



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:01 AM GMT

PDB ID : 2F9Y
Title : The Crystal Structure of The Carboxyltransferase Subunit of ACC from Escherichia coli
Authors : Bilder, P.W.
Deposited on : 2005-12-06
Resolution : 3.20 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

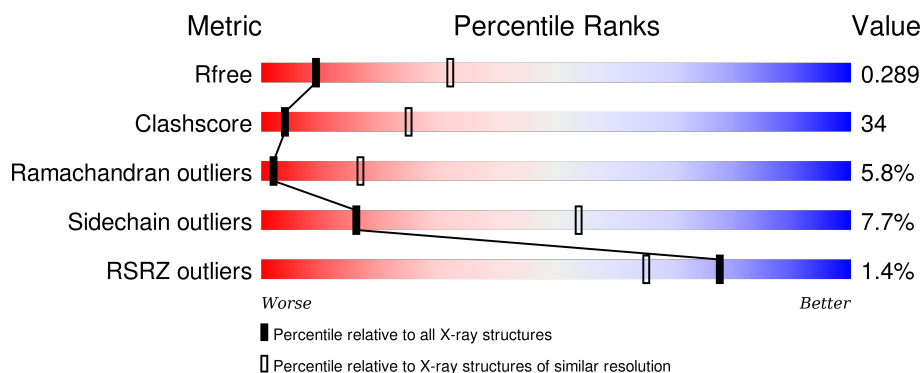
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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}	91344	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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	339	<div> <div> <div></div> <div>44%</div> <div>38%</div> <div>6%</div> <div>11%</div> </div> </div>
2	B	304	<div> <div> <div>2%</div> <div>35%</div> <div>40%</div> <div>9%</div> <div>15%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4266 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 Acetyl-CoA carboxylase, Carboxyltransferase alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	301	Total	C	N	O	S	35	0	0
			2302	1452	411	427	12			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	CLONING ARTIFACT	GB 4902926
A	-18	GLY	-	CLONING ARTIFACT	GB 4902926
A	-17	SER	-	CLONING ARTIFACT	GB 4902926
A	-16	SER	-	CLONING ARTIFACT	GB 4902926
A	-15	HIS	-	EXPRESSION TAG	GB 4902926
A	-14	HIS	-	EXPRESSION TAG	GB 4902926
A	-13	HIS	-	EXPRESSION TAG	GB 4902926
A	-12	HIS	-	EXPRESSION TAG	GB 4902926
A	-11	HIS	-	EXPRESSION TAG	GB 4902926
A	-10	HIS	-	EXPRESSION TAG	GB 4902926
A	-9	SER	-	CLONING ARTIFACT	GB 4902926
A	-8	SER	-	CLONING ARTIFACT	GB 4902926
A	-7	GLY	-	CLONING ARTIFACT	GB 4902926
A	-6	LEU	-	CLONING ARTIFACT	GB 4902926
A	-5	VAL	-	CLONING ARTIFACT	GB 4902926
A	-4	PRO	-	CLONING ARTIFACT	GB 4902926
A	-3	ARG	-	CLONING ARTIFACT	GB 4902926
A	-2	GLY	-	CLONING ARTIFACT	GB 4902926
A	-1	SER	-	CLONING ARTIFACT	GB 4902926
A	0	HIS	-	CLONING ARTIFACT	GB 4902926

- Molecule 2 is a protein called Acetyl-coenzyme A carboxylase carboxyl transferase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	257	Total	C	N	O	S	13	0	0
			1956	1233	343	358	22			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total 1	Zn 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	5	Total 5	O 5	0	0
4	B	2	Total 2	O 2	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	156.73Å 156.73Å 192.82Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.73 – 3.20 30.01 – 3.20	Depositor EDS
% Data completeness (in resolution range)	95.0 (29.73-3.20) 93.4 (30.01-3.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.78 (at 3.18Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.257 , 0.261 0.241 , 0.289	Depositor DCC
R_{free} test set	2170 reflections (9.85%)	DCC
Wilson B-factor (Å ²)	84.0	Xtriage
Anisotropy	0.099	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 96.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 23522 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4266	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.97% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.39	0/2341	0.68	0/3158
2	B	0.35	0/1985	0.64	2/2665 (0.1%)
All	All	0.37	0/4326	0.66	2/5823 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	49	CYS	CA-CB-SG	5.90	124.61	114.00
2	B	285	PRO	N-CA-CB	5.59	110.00	103.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2302	0	2304	140	0
2	B	1956	0	2006	158	0
3	B	1	0	0	0	0
4	A	5	0	0	0	0
4	B	2	0	0	0	0
All	All	4266	0	4310	289	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 34.

All (289) 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:254:ILE:HG22	1:A:256:PRO:HD2	1.37	1.06
1:A:118:HIS:HD2	1:A:162:ASP:H	1.12	0.95
1:A:6:LEU:HG	1:A:7:ASP:H	1.31	0.95
2:B:52:HIS:HE1	2:B:263:ARG:H	1.06	0.93
2:B:52:HIS:HB3	2:B:263:ARG:NH2	1.84	0.93
1:A:8:PHE:HZ	1:A:66:LEU:HD11	1.39	0.88
2:B:106:ASP:HA	2:B:129:PHE:CE1	2.13	0.83
1:A:190:ARG:HD2	2:B:185:ALA:HB1	1.59	0.83
2:B:139:GLY:O	2:B:143:VAL:HG23	1.79	0.83
1:A:49:GLU:O	1:A:50:LEU:HB3	1.78	0.82
1:A:134:MET:HG2	1:A:164:PRO:HG2	1.61	0.81
2:B:64:LEU:HD11	2:B:115:LEU:HD11	1.61	0.81
1:A:209:LEU:HD22	2:B:174:LEU:HG	1.61	0.80
2:B:165:ALA:HB1	2:B:173:SER:HB3	1.65	0.79
1:A:261:LEU:O	1:A:262:LYS:HG2	1.84	0.78
1:A:260:GLU:C	1:A:261:LEU:HD12	2.04	0.78
2:B:223:ALA:O	2:B:247:ARG:HA	1.84	0.78
1:A:118:HIS:CD2	1:A:162:ASP:H	1.98	0.77
2:B:43:LEU:O	2:B:45:VAL:HG23	1.84	0.77
2:B:52:HIS:CE1	2:B:263:ARG:H	1.98	0.76
1:A:244:ALA:N	1:A:245:PRO:HD2	2.00	0.75
2:B:52:HIS:HB3	2:B:263:ARG:HH21	1.47	0.75
2:B:106:ASP:HA	2:B:129:PHE:HE1	1.51	0.74
1:A:219:MET:CE	1:A:264:ILE:HG21	2.20	0.71
1:A:188:MET:HE2	1:A:213:VAL:O	1.91	0.71
2:B:81:ASP:OD1	2:B:84:LYS:N	2.21	0.70
2:B:120:VAL:HG23	2:B:155:PRO:HB2	1.74	0.69
1:A:234:ALA:HB2	1:A:244:ALA:HA	1.74	0.69
2:B:62:HIS:CE1	2:B:70:LEU:HD22	2.27	0.69
2:B:200:ASP:HB2	2:B:223:ALA:HA	1.75	0.69
1:A:157:ILE:HB	1:A:195:VAL:HG22	1.73	0.69
2:B:86:ARG:HG3	2:B:86:ARG:O	1.93	0.68
2:B:106:ASP:OD2	2:B:137:VAL:HB	1.93	0.68
1:A:135:PRO:HG2	1:A:140:TYR:CZ	2.28	0.68
1:A:237:LEU:CD1	2:B:168:GLN:HE22	2.06	0.68
2:B:34:LEU:HD13	2:B:39:LEU:HD11	1.75	0.68
1:A:61:TRP:CG	1:A:245:PRO:HB3	2.29	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:73:LEU:HD11	2:B:148:GLN:HG2	1.74	0.68
2:B:27:CYS:HA	2:B:53:MET:HE1	1.76	0.67
2:B:209:SER:O	2:B:213:LEU:HG	1.95	0.67
2:B:190:ARG:HG3	2:B:192:LEU:HD13	1.76	0.66
1:A:72:ARG:O	1:A:122:ARG:NH2	2.29	0.66
1:A:216:LYS:HA	1:A:265:ASP:OD1	1.96	0.66
2:B:52:HIS:HE1	2:B:263:ARG:N	1.88	0.65
1:A:206:GLY:HA2	1:A:209:LEU:HB3	1.78	0.65
2:B:61:LEU:HD13	2:B:111:MET:HB2	1.79	0.64
1:A:254:ILE:N	1:A:254:ILE:HD12	2.13	0.64
1:A:55:PHE:CD2	1:A:58:LEU:HD11	2.33	0.63
1:A:105:ILE:HG12	1:A:114:MET:HE1	1.79	0.63
1:A:6:LEU:HB3	1:A:9:GLU:HG3	1.81	0.63
1:A:271:PRO:HD2	1:A:275:ALA:HA	1.80	0.63
1:A:243:LYS:HB3	1:A:246:LEU:HD12	1.79	0.63
2:B:54:ARG:NE	2:B:54:ARG:HA	2.14	0.63
2:B:206:VAL:HG12	2:B:211:ALA:HB2	1.81	0.62
2:B:121:VAL:HG21	2:B:149:ALA:HB2	1.81	0.62
2:B:136:SER:HA	2:B:176:GLN:NE2	2.15	0.62
2:B:55:MET:HB2	2:B:263:ARG:NH1	2.15	0.62
2:B:203:MET:HE1	2:B:226:GLY:HA3	1.80	0.62
1:A:294:LEU:O	1:A:294:LEU:HG	2.00	0.61
2:B:112:LYS:HG2	2:B:113:GLY:H	1.65	0.61
1:A:219:MET:CE	1:A:258:LEU:HD12	2.29	0.61
2:B:34:LEU:CD1	2:B:39:LEU:HD11	2.31	0.60
2:B:270:LEU:O	2:B:274:LEU:HB2	2.01	0.60
2:B:35:TYR:HB3	2:B:38:GLU:HG3	1.84	0.60
1:A:199:VAL:HB	1:A:219:MET:HG3	1.82	0.60
1:A:6:LEU:HG	1:A:7:ASP:N	2.10	0.60
2:B:278:MET:CE	2:B:278:MET:HA	2.32	0.60
2:B:85:PHE:HB3	2:B:88:SER:HB3	1.84	0.59
1:A:152:ARG:HH11	1:A:152:ARG:HG3	1.68	0.59
1:A:198:THR:HG23	1:A:218:ASN:HB2	1.84	0.59
2:B:108:LEU:HG	2:B:109:VAL:N	2.17	0.59
2:B:141:ARG:O	2:B:144:ARG:HB3	2.02	0.59
2:B:35:TYR:CE2	2:B:37:ALA:HB3	2.37	0.59
1:A:134:MET:HG2	1:A:164:PRO:CG	2.31	0.58
2:B:198:LEU:HD23	2:B:225:ILE:HG21	1.84	0.58
2:B:190:ARG:HG3	2:B:192:LEU:CD1	2.32	0.58
1:A:193:VAL:HB	1:A:194:PRO:CD	2.33	0.58
1:A:10:GLN:HE21	1:A:10:GLN:HA	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:PHE:CZ	1:A:66:LEU:HD11	2.31	0.57
2:B:225:ILE:HB	2:B:251:LEU:CD2	2.34	0.57
2:B:176:GLN:O	2:B:180:THR:HG23	2.04	0.57
2:B:106:ASP:HB2	2:B:137:VAL:HG11	1.86	0.57
1:A:298:ASP:OD2	1:A:298:ASP:N	2.37	0.57
1:A:43:LEU:HD23	1:A:44:ARG:N	2.20	0.57
1:A:219:MET:HE3	1:A:258:LEU:HD12	1.87	0.56
2:B:69:SER:O	2:B:71:VAL:HG23	2.05	0.56
2:B:234:GLU:C	2:B:236:THR:H	2.09	0.56
2:B:190:ARG:CG	2:B:192:LEU:HD13	2.34	0.56
1:A:219:MET:HE3	1:A:264:ILE:HG21	1.87	0.56
1:A:122:ARG:HH11	1:A:122:ARG:HG3	1.71	0.56
1:A:122:ARG:NH1	1:A:122:ARG:HG3	2.19	0.56
2:B:229:GLY:O	2:B:233:ILE:HG13	2.06	0.56
1:A:237:LEU:HD11	2:B:168:GLN:HE22	1.71	0.55
2:B:29:SER:O	2:B:30:CYS:HB2	2.06	0.55
1:A:256:PRO:HG2	1:A:257:ARG:H	1.70	0.55
1:A:137:PRO:HB2	1:A:141:ARG:NH1	2.21	0.55
1:A:237:LEU:HD13	2:B:168:GLN:HE22	1.71	0.55
1:A:216:LYS:HG2	1:A:293:ASP:HB3	1.87	0.55
2:B:74:GLY:HA2	2:B:76:GLU:OE2	2.07	0.55
2:B:177:MET:HE1	2:B:205:GLY:HA2	1.89	0.55
2:B:52:HIS:CE1	2:B:263:ARG:HB2	2.42	0.54
2:B:107:ALA:HB1	2:B:132:GLY:HA2	1.88	0.54
1:A:92:ASP:O	1:A:93:ARG:HB2	2.07	0.54
2:B:79:PRO:O	2:B:80:LYS:HB3	2.07	0.54
1:A:103:GLY:HA2	1:A:115:ILE:O	2.08	0.54
1:A:219:MET:HE2	1:A:264:ILE:HG21	1.88	0.54
2:B:33:VAL:O	2:B:34:LEU:HD23	2.07	0.53
1:A:60:ALA:HB1	1:A:254:ILE:HD11	1.89	0.53
2:B:221:PRO:HG2	2:B:262:ARG:HG2	1.91	0.53
1:A:295:ALA:C	1:A:297:LEU:H	2.12	0.53
1:A:122:ARG:O	1:A:127:LYS:HE2	2.09	0.53
1:A:10:GLN:HA	1:A:10:GLN:NE2	2.24	0.52
2:B:85:PHE:HD1	2:B:86:ARG:H	1.58	0.52
2:B:112:LYS:HG2	2:B:113:GLY:N	2.25	0.52
2:B:245:PHE:O	2:B:246:GLN:HB2	2.10	0.52
1:A:308:ARG:O	1:A:312:ARG:HG3	2.10	0.52
1:A:71:GLN:C	1:A:122:ARG:HH22	2.14	0.51
2:B:35:TYR:O	2:B:38:GLU:HB2	2.10	0.51
1:A:169:GLY:O	1:A:173:GLU:HG2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:278:MET:HB3	2:B:280:LEU:HD13	1.93	0.51
2:B:225:ILE:HB	2:B:251:LEU:HD21	1.93	0.51
1:A:55:PHE:CE2	1:A:66:LEU:HD21	2.45	0.51
1:A:43:LEU:C	1:A:45:GLU:H	2.14	0.50
2:B:25:THR:HG22	2:B:26:LYS:N	2.26	0.50
1:A:143:ALA:O	1:A:147:MET:HG3	2.10	0.50
1:A:108:LEU:O	1:A:109:ASP:HB2	2.10	0.50
2:B:106:ASP:CG	2:B:166:ARG:HH22	2.15	0.50
1:A:231:GLU:OE1	1:A:231:GLU:N	2.37	0.50
1:A:55:PHE:HE2	1:A:66:LEU:HD21	1.76	0.50
1:A:313:LEU:HA	1:A:316:TYR:CD1	2.47	0.50
1:A:118:HIS:HD2	1:A:162:ASP:N	1.95	0.49
1:A:244:ALA:N	1:A:245:PRO:CD	2.73	0.49
1:A:170:VAL:HG12	2:B:245:PHE:CE1	2.47	0.49
1:A:242:ASP:O	1:A:245:PRO:HD2	2.12	0.49
2:B:206:VAL:O	2:B:211:ALA:HB3	2.11	0.49
1:A:256:PRO:O	1:A:258:LEU:N	2.46	0.49
1:A:293:ASP:C	1:A:295:ALA:H	2.16	0.49
2:B:247:ARG:O	2:B:250:PHE:N	2.44	0.49
2:B:146:VAL:HG13	2:B:187:MET:CE	2.42	0.49
1:A:44:ARG:O	1:A:48:VAL:HG23	2.13	0.49
1:A:15:LEU:HD13	1:A:47:SER:HA	1.94	0.49
1:A:66:LEU:O	1:A:69:HIS:HB2	2.12	0.48
1:A:240:SER:OG	1:A:243:LYS:HB2	2.13	0.48
2:B:136:SER:HB2	2:B:169:GLU:OE1	2.13	0.48
1:A:152:ARG:HD2	1:A:153:PHE:CZ	2.47	0.48
2:B:77:LEU:O	2:B:105:LYS:HG2	2.13	0.48
2:B:28:ASP:H	2:B:53:MET:HE3	1.78	0.48
1:A:261:LEU:N	1:A:261:LEU:HD12	2.27	0.48
2:B:223:ALA:O	2:B:247:ARG:HD3	2.13	0.48
1:A:231:GLU:O	1:A:234:ALA:HB3	2.13	0.48
1:A:222:TYR:N	1:A:270:GLU:OE2	2.45	0.48
1:A:10:GLN:HB3	1:A:11:PRO:CD	2.44	0.48
1:A:317:GLY:HA2	2:B:172:MET:CE	2.44	0.48
1:A:153:PHE:O	1:A:154:LYS:C	2.52	0.48
2:B:156:LEU:HD12	2:B:157:ILE:H	1.79	0.48
1:A:237:LEU:HD13	2:B:168:GLN:NE2	2.29	0.47
1:A:8:PHE:C	1:A:8:PHE:CD1	2.88	0.47
2:B:106:ASP:OD2	2:B:138:VAL:HG23	2.13	0.47
1:A:222:TYR:HB2	1:A:270:GLU:OE2	2.13	0.47
2:B:172:MET:O	2:B:176:GLN:HG2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:268:ILE:CD1	1:A:286:LEU:HD12	2.45	0.47
1:A:253:ILE:HG12	1:A:254:ILE:HD12	1.96	0.47
2:B:59:ASN:O	2:B:62:HIS:HB2	2.14	0.47
2:B:170:ALA:O	2:B:172:MET:N	2.47	0.47
2:B:129:PHE:HD2	2:B:130:MET:HG3	1.80	0.47
1:A:205:SER:OG	1:A:206:GLY:N	2.47	0.47
2:B:203:MET:HE1	2:B:226:GLY:CA	2.43	0.47
1:A:10:GLN:O	1:A:13:ALA:N	2.47	0.47
2:B:161:ALA:HA	2:B:201:PRO:O	2.14	0.47
2:B:247:ARG:O	2:B:248:SER:C	2.53	0.47
2:B:86:ARG:O	2:B:87:ASP:HB2	2.15	0.47
2:B:32:GLN:HE21	2:B:32:GLN:HA	1.80	0.47
2:B:49:CYS:C	2:B:51:HIS:H	2.19	0.47
1:A:76:LEU:HD21	1:A:102:VAL:HG11	1.96	0.47
1:A:281:ALA:O	1:A:284:ALA:HB3	2.14	0.47
2:B:156:LEU:HD12	2:B:157:ILE:N	2.29	0.47
1:A:254:ILE:HG22	1:A:255:ARG:N	2.29	0.46
2:B:105:LYS:O	2:B:129:PHE:HE1	1.99	0.46
1:A:303:GLU:O	1:A:307:ASN:ND2	2.48	0.46
2:B:105:LYS:HB2	2:B:141:ARG:NH2	2.29	0.46
2:B:56:THR:HG23	2:B:59:ASN:HB2	1.98	0.46
1:A:10:GLN:HE21	1:A:10:GLN:CA	2.27	0.46
2:B:38:GLU:O	2:B:40:GLU:N	2.40	0.46
2:B:236:THR:O	2:B:238:ARG:HG3	2.15	0.46
1:A:116:ILE:O	1:A:159:THR:HA	2.14	0.46
2:B:76:GLU:HG2	2:B:76:GLU:O	2.15	0.46
1:A:220:LEU:HD23	1:A:268:ILE:HB	1.97	0.46
2:B:219:ALA:O	2:B:260:ILE:HG23	2.16	0.46
2:B:62:HIS:ND1	2:B:70:LEU:HD22	2.30	0.46
1:A:317:GLY:HA2	2:B:172:MET:HE2	1.97	0.46
2:B:278:MET:HA	2:B:278:MET:HE3	1.98	0.46
2:B:43:LEU:O	2:B:45:VAL:N	2.49	0.45
1:A:152:ARG:HH11	1:A:152:ARG:CG	2.29	0.45
1:A:219:MET:HE2	1:A:219:MET:HB2	1.84	0.45
1:A:221:GLN:HB2	1:A:269:PRO:HA	1.99	0.45
1:A:259:LYS:C	1:A:261:LEU:H	2.18	0.45
2:B:106:ASP:CG	2:B:137:VAL:HB	2.35	0.45
1:A:171:GLY:O	1:A:175:ARG:HB2	2.17	0.45
2:B:32:GLN:O	2:B:34:LEU:HG	2.17	0.45
1:A:50:LEU:C	1:A:52:ARG:N	2.69	0.45
2:B:135:GLY:O	2:B:137:VAL:N	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:184:LEU:O	2:B:187:MET:HB3	2.17	0.45
1:A:152:ARG:CZ	1:A:152:ARG:HB2	2.47	0.45
2:B:73:LEU:CD1	2:B:148:GLN:HG2	2.43	0.45
2:B:121:VAL:CG2	2:B:149:ALA:HB2	2.47	0.44
2:B:198:LEU:N	2:B:198:LEU:HD12	2.32	0.44
2:B:182:ALA:O	2:B:185:ALA:HB3	2.17	0.44
1:A:293:ASP:O	1:A:295:ALA:N	2.50	0.44
1:A:273:GLY:O	1:A:277:ARG:HG3	2.17	0.44
1:A:216:LYS:HG2	1:A:293:ASP:CG	2.38	0.44
2:B:192:LEU:HD12	2:B:192:LEU:N	2.31	0.44
1:A:39:GLU:C	1:A:41:HIS:N	2.71	0.44
1:A:132:PHE:N	1:A:132:PHE:CD1	2.85	0.44
2:B:61:LEU:HD22	2:B:111:MET:HB3	1.98	0.44
2:B:207:SER:O	2:B:212:MET:HB2	2.18	0.44
2:B:192:LEU:CD1	2:B:192:LEU:N	2.80	0.44
2:B:206:VAL:HG13	2:B:210:PHE:CZ	2.52	0.44
2:B:146:VAL:HG13	2:B:187:MET:HE2	1.99	0.44
2:B:177:MET:CE	2:B:205:GLY:HA2	2.47	0.44
2:B:150:LEU:O	2:B:190:ARG:NH2	2.51	0.43
2:B:43:LEU:O	2:B:44:GLU:C	2.56	0.43
2:B:237:VAL:O	2:B:239:GLU:HG3	2.17	0.43
2:B:194:TYR:CD2	2:B:214:GLY:HA2	2.53	0.43
1:A:270:GLU:OE1	1:A:275:ALA:HB2	2.18	0.43
2:B:107:ALA:O	2:B:125:PHE:HA	2.18	0.43
1:A:216:LYS:HG2	1:A:293:ASP:CB	2.49	0.43
2:B:51:HIS:CD2	2:B:53:MET:HG2	2.54	0.43
1:A:288:ALA:O	1:A:291:LEU:HB2	2.18	0.43
2:B:106:ASP:HB2	2:B:137:VAL:CG1	2.49	0.43
1:A:261:LEU:O	1:A:263:LEU:N	2.51	0.43
2:B:274:LEU:HD12	2:B:274:LEU:HA	1.82	0.43
2:B:249:GLU:H	2:B:249:GLU:CD	2.23	0.43
1:A:259:LYS:O	1:A:259:LYS:HD2	2.18	0.43
2:B:184:LEU:HD12	2:B:184:LEU:HA	1.92	0.43
1:A:40:VAL:O	1:A:42:ARG:N	2.52	0.43
2:B:106:ASP:OD2	2:B:166:ARG:NH2	2.52	0.42
2:B:129:PHE:CD2	2:B:130:MET:HG3	2.54	0.42
2:B:121:VAL:HG23	2:B:154:CYS:SG	2.59	0.42
1:A:254:ILE:CD1	1:A:254:ILE:N	2.81	0.42
2:B:244:GLY:O	2:B:245:PHE:C	2.58	0.42
2:B:65:LEU:O	2:B:66:ASP:O	2.37	0.42
2:B:198:LEU:HD23	2:B:225:ILE:CG2	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:206:GLY:HA2	1:A:209:LEU:CB	2.48	0.42
2:B:34:LEU:HD11	2:B:47:PRO:HD2	2.02	0.42
1:A:256:PRO:O	1:A:257:ARG:C	2.58	0.42
2:B:110:VAL:HG21	2:B:141:ARG:O	2.20	0.42
2:B:129:PHE:O	2:B:130:MET:HB2	2.20	0.42
2:B:153:ASN:HA	2:B:192:LEU:HD21	2.01	0.42
1:A:102:VAL:HG23	1:A:119:GLN:HB2	2.01	0.42
1:A:232:GLY:O	1:A:233:CYS:C	2.58	0.42
2:B:146:VAL:HA	2:B:187:MET:HE1	2.00	0.42
1:A:122:ARG:HH11	1:A:122:ARG:CG	2.33	0.42
1:A:77:ASP:O	1:A:80:ARG:HG2	2.19	0.42
1:A:50:LEU:C	1:A:52:ARG:H	2.21	0.42
1:A:147:MET:HB3	1:A:157:ILE:HD13	2.01	0.42
1:A:219:MET:HG2	1:A:223:SER:OG	2.18	0.42
1:A:114:MET:HB2	1:A:155:MET:HE2	2.01	0.42
2:B:234:GLU:O	2:B:236:THR:N	2.50	0.42
1:A:12:ILE:HD11	1:A:54:ILE:HD12	2.00	0.42
2:B:42:ASN:HA	2:B:42:ASN:HD22	1.67	0.42
2:B:61:LEU:HD11	2:B:109:VAL:O	2.20	0.41
2:B:237:VAL:HG12	2:B:237:VAL:O	2.20	0.41
2:B:114:THR:HA	2:B:118:MET:O	2.20	0.41
1:A:256:PRO:C	1:A:258:LEU:N	2.73	0.41
1:A:156:PRO:HB3	1:A:294:LEU:HD11	2.01	0.41
2:B:157:ILE:HD11	2:B:274:LEU:CD2	2.50	0.41
2:B:198:LEU:CD2	2:B:225:ILE:HG21	2.50	0.41
1:A:39:GLU:C	1:A:41:HIS:H	2.23	0.41
2:B:222:LYS:HA	2:B:249:GLU:OE1	2.21	0.41
2:B:150:LEU:HD23	2:B:190:ARG:NH1	2.34	0.41
2:B:156:LEU:CD2	2:B:187:MET:HE3	2.50	0.41
1:A:40:VAL:C	1:A:42:ARG:H	2.24	0.41
2:B:106:ASP:OD1	2:B:166:ARG:NH2	2.52	0.41
2:B:224:LEU:HD13	2:B:224:LEU:C	2.40	0.41
1:A:5:PHE:HD1	1:A:6:LEU:O	2.04	0.41
2:B:55:MET:HB2	2:B:263:ARG:HH12	1.85	0.41
2:B:200:ASP:HA	2:B:201:PRO:HA	1.75	0.41
1:A:102:VAL:CG2	1:A:119:GLN:HB2	2.51	0.41
2:B:106:ASP:CG	2:B:166:ARG:NH2	2.74	0.41
1:A:293:ASP:HA	1:A:296:ASP:OD1	2.20	0.41
2:B:234:GLU:C	2:B:236:THR:N	2.74	0.41
1:A:59:GLY:O	1:A:60:ALA:C	2.58	0.41
2:B:176:GLN:HA	2:B:176:GLN:OE1	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238:TRP:O	1:A:240:SER:N	2.54	0.40
1:A:313:LEU:HA	1:A:316:TYR:CE1	2.55	0.40
1:A:40:VAL:C	1:A:42:ARG:N	2.74	0.40
1:A:62:GLN:HA	1:A:65:GLN:OE1	2.21	0.40
2:B:74:GLY:O	2:B:77:LEU:HD12	2.22	0.40
2:B:107:ALA:CB	2:B:132:GLY:HA2	2.51	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	297/339 (88%)	237 (80%)	45 (15%)	15 (5%)	2	20
2	B	253/304 (83%)	194 (77%)	42 (17%)	17 (7%)	1	12
All	All	550/643 (86%)	431 (78%)	87 (16%)	32 (6%)	2	17

All (32) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	125	LYS
1	A	131	ASN
2	B	44	GLU
2	B	66	ASP
2	B	80	LYS
2	B	136	SER
1	A	5	PHE
1	A	7	ASP
1	A	8	PHE
1	A	205	SER
1	A	294	LEU
2	B	30	CYS

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Mol	Chain	Res	Type
2	B	43	LEU
2	B	88	SER
2	B	107	ALA
2	B	171	LEU
2	B	235	GLN
1	A	38	GLU
1	A	41	HIS
1	A	154	LYS
1	A	239	LYS
1	A	257	ARG
1	A	262	LYS
2	B	28	ASP
2	B	231	ARG
2	B	245	PHE
1	A	235	SER
2	B	79	PRO
2	B	170	ALA
2	B	242	PRO
2	B	205	GLY
1	A	271	PRO

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	230/273 (84%)	212 (92%)	18 (8%)	16	53
2	B	209/251 (83%)	193 (92%)	16 (8%)	16	54
All	All	439/524 (84%)	405 (92%)	34 (8%)	16	54

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	LEU
1	A	40	VAL
1	A	52	ARG

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Mol	Chain	Res	Type
1	A	84	ASP
1	A	122	ARG
1	A	125	LYS
1	A	148	GLN
1	A	170	VAL
1	A	188	MET
1	A	198	THR
1	A	218	ASN
1	A	225	TYR
1	A	238	TRP
1	A	257	ARG
1	A	296	ASP
1	A	298	ASP
1	A	305	LEU
1	A	318	TYR
2	B	42	ASN
2	B	55	MET
2	B	56	THR
2	B	76	GLU
2	B	86	ARG
2	B	87	ASP
2	B	106	ASP
2	B	129	PHE
2	B	166	ARG
2	B	184	LEU
2	B	201	PRO
2	B	203	MET
2	B	216	LEU
2	B	232	VAL
2	B	274	LEU
2	B	278	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	118	HIS
1	A	148	GLN
1	A	218	ASN
1	A	221	GLN
2	B	32	GLN
2	B	42	ASN

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Mol	Chain	Res	Type
2	B	52	HIS
2	B	99	GLN
2	B	153	ASN
2	B	168	GLN
2	B	217	ASN
2	B	279	ASN

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](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	301/339 (88%)	-0.41	3 (0%) 84 75	25, 61, 138, 174	27 (8%)
2	B	257/304 (84%)	-0.18	5 (1%) 70 55	32, 83, 160, 201	9 (3%)
All	All	558/643 (86%)	-0.31	8 (1%) 78 65	25, 73, 151, 201	36 (6%)

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	24	TRP	5.5
2	B	35	TYR	2.9
1	A	299	VAL	2.6
2	B	238	ARG	2.5
1	A	205	SER	2.4
2	B	100	LYS	2.4
1	A	226	SER	2.2
2	B	23	VAL	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 carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains.

The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	ZN	B	305	1/1	0.98	0.02	-2.21	81,81,81,81	0

6.5 Other polymers [i](#)

There are no such residues in this entry.