



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:52 PM GMT

PDB ID : 4GGO
Title : Crystal Structure of Trans-2-Enoyl-CoA Reductase from *Treponema denticola*
Authors : Bond-Watts, B.B.; Weeks, A.M.; Chang, M.C.Y.
Deposited on : 2012-08-06
Resolution : 2.00 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

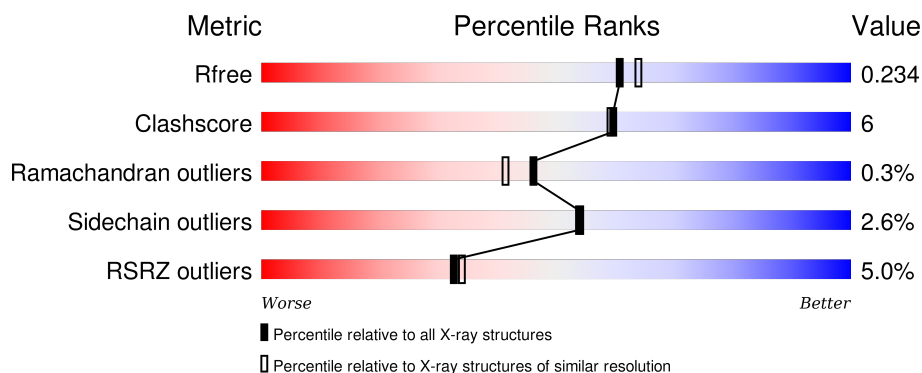
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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	401	<div> <div>2%</div> <div>88%</div> <div>11%</div> <div>.</div> </div>
1	B	401	<div> <div>4%</div> <div>84%</div> <div>14%</div> <div>..</div> </div>
1	C	401	<div> <div>3%</div> <div>89%</div> <div>9%</div> <div>.</div> </div>
1	D	401	<div> <div>10%</div> <div>80%</div> <div>16%</div> <div>..</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12957 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 Trans-2-enoyl-CoA reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	397	Total	C	N	O	S	0	0	0
			3079	1946	525	597	11			
1	B	397	Total	C	N	O	S	0	0	0
			3079	1946	525	597	11			
1	C	397	Total	C	N	O	S	0	0	0
			3079	1946	525	597	11			
1	D	397	Total	C	N	O	S	0	0	0
			3079	1946	525	597	11			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	EXPRESSION TAG	UNP Q73Q47
A	-2	THR	-	EXPRESSION TAG	UNP Q73Q47
A	-1	GLY	-	EXPRESSION TAG	UNP Q73Q47
A	0	ALA	-	EXPRESSION TAG	UNP Q73Q47
B	-3	GLY	-	EXPRESSION TAG	UNP Q73Q47
B	-2	THR	-	EXPRESSION TAG	UNP Q73Q47
B	-1	GLY	-	EXPRESSION TAG	UNP Q73Q47
B	0	ALA	-	EXPRESSION TAG	UNP Q73Q47
C	-3	GLY	-	EXPRESSION TAG	UNP Q73Q47
C	-2	THR	-	EXPRESSION TAG	UNP Q73Q47
C	-1	GLY	-	EXPRESSION TAG	UNP Q73Q47
C	0	ALA	-	EXPRESSION TAG	UNP Q73Q47
D	-3	GLY	-	EXPRESSION TAG	UNP Q73Q47
D	-2	THR	-	EXPRESSION TAG	UNP Q73Q47
D	-1	GLY	-	EXPRESSION TAG	UNP Q73Q47
D	0	ALA	-	EXPRESSION TAG	UNP Q73Q47

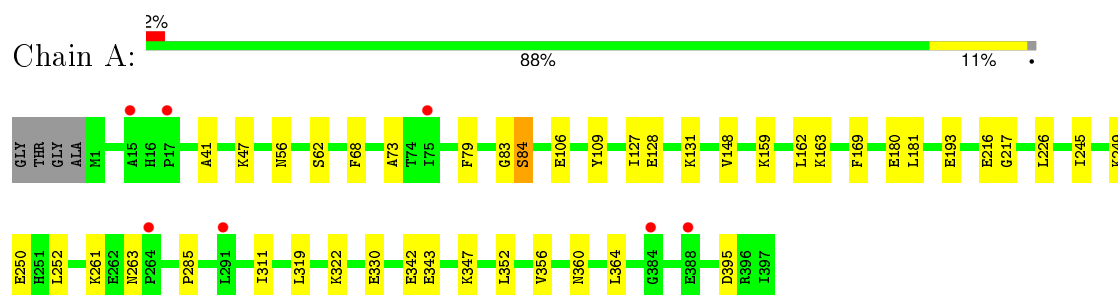
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	194	Total 194	O 194	0	0
2	B	143	Total 143	O 143	0	0
2	C	193	Total 193	O 193	0	0
2	D	111	Total 111	O 111	0	0

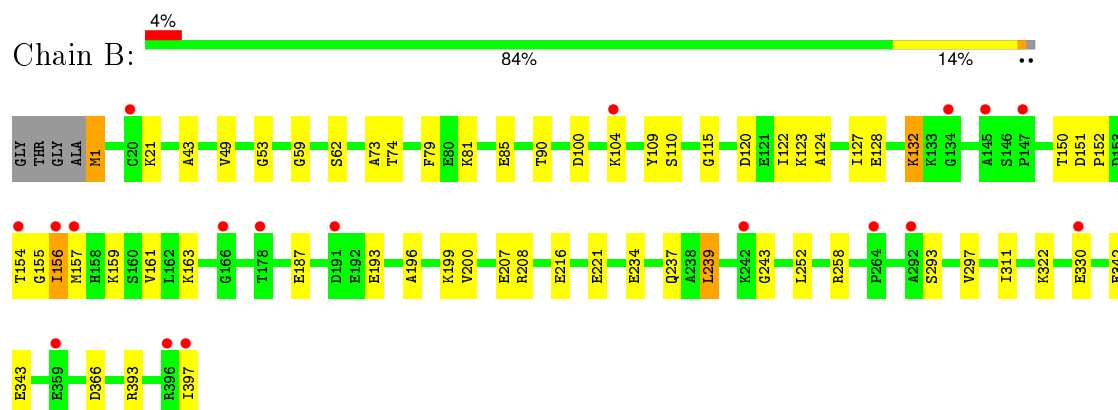
3 Residue-property plots [i](#)

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.

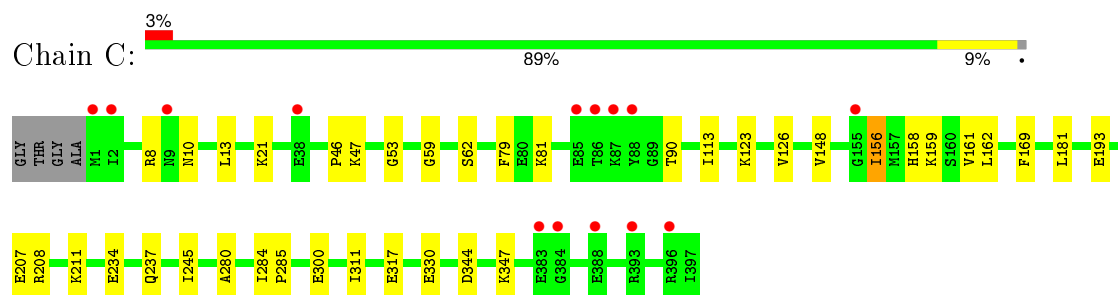
- Molecule 1: Trans-2-enoyl-CoA reductase



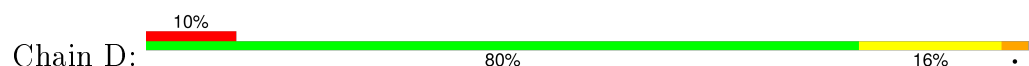
- Molecule 1: Trans-2-enoyl-CoA reductase

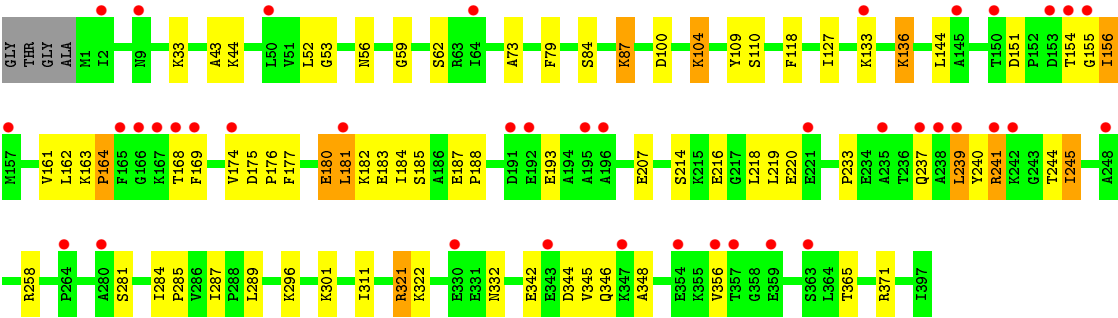


- Molecule 1: Trans-2-enoyl-CoA reductase



- Molecule 1: Trans-2-enoyl-CoA reductase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	62.05Å 87.05Å 91.46Å 106.14° 109.72° 98.52°	Depositor
Resolution (Å)	19.68 – 2.00 19.68 – 2.00	Depositor EDS
% Data completeness (in resolution range)	94.0 (19.68-2.00) 82.5 (19.68-2.00)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 2.01Å)	Xtriage
Refinement program	PHENIX AUTOREFINE	Depositor
R, R_{free}	0.206 , 0.235 0.204 , 0.234	Depositor DCC
R_{free} test set	5290 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	31.0	Xtriage
Anisotropy	0.638	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 52.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 105852 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12957	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.15% 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

5.1 Standard geometry

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	0.22	0/3134	0.39	0/4229
1	B	0.23	0/3134	0.40	0/4229
1	C	0.22	0/3134	0.38	0/4229
1	D	0.25	0/3134	0.44	1/4229 (0.0%)
All	All	0.23	0/12536	0.40	1/16916 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	181	LEU	N-CA-C	5.11	124.81	111.00

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	3079	0	3080	28	0
1	B	3079	0	3080	37	0
1	C	3079	0	3080	24	0
1	D	3079	0	3080	52	0
2	A	194	0	0	9	0
2	B	143	0	0	8	0
2	C	193	0	0	6	1
2	D	111	0	0	8	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	12957	0	12320	140	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (140) 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:150:THR:O	2:B:424:HOH:O	1.93	0.87
1:B:152:PRO:O	2:B:444:HOH:O	1.94	0.83
1:C:46:PRO:O	2:C:491:HOH:O	1.96	0.82
1:A:249:LYS:NZ	2:A:516:HOH:O	2.16	0.79
1:B:330:GLU:OE2	2:B:542:HOH:O	2.01	0.78
1:D:154:THR:OG1	2:D:504:HOH:O	2.00	0.78
1:B:120:ASP:O	2:B:536:HOH:O	2.03	0.77
1:B:1:MET:N	2:B:517:HOH:O	2.17	0.76
1:D:180:GLU:C	1:D:181:LEU:HD12	2.07	0.75
1:D:118:PHE:O	2:D:493:HOH:O	2.05	0.74
1:D:174:VAL:HG22	1:D:181:LEU:HG	1.71	0.73
1:D:181:LEU:N	1:D:181:LEU:HD12	2.05	0.71
1:A:128:GLU:OE1	2:A:510:HOH:O	2.09	0.70
1:A:342:GLU:OE2	2:A:404:HOH:O	2.10	0.69
1:C:8:ARG:HB3	1:C:13:LEU:HD11	1.76	0.68
1:B:124:ALA:N	2:B:536:HOH:O	2.21	0.67
1:A:250:GLU:OE2	2:A:507:HOH:O	2.13	0.65
1:D:348:ALA:O	2:D:478:HOH:O	2.14	0.65
1:C:10:ASN:O	2:C:458:HOH:O	2.14	0.64
1:B:150:THR:HG22	1:B:157:MET:HG2	1.80	0.64
1:C:237:GLN:NE2	2:C:586:HOH:O	2.30	0.64
1:D:301:LYS:NZ	1:D:342:GLU:OE1	2.29	0.63
1:A:250:GLU:OE1	2:A:445:HOH:O	2.16	0.62
1:D:240:TYR:CD2	1:D:245:ILE:HG21	2.35	0.62
1:B:104:LYS:NZ	1:B:110:SER:O	2.33	0.61
1:A:181:LEU:HD11	1:A:356:VAL:HG12	1.83	0.61
1:C:47:LYS:NZ	2:C:578:HOH:O	2.28	0.61
1:B:128:GLU:OE1	1:B:132:LYS:NZ	2.35	0.60
1:D:207:GLU:OE1	1:D:258:ARG:NH2	2.33	0.60
1:C:344:ASP:HA	1:C:347:LYS:HE2	1.85	0.58
1:B:151:ASP:OD2	1:B:199:LYS:NZ	2.36	0.58
1:C:148:VAL:HG12	1:C:159:LYS:HG2	1.86	0.58
1:B:239:LEU:HA	1:B:243:GLY:HA3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:344:ASP:OD1	1:D:345:VAL:N	2.37	0.58
1:D:33:LYS:NZ	2:D:464:HOH:O	2.37	0.56
1:D:296:LYS:HG3	1:D:365:THR:HA	1.86	0.56
1:A:360:ASN:HB2	1:A:364:LEU:HD13	1.88	0.56
1:B:342:GLU:OE2	2:B:429:HOH:O	2.17	0.56
1:D:163:LYS:NZ	1:D:193:GLU:OE1	2.38	0.56
1:D:289:LEU:HD11	1:D:356:VAL:HG11	1.89	0.55
1:C:234:GLU:HA	1:C:237:GLN:HG3	1.88	0.55
1:B:62:SER:HA	1:B:311:ILE:HG21	1.88	0.55
1:D:154:THR:O	1:D:156:ILE:N	2.36	0.54
1:C:62:SER:HA	1:C:311:ILE:HG21	1.88	0.54
1:C:300:GLU:O	1:D:321:ARG:NH2	2.39	0.54
1:B:124:ALA:HA	1:B:127:ILE:HB	1.89	0.54
1:A:131:LYS:NZ	1:A:216:GLU:OE1	2.41	0.54
1:D:44:LYS:NZ	2:D:503:HOH:O	2.07	0.54
1:A:343:GLU:HG2	1:A:347:LYS:HE2	1.89	0.54
1:A:62:SER:HA	1:A:311:ILE:HG21	1.90	0.53
1:D:214:SER:HB2	1:D:219:LEU:HD12	1.90	0.53
1:B:127:ILE:HD13	1:B:216:GLU:HG3	1.90	0.53
1:B:85:GLU:HA	1:B:393:ARG:HG3	1.91	0.52
1:A:47:LYS:HE2	2:A:475:HOH:O	2.09	0.52
1:D:62:SER:HA	1:D:311:ILE:HG21	1.91	0.52
1:D:240:TYR:HD2	1:D:245:ILE:HG21	1.71	0.52
1:D:181:LEU:CD1	1:D:181:LEU:N	2.72	0.52
1:A:56:ASN:ND2	2:A:552:HOH:O	2.43	0.51
1:D:180:GLU:CA	1:D:181:LEU:HD12	2.41	0.51
1:B:366:ASP:OD1	2:B:428:HOH:O	2.19	0.50
1:B:100:ASP:O	1:B:104:LYS:HG2	2.11	0.50
1:B:163:LYS:NZ	1:B:193:GLU:OE2	2.39	0.49
1:C:207:GLU:HG2	1:C:211:LYS:HE2	1.92	0.49
1:B:1:MET:N	1:B:1:MET:SD	2.82	0.49
1:D:281:SER:HA	1:D:284:ILE:HD11	1.93	0.49
1:A:330:GLU:H	1:A:330:GLU:CD	2.15	0.49
1:C:21:LYS:NZ	2:C:512:HOH:O	2.46	0.49
1:B:21:LYS:HD3	1:B:397:ILE:HG21	1.93	0.49
1:D:136:LYS:HG2	1:D:220:GLU:HB2	1.93	0.48
1:A:127:ILE:HD13	1:A:216:GLU:HG3	1.95	0.48
1:D:233:PRO:HB3	1:D:346:GLN:OE1	2.14	0.48
1:A:162:LEU:HD22	1:A:245:ILE:HD11	1.96	0.47
1:A:226:LEU:HD11	1:A:319:LEU:HD21	1.95	0.47
1:D:184:ILE:HG23	1:D:285:PRO:HG3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:240:TYR:O	1:D:241:ARG:HB3	2.15	0.47
1:B:73:ALA:HB1	1:B:109:TYR:HD2	1.78	0.46
1:B:151:ASP:HB3	1:B:154:THR:O	2.15	0.46
1:B:330:GLU:H	1:B:330:GLU:CD	2.19	0.46
1:B:154:THR:O	1:B:156:ILE:N	2.39	0.46
1:D:164:PRO:HG2	1:D:169:PHE:HB3	1.97	0.46
1:A:73:ALA:HB1	1:A:109:TYR:HD2	1.79	0.46
1:D:56:ASN:ND2	2:D:438:HOH:O	2.28	0.46
1:D:73:ALA:HB1	1:D:109:TYR:HD2	1.81	0.46
1:B:43:ALA:HB2	1:B:322:LYS:HD3	1.97	0.46
1:D:163:LYS:HA	1:D:164:PRO:HD3	1.72	0.46
1:D:104:LYS:NZ	1:D:110:SER:O	2.49	0.46
1:C:113:ILE:HD12	1:C:126:VAL:HG22	1.98	0.45
1:D:100:ASP:O	1:D:104:LYS:HE2	2.16	0.45
1:B:293:SER:O	1:B:297:VAL:HG12	2.15	0.45
1:D:168:THR:HA	1:D:188:PRO:HD3	1.98	0.45
1:B:207:GLU:OE1	1:B:258:ARG:NH2	2.42	0.45
1:D:175:ASP:HA	1:D:176:PRO:HD3	1.82	0.45
1:B:234:GLU:HA	1:B:237:GLN:HG3	2.00	0.44
1:A:68:PHE:HE1	1:A:106:GLU:HG3	1.82	0.44
1:D:177:PHE:HA	1:D:371:ARG:HD3	1.99	0.44
1:A:180:GLU:OE2	2:A:514:HOH:O	2.21	0.43
1:C:123:LYS:HE2	1:C:208:ARG:HG2	2.00	0.43
1:D:43:ALA:HB2	1:D:322:LYS:HD3	2.00	0.43
1:B:123:LYS:HE2	1:B:208:ARG:HG2	1.99	0.43
1:D:182:LYS:NZ	2:D:496:HOH:O	2.52	0.43
1:D:84:SER:OG	1:D:87:LYS:HB2	2.18	0.43
1:B:115:GLY:HA3	1:B:122:ILE:HD11	2.01	0.43
1:C:161:VAL:HG21	1:C:193:GLU:HB3	1.99	0.43
1:C:317:GLU:OE1	2:C:572:HOH:O	2.21	0.43
1:D:284:ILE:HA	1:D:285:PRO:HD3	1.91	0.43
1:D:127:ILE:HG23	1:D:218:LEU:HD11	2.01	0.43
1:D:154:THR:O	1:D:154:THR:OG1	2.27	0.42
1:D:127:ILE:HD13	1:D:216:GLU:HG3	2.00	0.42
1:D:284:ILE:HD12	1:D:287:ILE:HG13	2.01	0.42
1:D:163:LYS:H	1:D:244:THR:HG21	1.84	0.42
1:C:162:LEU:HD22	1:C:245:ILE:HD11	2.00	0.42
1:C:330:GLU:CD	1:C:330:GLU:H	2.22	0.42
1:D:53:GLY:O	1:D:59:GLY:HA3	2.19	0.42
1:A:163:LYS:NZ	1:A:193:GLU:OE2	2.40	0.42
1:D:163:LYS:HB3	1:D:187:GLU:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:73:ALA:HB1	1:D:109:TYR:CD2	2.54	0.42
1:A:395:ASP:N	1:A:395:ASP:OD1	2.50	0.42
1:B:85:GLU:HG3	1:B:393:ARG:HD3	2.01	0.42
1:B:161:VAL:HG21	1:B:193:GLU:HB3	2.01	0.42
1:C:81:LYS:HD2	1:C:81:LYS:N	2.35	0.42
1:A:352:LEU:O	1:A:356:VAL:HG23	2.19	0.41
1:B:49:VAL:HB	1:B:74:THR:HG22	2.02	0.41
1:D:151:ASP:HB3	1:D:154:THR:O	2.20	0.41
1:D:239:LEU:HA	1:D:239:LEU:HD12	1.63	0.41
1:A:68:PHE:CE1	1:A:106:GLU:HG3	2.56	0.41
1:C:53:GLY:O	1:C:59:GLY:HA3	2.21	0.41
1:A:41:ALA:O	1:A:322:LYS:NZ	2.46	0.41
1:B:53:GLY:O	1:B:59:GLY:HA3	2.20	0.41
1:C:81:LYS:HB2	1:C:90:THR:HG23	2.03	0.41
1:B:196:ALA:O	1:B:200:VAL:HG23	2.21	0.41
1:C:169:PHE:CE2	1:C:285:PRO:HB2	2.56	0.41
1:D:237:GLN:NE2	2:D:448:HOH:O	2.53	0.41
1:A:83:GLY:O	1:A:84:SER:HB3	2.21	0.41
1:D:184:ILE:HG12	1:D:185:SER:N	2.36	0.40
1:A:169:PHE:CE2	1:A:285:PRO:HB2	2.56	0.40
1:A:217:GLY:HA2	2:A:594:HOH:O	2.21	0.40
1:A:148:VAL:HG12	1:A:159:LYS:HG2	2.03	0.40
1:B:81:LYS:HB2	1:B:90:THR:HG23	2.02	0.40
1:C:156:ILE:HD11	1:C:158:HIS:NE2	2.36	0.40
1:C:280:ALA:O	1:C:284:ILE:HG13	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:554:HOH:O	2:D:484:HOH:O[1_454]	2.06	0.14

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	395/401 (98%)	386 (98%)	9 (2%)	0	100	100
1	B	395/401 (98%)	382 (97%)	12 (3%)	1 (0%)	46	41
1	C	395/401 (98%)	388 (98%)	7 (2%)	0	100	100
1	D	395/401 (98%)	373 (94%)	19 (5%)	3 (1%)	24	15
All	All	1580/1604 (98%)	1529 (97%)	47 (3%)	4 (0%)	46	41

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	241	ARG
1	D	155	GLY
1	D	164	PRO
1	B	155	GLY

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	323/324 (100%)	318 (98%)	5 (2%)	72	75
1	B	323/324 (100%)	313 (97%)	10 (3%)	47	46
1	C	323/324 (100%)	320 (99%)	3 (1%)	84	88
1	D	323/324 (100%)	307 (95%)	16 (5%)	30	24
All	All	1292/1296 (100%)	1258 (97%)	34 (3%)	54	54

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

Mol	Chain	Res	Type
1	A	79	PHE
1	A	84	SER
1	A	252	LEU
1	A	261	LYS

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Mol	Chain	Res	Type
1	A	263	ASN
1	B	1	MET
1	B	79	PHE
1	B	132	LYS
1	B	156	ILE
1	B	159	LYS
1	B	187	GLU
1	B	221	GLU
1	B	239	LEU
1	B	252	LEU
1	B	343	GLU
1	C	79	PHE
1	C	156	ILE
1	C	181	LEU
1	D	52	LEU
1	D	79	PHE
1	D	87	LYS
1	D	104	LYS
1	D	133	LYS
1	D	136	LYS
1	D	144	LEU
1	D	156	ILE
1	D	161	VAL
1	D	162	LEU
1	D	180	GLU
1	D	183	GLU
1	D	239	LEU
1	D	245	ILE
1	D	321	ARG
1	D	332	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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	397/401 (99%)	0.04	7 (1%) 71 72	17, 32, 51, 70	0
1	B	397/401 (99%)	0.22	18 (4%) 37 38	16, 36, 59, 77	0
1	C	397/401 (99%)	0.10	14 (3%) 48 49	19, 34, 56, 80	0
1	D	397/401 (99%)	0.51	40 (10%) 9 10	19, 40, 70, 85	0
All	All	1588/1604 (99%)	0.22	79 (4%) 32 34	16, 35, 61, 85	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	166	GLY	6.6
1	D	166	GLY	6.6
1	D	165	PHE	5.3
1	D	167	LYS	4.6
1	D	154	THR	4.5
1	D	155	GLY	4.5
1	D	264	PRO	4.1
1	C	2	ILE	4.0
1	D	191	ASP	3.7
1	B	264	PRO	3.6
1	A	388	GLU	3.5
1	D	357	THR	3.5
1	D	9	ASN	3.5
1	B	191	ASP	3.4
1	D	169	PHE	3.2
1	A	15	ALA	3.2
1	B	145	ALA	3.2
1	D	237	GLN	3.2
1	B	396	ARG	3.1
1	D	192	GLU	3.1
1	B	359	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	241	ARG	3.0
1	D	196	ALA	3.0
1	D	248	ALA	3.0
1	D	242	LYS	3.0
1	C	1	MET	3.0
1	D	330	GLU	2.9
1	D	168	THR	2.9
1	C	86	THR	2.9
1	A	264	PRO	2.9
1	D	145	ALA	2.8
1	D	195	ALA	2.8
1	D	157	MET	2.8
1	D	354	GLU	2.7
1	D	359	GLU	2.7
1	C	384	GLY	2.7
1	B	104	LYS	2.6
1	C	85	GLU	2.6
1	C	9	ASN	2.6
1	D	153	ASP	2.6
1	C	155	GLY	2.5
1	D	2	ILE	2.5
1	D	238	ALA	2.5
1	B	330	GLU	2.5
1	C	88	TYR	2.5
1	D	356	VAL	2.5
1	B	157	MET	2.5
1	C	388	GLU	2.5
1	D	363	SER	2.4
1	A	384	GLY	2.4
1	D	347	LYS	2.4
1	D	133	LYS	2.4
1	C	393	ARG	2.4
1	D	280	ALA	2.4
1	C	396	ARG	2.3
1	B	154	THR	2.3
1	C	38	GLU	2.3
1	D	221	GLU	2.3
1	D	239	LEU	2.3
1	B	397	ILE	2.2
1	D	50	LEU	2.2
1	B	292	ALA	2.2
1	C	383	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	343	GLU	2.2
1	A	75	ILE	2.2
1	A	291	LEU	2.2
1	B	156	ILE	2.2
1	D	181	LEU	2.1
1	B	242	LYS	2.1
1	D	174	VAL	2.1
1	B	147	PRO	2.1
1	D	235	ALA	2.1
1	B	134	GLY	2.1
1	C	87	LYS	2.1
1	A	17	PRO	2.0
1	B	20	CYS	2.0
1	D	64	ILE	2.0
1	B	178	THR	2.0
1	D	150	THR	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.