



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 12:49 am BST

PDB ID : 3ZGY
Title : Apo-structure of R-selective imine reductase from *Streptomyces kanamyceticus*
Authors : Rodriguez Mata, M.; Frank, A.; Grogan, G.
Deposited on : 2012-12-19
Resolution : 2.71 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

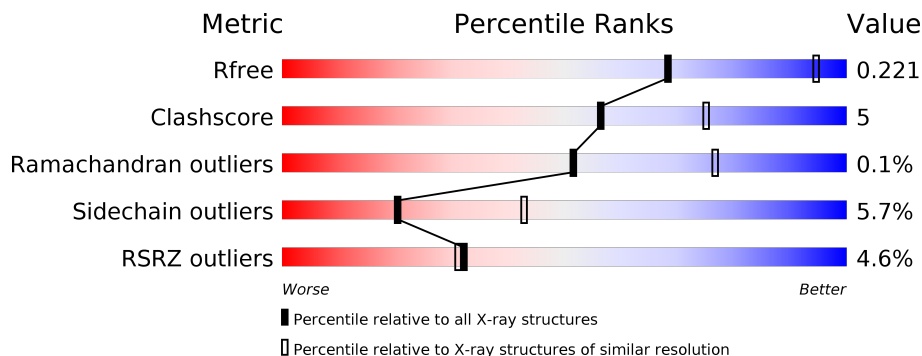
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.71 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}	130704	3359 (2.74-2.70)
Clashscore	141614	3686 (2.74-2.70)
Ramachandran outliers	138981	3622 (2.74-2.70)
Sidechain outliers	138945	3623 (2.74-2.70)
RSRZ outliers	127900	3276 (2.74-2.70)

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 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	309	 3% 80% 11% • 8%
1	B	309	 8% 78% 9% • 12%
1	C	309	 4% 76% 11% • 11%
1	D	309	 3% 83% 11% • 5%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8019 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 R-IMINE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	285	1998	1263	345	385	5	0	0	0
1	B	273	1914	1211	330	368	5	0	0	0
1	C	274	1923	1220	329	369	5	0	0	0
1	D	294	2077	1311	363	397	6	0	0	0

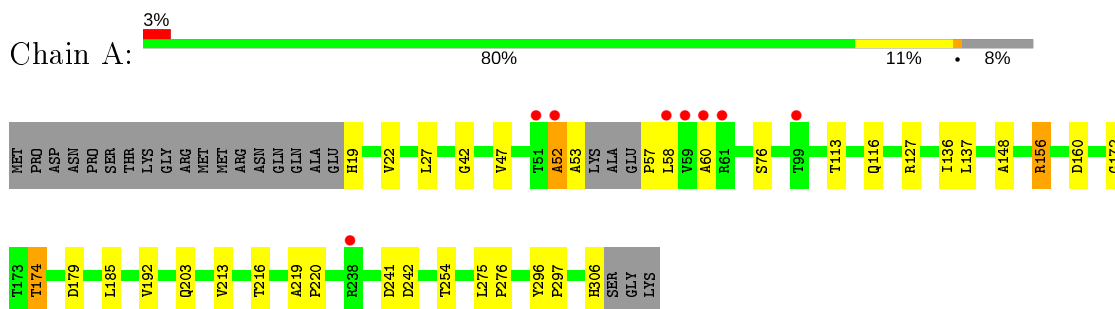
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	33	Total	O	0	0
			33	33		
2	B	19	Total	O	0	0
			19	19		
2	C	19	Total	O	0	0
			19	19		
2	D	36	Total	O	0	0
			36	36		

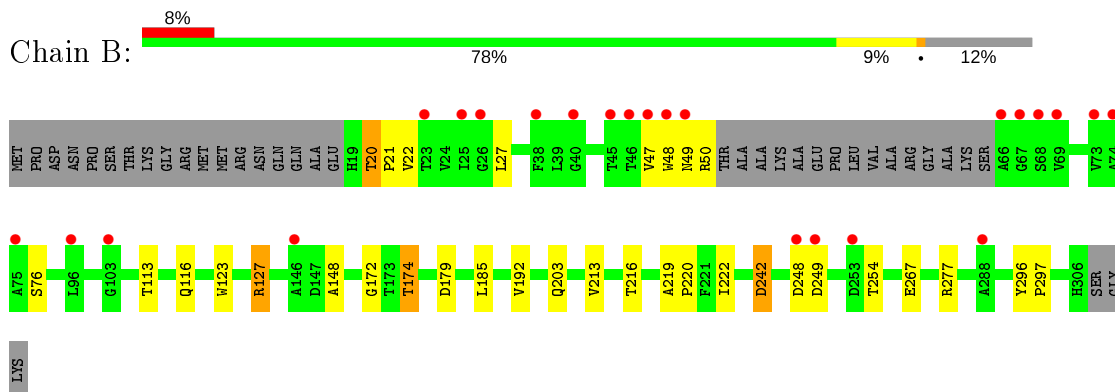
3 Residue-property plots [i](#)

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

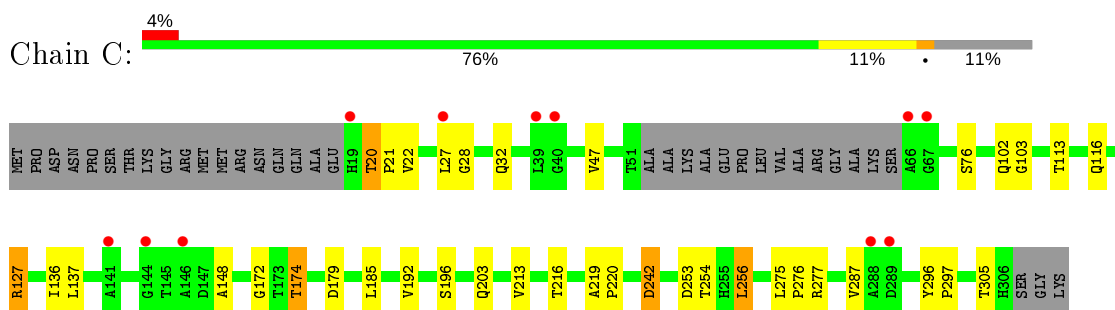
- Molecule 1: R-IMINE REDUCTASE



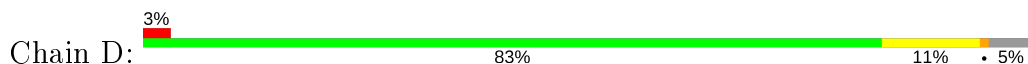
- Molecule 1: R-IMINE REDUCTASE

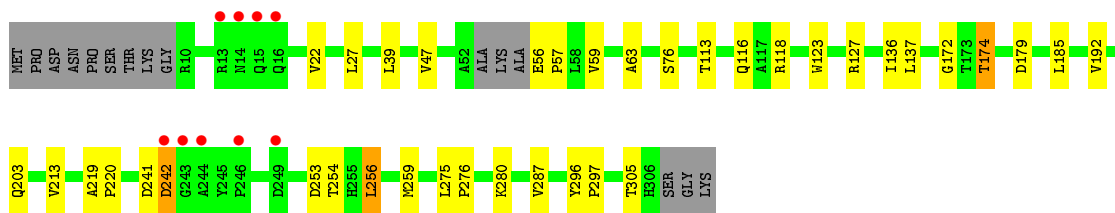


- Molecule 1: R-IMINE REDUCTASE



- Molecule 1: R-IMINE REDUCTASE





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	203.41Å 131.11Å 77.48Å 90.00° 107.22° 90.00°	Depositor
Resolution (Å)	36.23 – 2.71 36.23 – 2.62	Depositor EDS
% Data completeness (in resolution range)	99.3 (36.23-2.71) 99.4 (36.23-2.62)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 2.61Å)	Xtrriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.185 , 0.214 0.194 , 0.221	Depositor DCC
R_{free} test set	2943 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	60.3	Xtrriage
Anisotropy	0.565	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 50.1	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	0.95	EDS
Total number of atoms	8019	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.50% 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	0.77	0/2041	0.82	2/2796 (0.1%)
1	B	0.73	1/1956 (0.1%)	0.81	2/2680 (0.1%)
1	C	0.68	0/1964	0.79	2/2690 (0.1%)
1	D	0.80	1/2121 (0.0%)	0.85	2/2902 (0.1%)
All	All	0.75	2/8082 (0.0%)	0.82	8/11068 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	123	TRP	CB-CG	-5.59	1.40	1.50
1	D	123	TRP	CG-CD1	-5.08	1.29	1.36

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	179	ASP	CB-CG-OD1	8.30	125.77	118.30
1	D	179	ASP	CB-CG-OD1	7.41	124.97	118.30
1	C	179	ASP	CB-CG-OD1	7.06	124.65	118.30
1	B	179	ASP	CB-CG-OD1	6.88	124.49	118.30
1	A	156	ARG	NE-CZ-NH1	-6.11	117.25	120.30
1	B	127	ARG	NE-CZ-NH1	5.72	123.16	120.30
1	C	127	ARG	NE-CZ-NH1	5.64	123.12	120.30
1	D	118	ARG	NE-CZ-NH1	5.04	122.82	120.30

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	1998	0	1919	20	0
1	B	1914	0	1825	22	0
1	C	1923	0	1848	21	0
1	D	2077	0	2001	21	0
2	A	33	0	0	0	0
2	B	19	0	0	1	0
2	C	19	0	0	1	0
2	D	36	0	0	0	0
All	All	8019	0	7593	78	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 5.

All (78) 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:172:GLY:O	1:B:174:THR:HG22	1.82	0.80
1:C:172:GLY:O	1:C:174:THR:HG22	1.83	0.79
1:D:172:GLY:O	1:D:174:THR:HG22	1.83	0.78
1:A:172:GLY:O	1:A:174:THR:HG22	1.86	0.75
1:D:56:GLU:HB3	1:D:57:PRO:HD3	1.70	0.71
1:D:56:GLU:HB3	1:D:57:PRO:CD	2.22	0.70
1:D:113:THR:H	1:D:116:GLN:HE21	1.38	0.69
1:C:253:ASP:O	1:C:256:LEU:HB2	1.93	0.69
1:B:113:THR:H	1:B:116:GLN:HE21	1.38	0.69
1:C:28:GLY:O	1:C:32:GLN:HG3	1.94	0.67
1:D:253:ASP:O	1:D:256:LEU:HB2	1.95	0.67
1:C:253:ASP:HA	1:C:256:LEU:HD22	1.77	0.66
1:D:253:ASP:HA	1:D:256:LEU:HD22	1.77	0.65
1:D:27:LEU:HD11	1:D:47:VAL:CG1	2.27	0.64
1:A:113:THR:H	1:A:116:GLN:HE21	1.43	0.64
1:C:103:GLY:C	1:C:103:GLY:N	2.53	0.62
1:A:306:HIS:HB2	2:B:2018:HOH:O	2.00	0.61
1:C:27:LEU:HD11	1:C:47:VAL:CG1	2.31	0.61
1:A:57:PRO:O	1:A:60:ALA:N	2.34	0.61
1:D:27:LEU:HD11	1:D:47:VAL:HG13	1.83	0.60
1:B:27:LEU:HD11	1:B:47:VAL:CG1	2.31	0.60
1:C:113:THR:H	1:C:116:GLN:HE21	1.47	0.60
1:A:27:LEU:HD11	1:A:47:VAL:CG1	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:267:GLU:OE2	1:B:277:ARG:NH1	2.34	0.60
1:B:49:ASN:CG	1:B:50:ARG:H	2.06	0.59
1:C:256:LEU:HD21	1:C:287:VAL:HG21	1.85	0.57
1:B:242:ASP:N	1:B:242:ASP:OD1	2.37	0.56
1:D:256:LEU:HD21	1:D:287:VAL:HG21	1.86	0.56
1:A:27:LEU:HD11	1:A:47:VAL:HG13	1.87	0.56
1:C:27:LEU:HD11	1:C:47:VAL:HG13	1.88	0.56
1:B:27:LEU:HD11	1:B:47:VAL:HG13	1.88	0.54
1:B:48:TRP:CE3	1:B:49:ASN:N	2.78	0.52
1:A:219:ALA:HB3	1:A:220:PRO:HD3	1.92	0.51
1:B:20:THR:HG23	1:B:21:PRO:HD2	1.92	0.51
1:B:113:THR:H	1:B:116:GLN:NE2	2.08	0.51
1:B:48:TRP:CE3	1:B:48:TRP:C	2.85	0.50
1:B:219:ALA:HB3	1:B:220:PRO:HD3	1.94	0.50
1:C:20:THR:HG22	1:C:21:PRO:HD2	1.93	0.50
1:D:259:MET:HB2	1:D:280:LYS:HD3	1.94	0.49
1:D:296:TYR:N	1:D:297:PRO:CD	2.76	0.49
1:D:113:THR:H	1:D:116:GLN:NE2	2.08	0.48
1:A:113:THR:H	1:A:116:GLN:NE2	2.08	0.48
1:A:296:TYR:N	1:A:297:PRO:CD	2.77	0.48
1:B:48:TRP:CE3	1:B:49:ASN:HA	2.48	0.48
1:C:219:ALA:HB3	1:C:220:PRO:HD3	1.95	0.48
1:B:296:TYR:N	1:B:297:PRO:CD	2.77	0.48
1:C:192:VAL:HG21	1:D:203:GLN:HG2	1.96	0.47
1:D:219:ALA:HB3	1:D:220:PRO:HD3	1.95	0.47
1:B:48:TRP:CE3	1:B:49:ASN:CA	2.97	0.47
1:C:113:THR:H	1:C:116:GLN:NE2	2.12	0.47
1:D:56:GLU:CB	1:D:57:PRO:CD	2.86	0.47
1:A:52:ALA:O	1:A:53:ALA:HB3	2.14	0.46
1:C:216:THR:HG23	1:D:241:ASP:OD2	2.15	0.46
1:A:57:PRO:O	1:A:60:ALA:HB3	2.16	0.46
1:B:49:ASN:ND2	1:B:50:ARG:H	2.13	0.46
1:C:296:TYR:N	1:C:297:PRO:CD	2.79	0.45
1:C:275:LEU:HB3	1:C:276:PRO:HD3	1.99	0.44
1:D:242:ASP:OD1	1:D:242:ASP:N	2.50	0.44
1:A:241:ASP:OD2	1:B:216:THR:HG23	2.17	0.44
1:A:57:PRO:O	1:A:58:LEU:C	2.55	0.43
1:A:19:HIS:HA	1:A:42:GLY:O	2.19	0.43
1:C:277:ARG:NE	2:C:2017:HOH:O	2.41	0.43
1:C:203:GLN:HG2	1:D:192:VAL:HG21	2.01	0.43
1:C:136:ILE:C	1:C:137:LEU:HD12	2.39	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:148:ALA:O	1:C:172:GLY:HA3	2.20	0.42
1:D:136:ILE:C	1:D:137:LEU:HD12	2.40	0.42
1:A:156:ARG:NH1	1:A:160:ASP:OD1	2.52	0.42
1:B:148:ALA:O	1:B:172:GLY:HA3	2.20	0.42
1:D:39:LEU:HD11	1:D:63:ALA:HB2	2.02	0.42
1:B:222:ILE:HD12	1:B:222:ILE:HA	1.97	0.42
1:A:192:VAL:HG21	1:B:203:GLN:HG2	2.01	0.42
1:A:148:ALA:O	1:A:172:GLY:HA3	2.20	0.41
1:A:203:GLN:HG2	1:B:192:VAL:HG21	2.02	0.41
1:A:275:LEU:HB3	1:A:276:PRO:HD3	2.02	0.41
1:C:242:ASP:OD1	1:C:242:ASP:N	2.53	0.41
1:D:275:LEU:HB3	1:D:276:PRO:HD3	2.03	0.41
1:B:248:ASP:O	1:B:249:ASP:HB2	2.21	0.40
1:A:136:ILE:C	1:A:137:LEU:HD12	2.42	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	281/309 (91%)	276 (98%)	4 (1%)	1 (0%)	34	58
1	B	269/309 (87%)	261 (97%)	8 (3%)	0	100	100
1	C	269/309 (87%)	260 (97%)	9 (3%)	0	100	100
1	D	290/309 (94%)	280 (97%)	10 (3%)	0	100	100
All	All	1109/1236 (90%)	1077 (97%)	31 (3%)	1 (0%)	51	77

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	52	ALA

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	185/214 (86%)	176 (95%)	9 (5%)	25	50
1	B	176/214 (82%)	167 (95%)	9 (5%)	24	48
1	C	179/214 (84%)	166 (93%)	13 (7%)	14	31
1	D	193/214 (90%)	182 (94%)	11 (6%)	20	43
All	All	733/856 (86%)	691 (94%)	42 (6%)	20	43

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	VAL
1	A	76	SER
1	A	127	ARG
1	A	174	THR
1	A	185	LEU
1	A	213	VAL
1	A	216	THR
1	A	242	ASP
1	A	254	THR
1	B	20	THR
1	B	22	VAL
1	B	76	SER
1	B	127	ARG
1	B	174	THR
1	B	185	LEU
1	B	213	VAL
1	B	242	ASP
1	B	254	THR
1	C	20	THR
1	C	22	VAL
1	C	76	SER
1	C	102	GLN
1	C	127	ARG
1	C	174	THR

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Mol	Chain	Res	Type
1	C	185	LEU
1	C	196	SER
1	C	213	VAL
1	C	242	ASP
1	C	254	THR
1	C	256	LEU
1	C	305	THR
1	D	22	VAL
1	D	59	VAL
1	D	76	SER
1	D	127	ARG
1	D	174	THR
1	D	185	LEU
1	D	213	VAL
1	D	242	ASP
1	D	254	THR
1	D	256	LEU
1	D	305	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	116	GLN
1	A	203	GLN
1	B	49	ASN
1	B	116	GLN
1	B	203	GLN
1	C	116	GLN
1	C	203	GLN
1	D	116	GLN
1	D	203	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 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	285/309 (92%)	0.09	8 (2%) 53 54	37, 57, 97, 130	0
1	B	273/309 (88%)	0.36	24 (8%) 10 8	34, 62, 112, 134	0
1	C	274/309 (88%)	0.20	11 (4%) 38 37	41, 71, 112, 126	0
1	D	294/309 (95%)	-0.23	9 (3%) 49 50	31, 52, 83, 113	0
All	All	1126/1236 (91%)	0.10	52 (4%) 32 31	31, 60, 107, 134	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	60	ALA	4.8
1	B	67	GLY	4.7
1	A	51	THR	4.3
1	A	58	LEU	4.0
1	B	66	ALA	3.9
1	B	26	GLY	3.8
1	B	73	VAL	3.7
1	B	249	ASP	3.6
1	B	69	VAL	3.6
1	B	74	ALA	3.3
1	C	141	ALA	3.3
1	D	244	ALA	3.2
1	B	40	GLY	3.2
1	C	40	GLY	3.2
1	B	103	GLY	3.2
1	C	39	LEU	3.1
1	A	59	VAL	3.0
1	C	146	ALA	2.7
1	B	38	PHE	2.7
1	B	48	TRP	2.7
1	B	45	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	14	ASN	2.6
1	B	146	ALA	2.6
1	B	253	ASP	2.6
1	D	243	GLY	2.6
1	D	16	GLN	2.5
1	B	68	SER	2.5
1	D	249	ASP	2.5
1	B	47	VAL	2.5
1	B	25	ILE	2.5
1	B	96	LEU	2.5
1	A	99	THR	2.5
1	C	288	ALA	2.4
1	B	288	ALA	2.4
1	B	248	ASP	2.4
1	A	238	ARG	2.4
1	B	75	ALA	2.3
1	C	66	ALA	2.3
1	C	67	GLY	2.3
1	C	27	LEU	2.3
1	A	61	ARG	2.3
1	D	13	ARG	2.3
1	D	242	ASP	2.3
1	A	52	ALA	2.2
1	C	144	GLY	2.2
1	C	289	ASP	2.2
1	B	23	THR	2.1
1	C	19	HIS	2.1
1	D	246	PRO	2.1
1	B	46	THR	2.0
1	D	15	GLN	2.0
1	B	49	ASN	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

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.