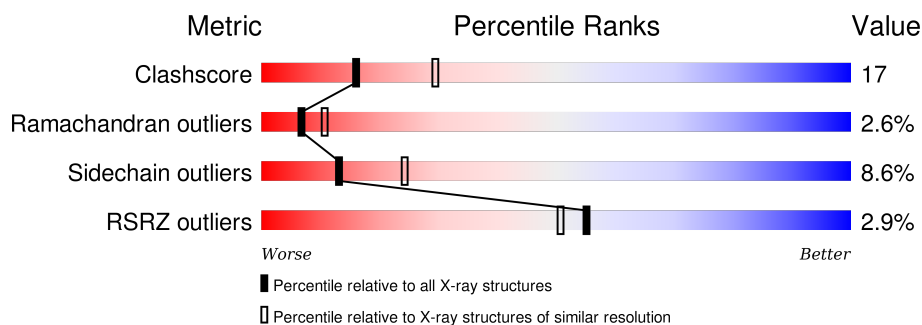
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.58 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	102246	3003 (2.60-2.56)
Ramachandran outliers	100387	2956 (2.60-2.56)
Sidechain outliers	100360	2956 (2.60-2.56)
RSRZ outliers	91569	2642 (2.60-2.56)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	465	<div> <div>2%</div> <div> <div></div> <div>55%</div> <div>25%</div> <div>•</div> <div>18%</div> </div> </div>
1	B	465	<div> <div>3%</div> <div> <div></div> <div>56%</div> <div>22%</div> <div>•</div> <div>18%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6181 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 Putative uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	383	Total	C	N	O	S	0	0	0
			2884	1814	503	548	19			
1	B	383	Total	C	N	O	S	0	0	0
			2884	1814	503	548	19			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP Q7S6N4
A	-18	GLY	-	EXPRESSION TAG	UNP Q7S6N4
A	-17	SER	-	EXPRESSION TAG	UNP Q7S6N4
A	-16	SER	-	EXPRESSION TAG	UNP Q7S6N4
A	-15	HIS	-	EXPRESSION TAG	UNP Q7S6N4
A	-14	HIS	-	EXPRESSION TAG	UNP Q7S6N4
A	-13	HIS	-	EXPRESSION TAG	UNP Q7S6N4
A	-12	HIS	-	EXPRESSION TAG	UNP Q7S6N4
A	-11	HIS	-	EXPRESSION TAG	UNP Q7S6N4
A	-10	HIS	-	EXPRESSION TAG	UNP Q7S6N4
A	-9	SER	-	EXPRESSION TAG	UNP Q7S6N4
A	-8	SER	-	EXPRESSION TAG	UNP Q7S6N4
A	-7	GLY	-	EXPRESSION TAG	UNP Q7S6N4
A	-6	LEU	-	EXPRESSION TAG	UNP Q7S6N4
A	-5	VAL	-	EXPRESSION TAG	UNP Q7S6N4
A	-4	PRO	-	EXPRESSION TAG	UNP Q7S6N4
A	-3	ARG	-	EXPRESSION TAG	UNP Q7S6N4
A	-2	GLY	-	EXPRESSION TAG	UNP Q7S6N4
A	-1	SER	-	EXPRESSION TAG	UNP Q7S6N4
A	0	HIS	-	EXPRESSION TAG	UNP Q7S6N4
B	-19	MET	-	EXPRESSION TAG	UNP Q7S6N4
B	-18	GLY	-	EXPRESSION TAG	UNP Q7S6N4
B	-17	SER	-	EXPRESSION TAG	UNP Q7S6N4
B	-16	SER	-	EXPRESSION TAG	UNP Q7S6N4
B	-15	HIS	-	EXPRESSION TAG	UNP Q7S6N4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	EXPRESSION TAG	UNP Q7S6N4
B	-13	HIS	-	EXPRESSION TAG	UNP Q7S6N4
B	-12	HIS	-	EXPRESSION TAG	UNP Q7S6N4
B	-11	HIS	-	EXPRESSION TAG	UNP Q7S6N4
B	-10	HIS	-	EXPRESSION TAG	UNP Q7S6N4
B	-9	SER	-	EXPRESSION TAG	UNP Q7S6N4
B	-8	SER	-	EXPRESSION TAG	UNP Q7S6N4
B	-7	GLY	-	EXPRESSION TAG	UNP Q7S6N4
B	-6	LEU	-	EXPRESSION TAG	UNP Q7S6N4
B	-5	VAL	-	EXPRESSION TAG	UNP Q7S6N4
B	-4	PRO	-	EXPRESSION TAG	UNP Q7S6N4
B	-3	ARG	-	EXPRESSION TAG	UNP Q7S6N4
B	-2	GLY	-	EXPRESSION TAG	UNP Q7S6N4
B	-1	SER	-	EXPRESSION TAG	UNP Q7S6N4
B	0	HIS	-	EXPRESSION TAG	UNP Q7S6N4

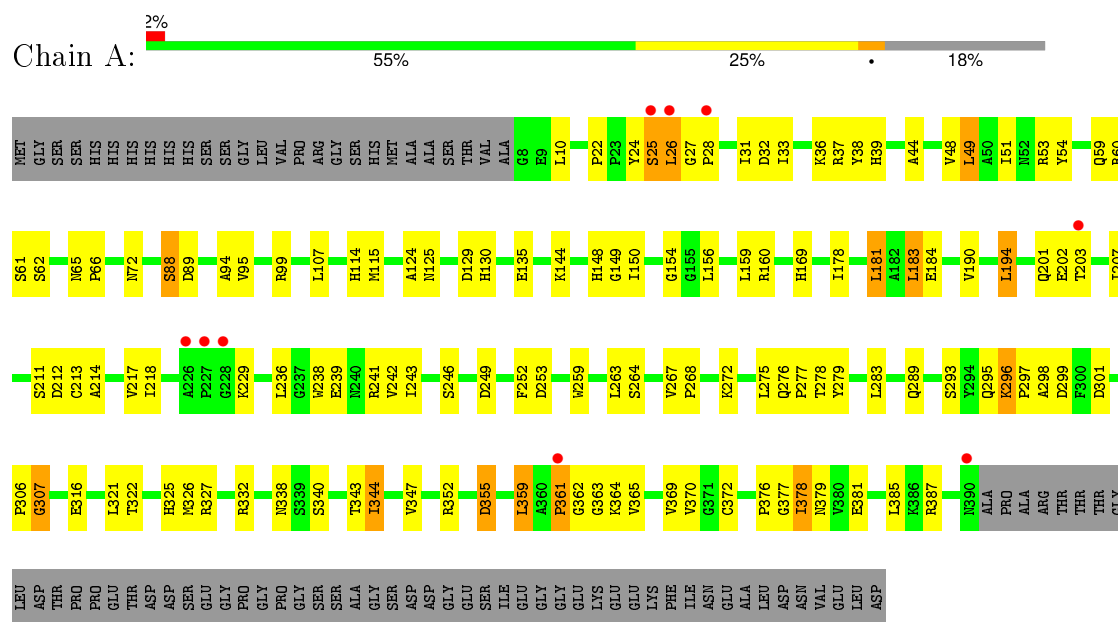
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	209	Total O 209 209	0	0
2	B	204	Total O 204 204	0	0

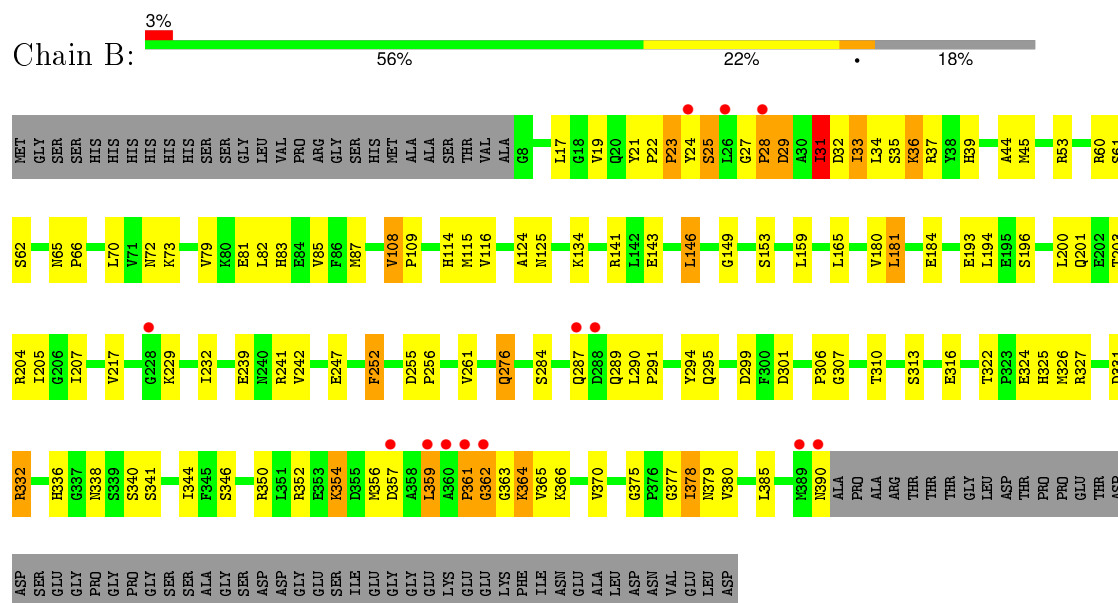
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Putative uncharacterized protein



• Molecule 1: Putative uncharacterized protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	70.02Å 104.96Å 105.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.58 24.83 – 2.59	Depositor EDS
% Data completeness (in resolution range)	98.0 (25.00-2.58) 98.0 (24.83-2.59)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 2.60Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.213 , 0.260 0.210 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	39.5	Xtriage
Anisotropy	0.313	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 43.0	EDS
Estimated twinning fraction	0.024 for -h,l,k	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 24322 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6181	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.18% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.45	0/2945	0.65	1/4001 (0.0%)
1	B	0.45	0/2945	0.66	0/4001
All	All	0.45	0/5890	0.65	1/8002 (0.0%)

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
All	All	0	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	181	LEU	CA-CB-CG	5.54	128.03	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	377	GLY	Peptide
1	B	377	GLY	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	2884	0	2894	106	0
1	B	2884	0	2894	98	0
2	A	209	0	0	5	0
2	B	204	0	0	6	0
All	All	6181	0	5788	195	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 17.

All (195) 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:B:344:ILE:HD12	1:B:344:ILE:H	1.26	1.00
1:B:344:ILE:H	1:B:344:ILE:CD1	1.86	0.87
1:A:144:LYS:HE3	2:A:516:HOH:O	1.79	0.82
1:A:114:HIS:HB2	1:A:178:ILE:HG12	1.63	0.81
1:A:37:ARG:HH11	1:A:72:ASN:HD21	1.28	0.78
1:B:37:ARG:HH21	1:B:72:ASN:HD21	1.30	0.78
1:A:361:PRO:HG2	1:A:364:LYS:HB2	1.65	0.78
1:A:307:GLY:HA3	1:A:338:ASN:ND2	1.98	0.78
1:B:344:ILE:HD12	1:B:344:ILE:N	1.98	0.77
1:A:28:PRO:O	1:A:31:ILE:HG12	1.87	0.74
1:A:327:ARG:HG3	1:A:355:ASP:OD1	1.88	0.74
1:A:236:LEU:HD22	1:A:289:GLN:HE22	1.52	0.74
1:A:307:GLY:HA3	1:A:338:ASN:HD22	1.53	0.72
1:B:184:GLU:HG2	1:B:340:SER:HB3	1.70	0.72
1:A:344:ILE:HD13	1:A:344:ILE:H	1.55	0.72
1:B:350:ARG:HH22	1:B:354:LYS:HE2	1.54	0.72
1:A:307:GLY:CA	1:A:338:ASN:HD22	2.04	0.71
1:B:361:PRO:HG2	1:B:364:LYS:HB2	1.73	0.70
1:B:25:SER:HB2	1:B:61:SER:HA	1.72	0.70
1:A:24:TYR:O	1:A:25:SER:C	2.29	0.70
1:B:17:LEU:HD23	1:B:217:VAL:HG12	1.73	0.69
1:A:31:ILE:HD12	1:A:60:ARG:HH21	1.57	0.68
1:B:322:THR:H	1:B:325:HIS:CD2	2.11	0.68
1:B:322:THR:OG1	1:B:325:HIS:HD2	1.76	0.68
1:A:33:ILE:HD12	1:A:53:ARG:HH21	1.60	0.67
1:B:331:ASP:OD2	1:B:332:ARG:NH1	2.29	0.66
1:B:31:ILE:HD12	1:B:60:ARG:NH2	2.11	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:THR:H	1:A:325:HIS:CD2	2.14	0.65
1:B:313:SER:HA	1:B:316:GLU:HG2	1.77	0.64
1:B:37:ARG:HH21	1:B:72:ASN:ND2	1.95	0.64
1:B:344:ILE:HG23	2:B:604:HOH:O	1.96	0.64
1:A:37:ARG:HH11	1:A:72:ASN:ND2	1.95	0.64
1:A:36:LYS:HD3	1:A:49:LEU:HD21	1.78	0.64
1:A:344:ILE:HD13	1:A:344:ILE:N	2.13	0.64
1:B:307:GLY:HA2	1:B:338:ASN:HB2	1.79	0.63
1:A:31:ILE:HG13	1:A:32:ASP:H	1.63	0.63
1:B:359:LEU:HD11	1:B:365:VAL:HA	1.79	0.63
1:A:39:HIS:HE1	2:A:455:HOH:O	1.82	0.63
1:B:23:PRO:C	1:B:25:SER:H	2.02	0.62
1:B:359:LEU:HD21	1:B:365:VAL:N	2.15	0.61
1:B:31:ILE:HD12	1:B:60:ARG:HH21	1.65	0.61
1:B:361:PRO:HG2	1:B:364:LYS:H	1.66	0.60
1:A:359:LEU:HD11	1:A:365:VAL:HG22	1.82	0.60
1:B:295:GLN:HG2	2:B:601:HOH:O	2.01	0.60
1:A:307:GLY:CA	1:A:338:ASN:ND2	2.65	0.60
1:A:38:TYR:OH	1:A:72:ASN:HA	2.02	0.59
1:B:108:VAL:HG22	1:B:109:PRO:HD2	1.83	0.59
1:B:25:SER:HA	1:B:62:SER:H	1.67	0.59
1:A:31:ILE:CD1	1:A:60:ARG:HH21	2.15	0.59
1:A:184:GLU:HG2	1:A:340:SER:HB3	1.84	0.59
1:B:359:LEU:HD22	1:B:365:VAL:CG2	2.32	0.59
1:B:359:LEU:CD1	1:B:365:VAL:HA	2.33	0.59
1:B:35:SER:C	1:B:37:ARG:H	2.07	0.58
1:A:190:VAL:HA	1:A:207:ILE:HD11	1.85	0.58
1:B:322:THR:OG1	1:B:325:HIS:CD2	2.56	0.58
1:B:203:THR:O	1:B:204:ARG:HB2	2.04	0.57
1:B:344:ILE:HG12	2:B:522:HOH:O	2.05	0.57
1:B:313:SER:HA	1:B:316:GLU:CG	2.34	0.57
1:A:33:ILE:CD1	1:A:53:ARG:HH21	2.17	0.57
1:A:190:VAL:HA	1:A:207:ILE:CD1	2.35	0.56
1:A:37:ARG:NH1	1:A:72:ASN:HD21	2.01	0.56
1:A:344:ILE:HG13	2:A:532:HOH:O	2.06	0.56
1:B:242:VAL:HG22	1:B:379:ASN:OD1	2.06	0.56
1:A:242:VAL:HG12	1:A:379:ASN:OD1	2.06	0.56
1:A:160:ARG:NH2	2:A:452:HOH:O	2.38	0.55
1:A:65:ASN:HB2	1:A:66:PRO:HD2	1.87	0.55
1:A:184:GLU:HG2	1:A:340:SER:CB	2.37	0.55
1:A:241:ARG:HE	1:B:141:ARG:NH1	2.04	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:241:ARG:HH22	1:A:277:PRO:HB2	1.70	0.55
1:A:362:GLY:C	1:A:364:LYS:H	2.10	0.55
1:B:252:PHE:HE2	1:B:261:VAL:HG12	1.70	0.54
1:B:361:PRO:CG	1:B:364:LYS:HB2	2.37	0.54
1:A:344:ILE:CD1	1:A:344:ILE:H	2.12	0.54
1:B:301:ASP:OD1	1:B:327:ARG:NH2	2.40	0.54
1:A:60:ARG:HG2	1:A:211:SER:HB3	1.88	0.54
1:B:241:ARG:HG2	1:B:242:VAL:N	2.22	0.54
1:A:44:ALA:H	1:A:201:GLN:HE22	1.56	0.54
1:B:361:PRO:HB3	2:B:629:HOH:O	2.07	0.54
1:B:359:LEU:HD21	1:B:364:LYS:C	2.27	0.54
1:B:33:ILE:HG13	1:B:53:ARG:HH21	1.72	0.54
1:A:246:SER:HB2	1:A:378:ILE:HG13	1.90	0.53
1:B:32:ASP:O	1:B:36:LYS:HB2	2.09	0.53
1:A:62:SER:HA	1:A:212:ASP:OD2	2.09	0.53
1:A:253:ASP:O	1:A:259:TRP:HA	2.09	0.53
1:A:39:HIS:CE1	2:A:455:HOH:O	2.60	0.53
1:A:148:HIS:HB3	1:B:146:LEU:HD13	1.90	0.53
1:B:359:LEU:HD22	1:B:365:VAL:HG23	1.91	0.52
1:B:324:GLU:O	1:B:327:ARG:HB3	2.10	0.52
1:A:283:LEU:HD21	1:A:295:GLN:O	2.09	0.52
1:B:17:LEU:HD23	1:B:217:VAL:CG1	2.40	0.52
1:A:31:ILE:CD1	1:A:60:ARG:NH2	2.73	0.52
1:B:375:GLY:O	1:B:378:ILE:HA	2.10	0.51
1:A:31:ILE:HD12	1:A:60:ARG:NH2	2.24	0.50
1:A:322:THR:H	1:A:325:HIS:HD2	1.58	0.50
1:A:37:ARG:HG3	1:A:38:TYR:CE1	2.46	0.50
1:B:116:VAL:HG23	1:B:116:VAL:O	2.11	0.50
1:A:60:ARG:NH1	1:A:207:ILE:O	2.35	0.50
1:A:301:ASP:OD1	1:A:327:ARG:NH2	2.45	0.50
1:A:343:THR:O	1:A:347:VAL:HG23	2.12	0.50
1:A:298:ALA:O	1:A:325:HIS:HE1	1.95	0.49
1:B:359:LEU:HD11	1:B:364:LYS:O	2.12	0.49
1:A:322:THR:OG1	1:A:325:HIS:HD2	1.94	0.49
1:A:276:GLN:HB3	1:A:277:PRO:HD3	1.93	0.49
1:B:70:LEU:HD22	1:B:85:VAL:HG21	1.95	0.49
1:A:296:LYS:HG3	1:A:299:ASP:OD1	2.12	0.49
1:B:65:ASN:HB3	1:B:66:PRO:HD2	1.95	0.49
1:A:51:ILE:HA	1:A:54:TYR:CD2	2.48	0.48
1:A:25:SER:HA	1:A:61:SER:HA	1.96	0.48
1:B:79:VAL:HA	1:B:82:LEU:HD12	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:193:GLU:CB	1:B:207:ILE:HG23	2.44	0.48
1:A:307:GLY:HA2	1:A:338:ASN:HD22	1.79	0.48
1:B:357:ASP:C	1:B:359:LEU:H	2.17	0.48
1:A:241:ARG:HG2	1:A:242:VAL:N	2.29	0.48
1:A:26:LEU:HD22	1:A:59:GLN:HG3	1.96	0.48
1:A:190:VAL:O	1:A:194:LEU:HB2	2.14	0.48
1:A:95:VAL:HG21	1:A:135:GLU:HG2	1.96	0.47
1:A:272:LYS:HA	1:A:275:LEU:HD12	1.95	0.47
1:B:196:SER:O	1:B:200:LEU:HB2	2.15	0.47
1:A:94:ALA:HB2	1:A:183:LEU:HG	1.97	0.47
1:B:361:PRO:CB	2:B:629:HOH:O	2.63	0.47
1:A:22:PRO:O	1:A:25:SER:HB2	2.14	0.47
1:B:23:PRO:C	1:B:25:SER:N	2.67	0.47
1:B:359:LEU:HD22	1:B:365:VAL:HG22	1.96	0.47
1:A:44:ALA:H	1:A:201:GLN:NE2	2.13	0.47
1:A:316:GLU:HG3	1:A:326:MET:CE	2.45	0.47
1:A:25:SER:O	1:A:27:GLY:N	2.48	0.46
1:B:299:ASP:O	1:B:366:LYS:HD2	2.14	0.46
1:B:291:PRO:HG2	1:B:294:TYR:HD1	1.81	0.46
1:B:350:ARG:HH22	1:B:354:LYS:CE	2.27	0.46
1:A:65:ASN:HB2	1:A:66:PRO:CD	2.45	0.46
1:A:279:TYR:CE1	1:A:297:PRO:HD3	2.51	0.46
1:A:296:LYS:O	1:A:297:PRO:C	2.55	0.46
1:A:359:LEU:HD22	1:A:359:LEU:H	1.80	0.46
1:A:344:ILE:CD1	1:A:344:ILE:N	2.76	0.46
1:A:124:ALA:O	1:B:149:GLY:HA2	2.16	0.46
1:B:19:VAL:HG21	1:B:346:SER:HA	1.98	0.45
1:A:88:SER:OG	1:A:89:ASP:OD1	2.28	0.45
1:B:29:ASP:OD1	1:B:29:ASP:N	2.50	0.45
1:A:316:GLU:HA	1:A:321:LEU:HB2	1.99	0.45
1:A:207:ILE:HA	1:A:207:ILE:HD12	1.83	0.45
1:B:359:LEU:HG	1:B:361:PRO:HD2	1.98	0.45
1:B:21:TYR:CE2	1:B:336:HIS:HB3	2.51	0.45
1:B:350:ARG:NH2	1:B:354:LYS:HE2	2.28	0.45
1:B:361:PRO:O	1:B:362:GLY:C	2.55	0.45
1:A:160:ARG:HG3	1:A:238:TRP:CD1	2.52	0.45
1:B:27:GLY:HA2	1:B:28:PRO:HD3	1.79	0.45
1:B:301:ASP:CG	1:B:327:ARG:HH22	2.20	0.44
1:B:327:ARG:NH2	1:B:359:LEU:HD12	2.32	0.44
1:A:316:GLU:O	1:A:321:LEU:N	2.44	0.44
1:A:252:PHE:HB2	1:B:125:ASN:OD1	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:HIS:HE2	1:B:247:GLU:CG	2.31	0.44
1:B:252:PHE:CE2	1:B:261:VAL:HG12	2.52	0.44
1:A:31:ILE:HG13	1:A:32:ASP:N	2.31	0.44
1:B:22:PRO:O	1:B:25:SER:HB3	2.18	0.44
1:B:23:PRO:O	1:B:25:SER:N	2.41	0.43
1:B:114:HIS:HA	1:B:143:GLU:O	2.18	0.43
1:B:255:ASP:CG	1:B:256:PRO:HD2	2.40	0.43
1:B:29:ASP:C	1:B:31:ILE:H	2.22	0.43
1:A:361:PRO:HB2	1:A:362:GLY:H	1.64	0.42
1:A:249:ASP:HA	1:A:264:SER:HB3	2.00	0.42
1:A:338:ASN:C	1:A:338:ASN:OD1	2.56	0.42
1:B:39:HIS:CE1	1:B:194:LEU:HB3	2.54	0.42
1:B:36:LYS:HE2	1:B:45:MET:HG2	2.00	0.42
1:A:239:GLU:HB3	1:B:141:ARG:NH2	2.35	0.42
1:B:115:MET:CE	1:B:181:LEU:HD12	2.49	0.42
1:B:44:ALA:H	1:B:201:GLN:HE22	1.67	0.42
1:B:341:SER:O	1:B:344:ILE:HD13	2.20	0.42
1:A:362:GLY:C	1:A:364:LYS:N	2.72	0.42
1:A:213:CYS:SG	1:A:214:ALA:N	2.93	0.42
1:A:322:THR:O	1:A:325:HIS:HB2	2.19	0.42
1:B:180:VAL:HB	1:B:217:VAL:HG22	2.01	0.41
1:B:73:LYS:HG2	2:B:561:HOH:O	2.19	0.41
1:A:372:CYS:HA	1:A:381:GLU:O	2.21	0.41
1:A:115:MET:HE1	1:A:129:ASP:HB2	2.02	0.41
1:A:33:ILE:HD11	1:A:53:ARG:HE	1.85	0.41
1:A:229:LYS:O	1:A:229:LYS:HG3	2.20	0.41
1:B:361:PRO:HG2	1:B:364:LYS:CB	2.48	0.41
1:A:267:VAL:N	1:A:268:PRO:HD2	2.35	0.41
1:A:22:PRO:HG3	1:A:212:ASP:HB2	2.03	0.41
1:A:107:LEU:CD1	1:A:218:ILE:HG21	2.50	0.41
1:B:276:GLN:HB2	1:B:276:GLN:HE21	1.71	0.41
1:B:35:SER:C	1:B:37:ARG:N	2.73	0.41
1:A:130:HIS:NE2	1:B:247:GLU:CG	2.84	0.41
1:A:362:GLY:O	1:A:364:LYS:N	2.42	0.41
1:A:125:ASN:ND2	1:B:252:PHE:HB3	2.36	0.41
1:B:232:ILE:HD12	1:B:232:ILE:N	2.36	0.41
1:A:149:GLY:HA2	1:B:124:ALA:O	2.21	0.41
1:B:287:GLN:C	1:B:289:GLN:H	2.24	0.41
1:A:44:ALA:O	1:A:48:VAL:HG23	2.21	0.40
1:B:205:ILE:HG22	1:B:205:ILE:O	2.22	0.40
1:B:83:HIS:O	1:B:87:MET:HG2	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:ILE:HD12	1:A:154:GLY:HA2	2.03	0.40
1:B:70:LEU:HD11	1:B:81:GLU:HB3	2.04	0.40
1:A:26:LEU:HB3	1:A:59:GLN:NE2	2.37	0.40
1:A:243:ILE:HD12	1:A:378:ILE:CG2	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	381/465 (82%)	347 (91%)	25 (7%)	9 (2%)	7	13
1	B	381/465 (82%)	340 (89%)	30 (8%)	11 (3%)	6	9
All	All	762/930 (82%)	687 (90%)	55 (7%)	20 (3%)	7	11

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	25	SER
1	A	26	LEU
1	A	306	PRO
1	A	361	PRO
1	B	28	PRO
1	B	36	LYS
1	A	203	THR
1	A	307	GLY
1	B	306	PRO
1	B	23	PRO
1	B	25	SER
1	B	363	GLY
1	B	361	PRO
1	A	378	ILE

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Mol	Chain	Res	Type
1	B	356	MET
1	B	378	ILE
1	B	362	GLY
1	A	363	GLY
1	A	376	PRO
1	B	31	ILE

5.3.2 Protein sidechains ⓘ

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	313/377 (83%)	288 (92%)	25 (8%)	15	28
1	B	313/377 (83%)	284 (91%)	29 (9%)	11	20
All	All	626/754 (83%)	572 (91%)	54 (9%)	13	24

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	LEU
1	A	49	LEU
1	A	88	SER
1	A	99	ARG
1	A	156	LEU
1	A	159	LEU
1	A	169	HIS
1	A	181	LEU
1	A	183	LEU
1	A	194	LEU
1	A	202	GLU
1	A	217	VAL
1	A	263	LEU
1	A	278	THR
1	A	293	SER
1	A	296	LYS
1	A	332	ARG

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Mol	Chain	Res	Type
1	A	344	ILE
1	A	352	ARG
1	A	355	ASP
1	A	359	LEU
1	A	369	VAL
1	A	370	VAL
1	A	385	LEU
1	A	387	ARG
1	B	24	TYR
1	B	29	ASP
1	B	31	ILE
1	B	33	ILE
1	B	34	LEU
1	B	108	VAL
1	B	134	LYS
1	B	146	LEU
1	B	153	SER
1	B	159	LEU
1	B	165	LEU
1	B	181	LEU
1	B	229	LYS
1	B	239	GLU
1	B	252	PHE
1	B	276	GLN
1	B	284	SER
1	B	290	LEU
1	B	310	THR
1	B	326	MET
1	B	332	ARG
1	B	352	ARG
1	B	354	LYS
1	B	359	LEU
1	B	364	LYS
1	B	370	VAL
1	B	380	VAL
1	B	385	LEU
1	B	390	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	HIS

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Mol	Chain	Res	Type
1	A	59	GLN
1	A	72	ASN
1	A	75	ASN
1	A	201	GLN
1	A	287	GLN
1	A	289	GLN
1	A	325	HIS
1	B	39	HIS
1	B	72	ASN
1	B	169	HIS
1	B	201	GLN
1	B	240	ASN
1	B	276	GLN
1	B	325	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 carbohydrates 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

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	383/465 (82%)	-0.21	9 (2%) 64 59	19, 30, 47, 59	0
1	B	383/465 (82%)	-0.14	13 (3%) 49 43	15, 29, 50, 56	0
All	All	766/930 (82%)	-0.17	22 (2%) 55 51	15, 30, 50, 59	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	362	GLY	4.5
1	A	26	LEU	4.5
1	A	227	PRO	3.9
1	A	228	GLY	3.6
1	B	390	ASN	3.5
1	B	228	GLY	3.4
1	B	359	LEU	3.2
1	A	28	PRO	3.1
1	B	28	PRO	2.9
1	B	288	ASP	2.9
1	A	390	ASN	2.8
1	A	203	THR	2.8
1	A	361	PRO	2.7
1	A	25	SER	2.7
1	B	24	TYR	2.5
1	B	361	PRO	2.4
1	B	389	MET	2.3
1	B	357	ASP	2.3
1	B	360	ALA	2.3
1	A	226	ALA	2.2
1	B	287	GLN	2.2
1	B	26	LEU	2.1

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 carbohydrates 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.