



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:54 PM UTC

PDB ID : 8WK9 / pdb_00008wk9
Title : Aldo-keto reductase KmAKR - Y28A/K29H/T63M/Q213A/Y296W/W297H
Authors : Xu, S.Y.; Zhou, L.; Xu, Y.; Wang, Y.J.; Zheng, Y.G.
Deposited on : 2023-09-27
Resolution : 2.39 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
Xtriage (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

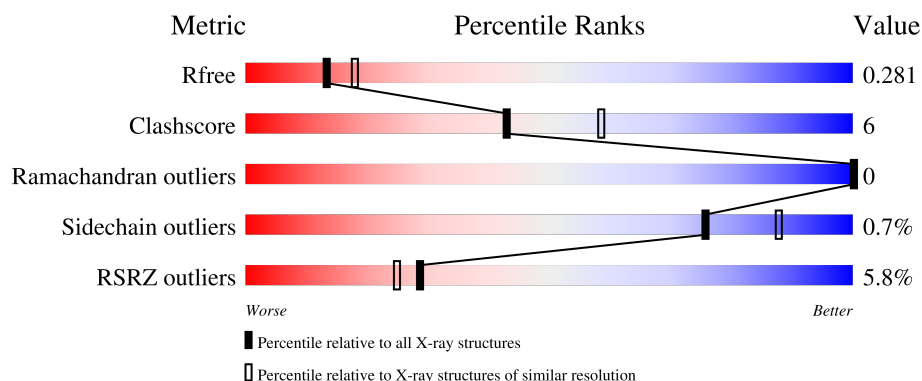
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.39 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	310	3% 83% 16% .
1	B	310	5% 83% 15% .
1	C	310	3% 83% 15% .
1	D	310	11% 78% 18% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9956 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 NADPH-dependent alpha-keto amide reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	308	Total	C	N	O	S	0	0	0
			2478	1602	397	478	1			
1	B	304	Total	C	N	O	S	0	0	0
			2447	1583	391	472	1			
1	C	302	Total	C	N	O	S	0	0	0
			2431	1575	390	465	1			
1	D	297	Total	C	N	O	S	0	0	0
			2389	1551	380	457	1			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	28	ALA	TYR	engineered mutation	UNP W0TDP7
A	29	HIS	LYS	engineered mutation	UNP W0TDP7
A	63	MET	THR	engineered mutation	UNP W0TDP7
A	109	ASN	LYS	conflict	UNP W0TDP7
A	213	ALA	GLN	engineered mutation	UNP W0TDP7
A	296	TRP	TYR	engineered mutation	UNP W0TDP7
A	297	HIS	TRP	engineered mutation	UNP W0TDP7
B	28	ALA	TYR	engineered mutation	UNP W0TDP7
B	29	HIS	LYS	engineered mutation	UNP W0TDP7
B	63	MET	THR	engineered mutation	UNP W0TDP7
B	109	ASN	LYS	conflict	UNP W0TDP7
B	213	ALA	GLN	engineered mutation	UNP W0TDP7
B	296	TRP	TYR	engineered mutation	UNP W0TDP7
B	297	HIS	TRP	engineered mutation	UNP W0TDP7
C	28	ALA	TYR	engineered mutation	UNP W0TDP7
C	29	HIS	LYS	engineered mutation	UNP W0TDP7
C	63	MET	THR	engineered mutation	UNP W0TDP7
C	109	ASN	LYS	conflict	UNP W0TDP7
C	213	ALA	GLN	engineered mutation	UNP W0TDP7
C	296	TRP	TYR	engineered mutation	UNP W0TDP7
C	297	HIS	TRP	engineered mutation	UNP W0TDP7

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	28	ALA	TYR	engineered mutation	UNP W0TDP7
D	29	HIS	LYS	engineered mutation	UNP W0TDP7
D	63	MET	THR	engineered mutation	UNP W0TDP7
D	109	ASN	LYS	conflict	UNP W0TDP7
D	213	ALA	GLN	engineered mutation	UNP W0TDP7
D	296	TRP	TYR	engineered mutation	UNP W0TDP7
D	297	HIS	TRP	engineered mutation	UNP W0TDP7

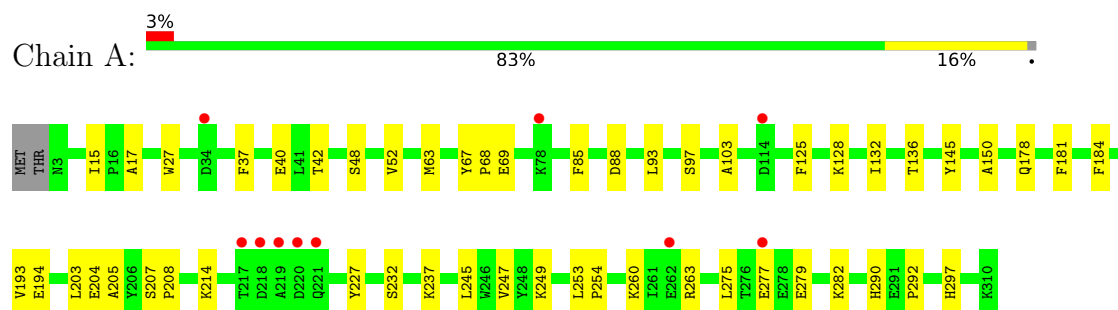
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	60	Total O 60 60	0	0
2	B	47	Total O 47 47	0	0
2	C	54	Total O 54 54	0	0
2	D	50	Total O 50 50	0	0

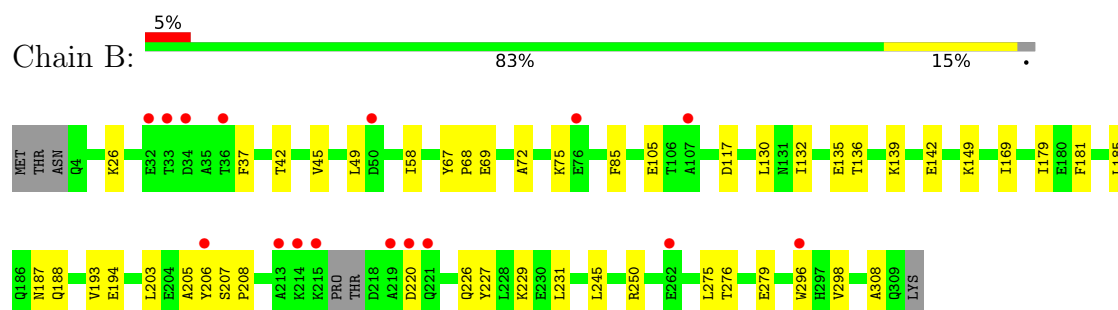
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry 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.

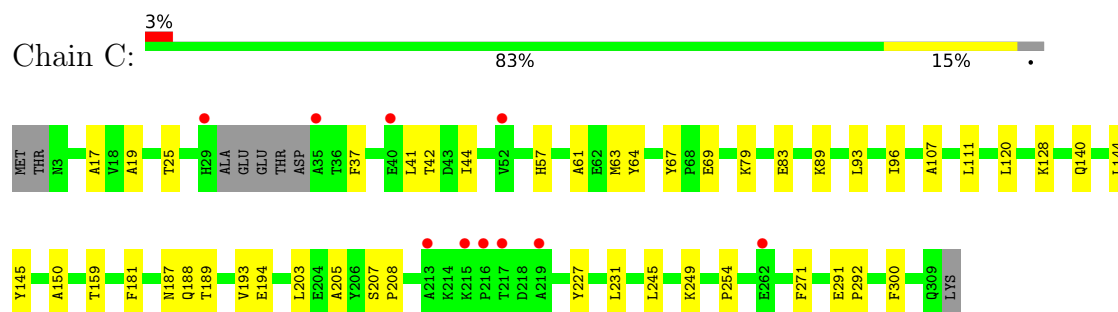
- Molecule 1: NADPH-dependent alpha-keto amide reductase



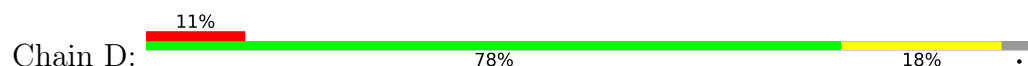
- Molecule 1: NADPH-dependent alpha-keto amide reductase

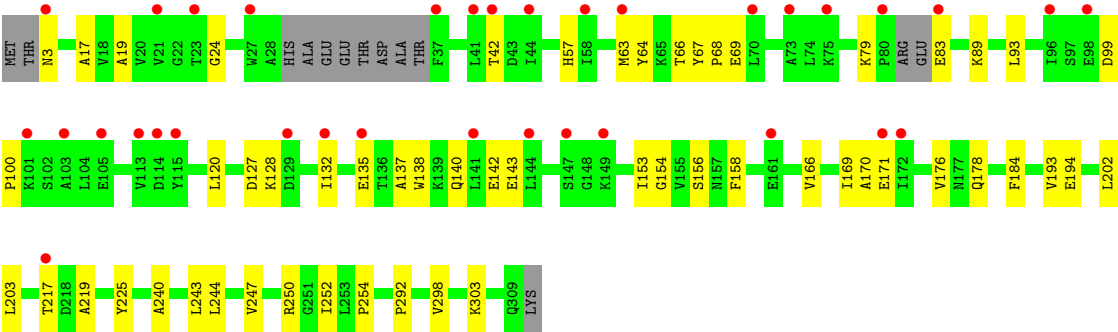


- Molecule 1: NADPH-dependent alpha-keto amide reductase



- Molecule 1: NADPH-dependent alpha-keto amide reductase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	68.28Å 114.83Å 83.07Å 90.00° 101.32° 90.00°	Depositor
Resolution (Å)	81.46 – 2.39 81.46 – 2.39	Depositor EDS
% Data completeness (in resolution range)	99.8 (81.46-2.39) 99.8 (81.46-2.39)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.32 (at 2.40Å)	Xtriage
Refinement program	REFMAC 5.8.0403	Depositor
R, R_{free}	0.220 , 0.283 0.221 , 0.281	Depositor DCC
R_{free} test set	2549 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	22.4	Xtriage
Anisotropy	1.066	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 20.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	9956	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.29% 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

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	0.66	0/2533	0.88	0/3434
1	B	0.63	0/2500	0.84	0/3387
1	C	0.64	0/2485	0.84	0/3368
1	D	0.63	0/2441	0.85	0/3307
All	All	0.64	0/9959	0.85	0/13496

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	B	0	1
1	D	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	250	ARG	Sidechain
1	D	250	ARG	Sidechain

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	2478	0	2489	38	0
1	B	2447	0	2455	30	0
1	C	2431	0	2447	26	0
1	D	2389	0	2408	33	0
2	A	60	0	0	0	0
2	B	47	0	0	0	0
2	C	54	0	0	0	0
2	D	50	0	0	1	0
All	All	9956	0	9799	121	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 6.

All (121) 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:49:LEU:HD21	1:B:58:ILE:HD11	1.44	0.99
1:B:298:VAL:HG11	1:C:128:LYS:HD3	1.73	0.71
1:C:42:THR:HG22	1:C:69:GLU:O	1.91	0.69
1:A:132:ILE:CG1	1:A:136:THR:HB	2.22	0.68
1:B:132:ILE:HD11	1:B:136:THR:HG22	1.77	0.66
1:B:206:TYR:HE1	1:B:296:TRP:CD1	2.15	0.64
1:C:89:LYS:HG2	1:C:120:LEU:HB2	1.79	0.64
1:B:206:TYR:CE1	1:B:296:TRP:CD1	2.88	0.62
1:B:194:GLU:H	1:B:194:GLU:CD	2.09	0.61
1:D:120:LEU:HA	1:D:154:GLY:O	2.02	0.60
1:A:132:ILE:HG12	1:A:136:THR:HB	1.85	0.59
1:B:42:THR:HG22	1:B:69:GLU:CA	2.33	0.59
1:D:135:GLU:HG3	1:D:169:ILE:HG12	1.84	0.59
1:A:125:PHE:CE1	1:D:303:LYS:HE2	2.37	0.58
1:A:260:LYS:HE3	1:A:263:ARG:NH2	2.19	0.58
1:A:42:THR:HG22	1:A:69:GLU:HA	1.86	0.57
1:A:42:THR:HG22	1:A:69:GLU:CA	2.34	0.57
1:D:132:ILE:HD11	1:D:137:ALA:HA	1.87	0.57
1:A:15:ILE:HG13	1:A:85:PHE:CZ	2.39	0.57
1:D:193:VAL:HA	1:D:203:LEU:HD11	1.86	0.57
1:D:67:TYR:N	1:D:68:PRO:CD	2.68	0.56
1:C:227:TYR:CE2	1:C:231:LEU:HD11	2.41	0.56
1:B:37:PHE:CZ	1:B:42:THR:HG21	2.39	0.56
1:D:79:LYS:HD2	1:D:83:GLU:HB3	1.88	0.56
1:D:24:GLY:HA2	1:D:64:TYR:CG	2.41	0.55
1:D:194:GLU:HB3	2:D:436:HOH:O	2.06	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:LEU:O	1:A:249:LYS:HG2	2.07	0.55
1:C:193:VAL:HA	1:C:203:LEU:HD11	1.89	0.55
1:B:220:ASP:O	1:B:226:GLN:NE2	2.27	0.54
1:B:42:THR:HG22	1:B:69:GLU:HA	1.90	0.54
1:B:245:LEU:HD23	1:B:275:LEU:HG	1.91	0.53
1:A:132:ILE:HG13	1:A:136:THR:HB	1.90	0.52
1:B:207:SER:N	1:B:208:PRO:CD	2.73	0.52
1:A:128:LYS:HB3	1:D:298:VAL:HG12	1.92	0.51
1:D:243:LEU:O	1:D:247:VAL:HG23	2.10	0.51
1:A:63:MET:HE1	1:A:93:LEU:HB3	1.91	0.51
1:A:37:PHE:CZ	1:A:42:THR:HG21	2.46	0.51
1:B:67:TYR:N	1:B:68:PRO:CD	2.73	0.51
1:B:181:PHE:HB3	1:B:205:ALA:CB	2.41	0.51
1:A:253:LEU:HD23	1:A:253:LEU:C	2.36	0.51
1:D:89:LYS:HE2	1:D:178:GLN:OE1	2.11	0.51
1:D:138:TRP:CZ2	1:D:153:ILE:HD12	2.46	0.51
1:D:42:THR:HG22	1:D:69:GLU:HA	1.93	0.51
1:C:159:THR:HG22	1:C:189:THR:HG21	1.94	0.50
1:B:45:VAL:O	1:B:49:LEU:HG	2.13	0.49
1:A:194:GLU:H	1:A:194:GLU:CD	2.20	0.49
1:D:184:PHE:O	1:D:292:PRO:HA	2.13	0.49
1:A:63:MET:CE	1:A:93:LEU:HB3	2.43	0.48
1:B:85:PHE:CE2	1:B:117:ASP:HB3	2.48	0.48
1:C:245:LEU:O	1:C:249:LYS:HG2	2.13	0.48
1:C:291:GLU:HG2	1:C:292:PRO:HD2	1.95	0.48
1:B:169:ILE:HD12	1:B:169:ILE:C	2.40	0.47
1:D:142:GLU:OE1	1:D:171:GLU:N	2.45	0.47
1:A:125:PHE:CE1	1:D:303:LYS:CE	2.98	0.47
1:A:27:TRP:CE2	1:C:25:THR:HG22	2.50	0.47
1:A:67:TYR:N	1:A:68:PRO:CD	2.78	0.46
1:A:67:TYR:OH	1:A:88:ASP:OD1	2.27	0.46
1:A:145:TYR:HA	1:A:150:ALA:O	2.15	0.46
1:C:107:ALA:O	1:C:111:LEU:HG	2.16	0.46
1:D:93:LEU:O	1:D:127:ASP:HB2	2.16	0.46
1:A:214:LYS:HD2	1:C:37:PHE:HB2	1.97	0.45
1:D:63:MET:CE	1:D:93:LEU:HB3	2.46	0.45
1:C:181:PHE:HB3	1:C:205:ALA:CB	2.47	0.45
1:D:17:ALA:O	1:D:254:PRO:HD2	2.16	0.45
1:B:227:TYR:CE2	1:B:231:LEU:HD11	2.51	0.45
1:D:169:ILE:HD12	1:D:169:ILE:O	2.17	0.45
1:C:19:ALA:HA	1:C:57:HIS:HB3	1.98	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:67:TYR:CE2	1:C:107:ALA:HB2	2.52	0.44
1:D:99:ASP:HB2	1:D:100:PRO:HD2	1.99	0.44
1:D:138:TRP:CH2	1:D:153:ILE:HD12	2.53	0.44
1:A:207:SER:N	1:A:208:PRO:CD	2.80	0.44
1:A:227:TYR:OH	1:A:279:GLU:HA	2.18	0.44
1:A:260:LYS:HD2	1:A:263:ARG:HG3	1.99	0.44
1:C:145:TYR:HA	1:C:150:ALA:O	2.17	0.44
1:A:48:SER:HA	1:A:52:VAL:HG23	1.99	0.44
1:C:61:ALA:HB3	1:C:64:TYR:CD2	2.53	0.44
1:D:156:SER:O	1:D:158:PHE:HD1	2.00	0.44
1:B:276:THR:OG1	1:B:279:GLU:HG3	2.18	0.44
1:C:194:GLU:H	1:C:194:GLU:CD	2.26	0.44
1:B:72:ALA:HA	1:B:75:LYS:HE3	2.00	0.43
1:A:181:PHE:HB3	1:A:205:ALA:CB	2.48	0.43
1:C:79:LYS:HE3	1:C:83:GLU:HB3	2.00	0.43
1:B:42:THR:HG22	1:B:69:GLU:O	2.18	0.43
1:B:139:LYS:O	1:B:142:GLU:HB2	2.18	0.43
1:C:41:LEU:O	1:C:44:ILE:HG22	2.18	0.43
1:A:97:SER:HB3	1:A:103:ALA:HB2	2.01	0.43
1:B:135:GLU:HG3	1:B:169:ILE:HG12	2.00	0.43
1:D:138:TRP:CD1	1:D:166:VAL:HG13	2.54	0.43
1:A:15:ILE:HG13	1:A:85:PHE:CE1	2.54	0.42
1:A:247:VAL:HG11	1:A:254:PRO:HB3	2.00	0.42
1:C:187:ASN:O	1:C:188:GLN:C	2.61	0.42
1:A:178:GLN:HG3	1:A:204:GLU:HB2	1.99	0.42
1:C:17:ALA:HB3	1:C:271:PHE:CZ	2.54	0.42
1:C:17:ALA:O	1:C:254:PRO:HD2	2.19	0.42
1:D:176:VAL:HA	1:D:202:LEU:O	2.19	0.42
1:B:42:THR:CG2	1:B:69:GLU:HA	2.48	0.42
1:B:179:ILE:O	1:B:205:ALA:HA	2.20	0.42
1:C:63:MET:HA	1:C:96:ILE:CD1	2.48	0.42
1:D:203:LEU:HB2	1:D:252:ILE:HG12	2.01	0.42
1:D:240:ALA:O	1:D:244:LEU:HG	2.20	0.42
1:A:227:TYR:CZ	1:A:282:LYS:HD3	2.54	0.42
1:C:93:LEU:HD11	1:C:300:PHE:CZ	2.55	0.42
1:C:207:SER:N	1:C:208:PRO:CD	2.82	0.42
1:D:64:TYR:O	1:D:66:THR:HG23	2.20	0.42
1:A:17:ALA:O	1:A:254:PRO:HD2	2.19	0.41
1:B:105:GLU:OE2	1:B:149:LYS:HD3	2.20	0.41
1:A:184:PHE:CZ	1:A:290:HIS:CD2	3.08	0.41
1:B:130:LEU:HD23	1:B:130:LEU:HA	1.91	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:185:LEU:HD12	1:B:308:ALA:HB1	2.02	0.41
1:D:140:GLN:O	1:D:143:GLU:HB2	2.21	0.41
1:D:166:VAL:O	1:D:170:ALA:HB2	2.20	0.41
1:A:194:GLU:CD	1:A:194:GLU:N	2.78	0.41
1:A:232:SER:HB2	1:A:237:LYS:O	2.21	0.41
1:A:245:LEU:HD23	1:A:275:LEU:HG	2.03	0.41
1:B:193:VAL:HA	1:B:203:LEU:HD11	2.03	0.41
1:D:19:ALA:HA	1:D:57:HIS:HB3	2.02	0.40
1:D:219:ALA:HB1	1:D:225:TYR:CD2	2.56	0.40
1:A:184:PHE:O	1:A:292:PRO:HA	2.20	0.40
1:A:193:VAL:HA	1:A:203:LEU:HD11	2.02	0.40
1:C:140:GLN:O	1:C:144:LEU:HG	2.21	0.40
1:B:187:ASN:O	1:B:188:GLN:C	2.63	0.40

There are no symmetry-related clashes.

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	306/310 (99%)	301 (98%)	5 (2%)	0	100	100
1	B	300/310 (97%)	294 (98%)	6 (2%)	0	100	100
1	C	298/310 (96%)	293 (98%)	5 (2%)	0	100	100
1	D	291/310 (94%)	281 (97%)	10 (3%)	0	100	100
All	All	1195/1240 (96%)	1169 (98%)	26 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	274/276 (99%)	271 (99%)	3 (1%)	65	82
1	B	270/276 (98%)	268 (99%)	2 (1%)	76	88
1	C	269/276 (98%)	269 (100%)	0	100	100
1	D	265/276 (96%)	262 (99%)	3 (1%)	65	82
All	All	1078/1104 (98%)	1070 (99%)	8 (1%)	76	88

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	GLU
1	A	277	GLU
1	A	297	HIS
1	B	26	LYS
1	B	229	LYS
1	D	3	ASN
1	D	128	LYS
1	D	217	THR

Sometimes 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	GLN
1	A	94	HIS
1	A	109	ASN
1	A	131	ASN
1	A	187	ASN
1	A	222	GLN
1	B	39	GLN
1	B	187	ASN
1	B	289	GLN
1	C	94	HIS
1	C	131	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	297	HIS
1	D	3	ASN
1	D	297	HIS
1	D	309	GLN

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 [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	308/310 (99%)	0.34	10 (3%) 50 46	10, 23, 46, 82	0
1	B	304/310 (98%)	0.57	16 (5%) 32 28	11, 28, 52, 101	0
1	C	302/310 (97%)	0.43	10 (3%) 49 45	9, 25, 51, 85	0
1	D	297/310 (95%)	0.76	34 (11%) 10 7	10, 29, 53, 81	0
All	All	1211/1240 (97%)	0.52	70 (5%) 29 25	9, 26, 52, 101	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	217	THR	4.5
1	C	29	HIS	4.3
1	A	218	ASP	3.8
1	C	35	ALA	3.5
1	C	40	GLU	3.5
1	D	70	LEU	3.4
1	D	147	SER	3.4
1	D	98	GLU	3.3
1	D	132	ILE	3.2
1	D	21	VAL	3.2
1	D	172	ILE	3.2
1	D	3	ASN	3.1
1	D	75	LYS	3.1
1	D	217	THR	3.0
1	C	217	THR	3.0
1	C	215	LYS	3.0
1	B	107	ALA	3.0
1	B	219	ALA	3.0
1	B	33	THR	2.9
1	D	135	GLU	2.9
1	C	213	ALA	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	219	ALA	2.7
1	A	221	GLN	2.7
1	D	44	ILE	2.7
1	B	36	THR	2.7
1	B	262	GLU	2.6
1	B	215	LYS	2.6
1	A	219	ALA	2.6
1	C	262	GLU	2.5
1	D	80	PRO	2.5
1	B	221	GLN	2.5
1	A	277	GLU	2.5
1	B	206	TYR	2.5
1	D	115	TYR	2.5
1	B	214	LYS	2.5
1	D	144	LEU	2.4
1	B	50	ASP	2.4
1	D	41	LEU	2.4
1	D	83	GLU	2.4
1	A	34	ASP	2.3
1	A	114	ASP	2.3
1	C	216	PRO	2.3
1	D	103	ALA	2.3
1	D	96	ILE	2.3
1	B	34	ASP	2.3
1	D	129	ASP	2.3
1	A	78	LYS	2.3
1	B	213	ALA	2.3
1	D	63	MET	2.3
1	D	171	GLU	2.3
1	D	27	TRP	2.3
1	D	105	GLU	2.3
1	D	58	ILE	2.3
1	D	149	LYS	2.3
1	D	73	ALA	2.2
1	B	220	ASP	2.2
1	D	42	THR	2.2
1	B	76	GLU	2.2
1	D	114	ASP	2.2
1	D	113	VAL	2.1
1	C	52	VAL	2.1
1	D	101	LYS	2.1
1	B	296	TRP	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	262	GLU	2.0
1	B	32	GLU	2.0
1	D	161	GLU	2.0
1	A	220	ASP	2.0
1	D	141	LEU	2.0
1	D	37	PHE	2.0
1	D	23	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 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.