



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:11 PM UTC

PDB ID : 1WCX / pdb_00001wcx
Title : Crystal Structure of Mutant Uroporphyrinogen III Synthase from an Extremely Thermophilic Bacterium *Thermus thermophilus* HB8 (L75M/I193M/L248M, SeMet derivative, Form-1 crystal)
Authors : Mizohata, E.; Matsuura, T.; Murayama, K.; Sakai, H.; Terada, T.; Shirouzu, M.; Kuramitsu, S.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2004-05-06
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

<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
Mogul	:	2022.3.0, CSD as543be (2022)
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)

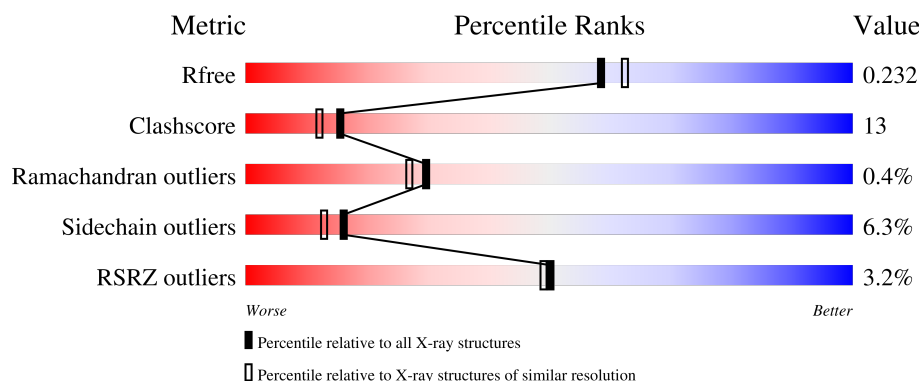
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

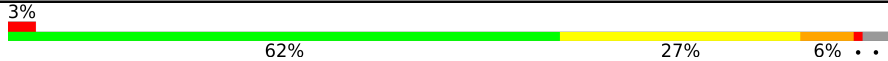
The reported resolution of this entry is 2.00 Å.

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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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 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	261	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2108 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 Uroporphyrinogen III Synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	252	Total	C	N	O	Se	0	0	0
			1930	1241	346	339	4			

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	cloning artifact	UNP Q5SKH2
A	2	ARG	-	cloning artifact	UNP Q5SKH2
A	3	ARG	-	cloning artifact	UNP Q5SKH2
A	4	LEU	-	cloning artifact	UNP Q5SKH2
A	5	GLU	-	cloning artifact	UNP Q5SKH2
A	6	GLU	-	cloning artifact	UNP Q5SKH2
A	7	ASP	-	cloning artifact	UNP Q5SKH2
A	8	ALA	-	cloning artifact	UNP Q5SKH2
A	9	VAL	-	cloning artifact	UNP Q5SKH2
A	75	MSE	LEU	engineered mutation	UNP Q5SKH2
A	161	MSE	MET	modified residue	UNP Q5SKH2
A	193	MSE	ILE	engineered mutation	UNP Q5SKH2
A	248	MSE	LEU	engineered mutation	UNP Q5SKH2

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		

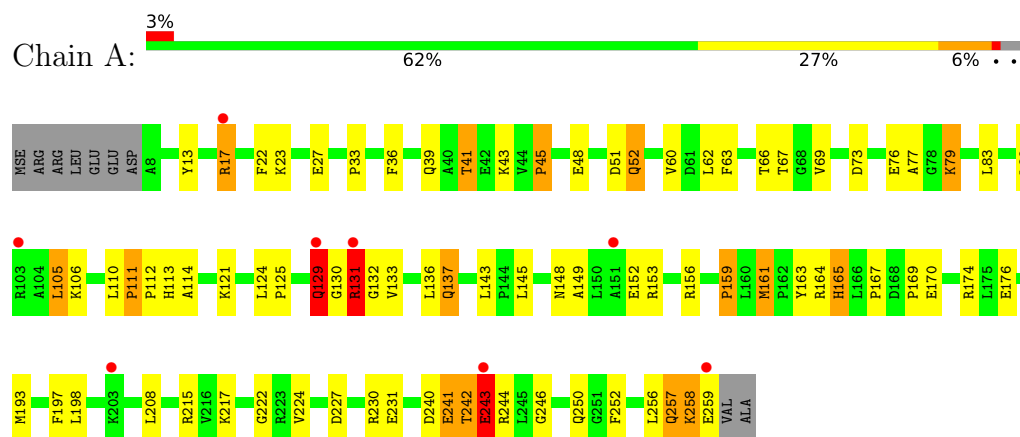
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	172	Total	O	0	0
			172	172		

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: Uroporphyrinogen III Synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	35.55Å 66.11Å 55.86Å 90.00° 91.40° 90.00°	Depositor
Resolution (Å)	42.66 – 2.00 42.66 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.0 (42.66-2.00) 99.1 (42.66-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.17 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.190 , 0.225 0.194 , 0.232	Depositor DCC
R_{free} test set	1704 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	20.9	Xtriage
Anisotropy	0.299	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.026 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2108	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.78% 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

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

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	1.88	26/1962 (1.3%)	1.35	21/2646 (0.8%)

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	2

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	129	GLN	C-N	-27.87	0.92	1.33
1	A	240	ASP	C-N	12.74	1.53	1.33
1	A	161	MSE	SE-CE	-8.36	1.70	1.95
1	A	132	GLY	C-N	-7.88	1.23	1.33
1	A	33	PRO	C-O	-7.60	1.15	1.23
1	A	252	PHE	C-O	-7.01	1.16	1.24
1	A	241	GLU	C-N	-6.70	1.23	1.33
1	A	79	LYS	N-CA	6.66	1.54	1.46
1	A	182	GLY	C-O	-6.57	1.15	1.23
1	A	230	ARG	NE-CZ	6.36	1.40	1.33
1	A	112	PRO	CA-C	6.08	1.60	1.52
1	A	136	LEU	C-N	-6.01	1.25	1.33
1	A	105	LEU	N-CA	5.99	1.53	1.46
1	A	161	MSE	CA-C	-5.91	1.46	1.52
1	A	165	HIS	C-N	5.88	1.39	1.33
1	A	45	PRO	N-CA	5.67	1.54	1.47
1	A	222	GLY	C-O	5.66	1.31	1.23
1	A	159	PRO	C-N	5.62	1.41	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	79	LYS	C-O	-5.50	1.17	1.24
1	A	77	ALA	C-O	5.41	1.30	1.24
1	A	137	GLN	CA-C	5.36	1.59	1.52
1	A	130	GLY	C-N	-5.29	1.26	1.33
1	A	36	PHE	C-O	-5.28	1.18	1.24
1	A	111	PRO	C-O	-5.25	1.18	1.24
1	A	60	VAL	C-O	-5.12	1.18	1.23
1	A	170	GLU	C-O	-5.04	1.18	1.24

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	183	GLU	N-CA-CB	-14.54	88.54	110.20
1	A	240	ASP	CA-C-N	-13.53	100.56	122.08
1	A	240	ASP	C-N-CA	-13.53	100.56	122.08
1	A	164	ARG	CB-CA-C	-7.08	95.33	109.35
1	A	132	GLY	O-C-N	6.74	131.46	122.70
1	A	161	MSE	CA-C-N	6.49	126.07	119.19
1	A	161	MSE	C-N-CA	6.49	126.07	119.19
1	A	244	ARG	N-CA-C	-6.00	99.87	109.96
1	A	143	LEU	CA-C-N	5.99	125.70	119.05
1	A	143	LEU	C-N-CA	5.99	125.70	119.05
1	A	242	THR	N-CA-C	5.98	120.31	112.89
1	A	133	VAL	N-CA-C	5.60	116.01	108.17
1	A	17	ARG	N-CA-C	5.56	118.27	111.82
1	A	131	ARG	NE-CZ-NH2	5.46	124.11	119.20
1	A	244	ARG	N-CA-CB	5.38	118.94	110.23
1	A	129	GLN	O-C-N	-5.27	116.87	123.04
1	A	165	HIS	N-CA-C	-5.20	100.69	109.07
1	A	93	PHE	N-CA-C	-5.12	101.61	109.76
1	A	182	GLY	CA-C-N	5.12	129.33	120.58
1	A	182	GLY	C-N-CA	5.12	129.33	120.58
1	A	243	GLU	N-CA-C	5.10	121.47	113.72

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	129	GLN	Mainchain
1	A	241	GLU	Mainchain

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	1930	0	2021	53	1
2	A	6	0	8	0	0
3	A	172	0	0	15	1
All	All	2108	0	2029	53	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 13.

All (53) 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:156:ARG:NH1	3:A:1145:HOH:O	1.87	1.04
1:A:94:ARG:H	1:A:113:HIS:HD2	1.12	0.95
1:A:242:THR:O	1:A:243:GLU:OE1	1.94	0.86
1:A:181:ARG:HD2	1:A:183:GLU:OE1	1.82	0.80
1:A:76:GLU:HG2	3:A:1099:HOH:O	1.83	0.78
1:A:94:ARG:H	1:A:113:HIS:CD2	2.02	0.74
1:A:67:THR:H	1:A:137:GLN:HE22	1.35	0.74
1:A:121:LYS:HA	1:A:124:LEU:HD23	1.73	0.70
1:A:257:GLN:HE21	1:A:257:GLN:HA	1.56	0.70
1:A:106:LYS:HD3	1:A:111:PRO:HB3	1.73	0.69
1:A:176:GLU:O	1:A:180:LEU:HD23	1.99	0.63
1:A:41:THR:CG2	1:A:165:HIS:HD2	2.12	0.61
1:A:45:PRO:HB2	1:A:48:GLU:HG3	1.83	0.61
1:A:131:ARG:HG2	3:A:1149:HOH:O	2.01	0.59
1:A:156:ARG:CZ	3:A:1145:HOH:O	2.43	0.58
1:A:259:GLU:OE2	3:A:1150:HOH:O	2.17	0.57
1:A:17:ARG:HG2	3:A:1125:HOH:O	2.05	0.57
1:A:149:ALA:O	1:A:153:ARG:HG2	2.04	0.57
1:A:193:MSE:HE2	1:A:197:PHE:CZ	2.40	0.56
1:A:198:LEU:HG	1:A:208:LEU:HD21	1.87	0.56
1:A:39:GLN:HE21	1:A:41:THR:CG2	2.20	0.55
1:A:165:HIS:HB2	1:A:193:MSE:HE1	1.88	0.55
1:A:246:GLY:O	1:A:250:GLN:HG3	2.08	0.54
1:A:161:MSE:HE2	3:A:1041:HOH:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:LEU:CD2	1:A:114:ALA:HB3	2.40	0.51
1:A:41:THR:HG22	1:A:165:HIS:HD2	1.73	0.51
1:A:180:LEU:HA	1:A:215:ARG:NH1	2.26	0.51
1:A:66:THR:H	1:A:137:GLN:NE2	2.09	0.51
1:A:13:TYR:HB2	1:A:22:PHE:CE2	2.46	0.51
1:A:148:ASN:O	1:A:152:GLU:HG3	2.11	0.51
1:A:39:GLN:HE21	1:A:41:THR:HG21	1.77	0.49
1:A:181:ARG:NH1	3:A:1110:HOH:O	2.30	0.49
1:A:217:LYS:NZ	3:A:1051:HOH:O	2.45	0.49
1:A:69:VAL:HG13	1:A:163:TYR:CZ	2.48	0.48
1:A:181:ARG:NE	3:A:1110:HOH:O	2.40	0.48
1:A:62:LEU:HD12	1:A:63:PHE:N	2.29	0.46
1:A:23:LYS:O	1:A:27:GLU:HG3	2.14	0.46
1:A:156:ARG:NE	3:A:1145:HOH:O	2.49	0.45
1:A:43:LYS:HE3	1:A:73:ASP:OD1	2.16	0.45
1:A:41:THR:CG2	1:A:165:HIS:CD2	2.97	0.45
1:A:79:LYS:HE2	3:A:1099:HOH:O	2.17	0.44
1:A:165:HIS:CB	1:A:193:MSE:HE1	2.47	0.44
1:A:124:LEU:N	1:A:125:PRO:HD2	2.32	0.44
1:A:257:GLN:C	1:A:259:GLU:H	2.25	0.44
1:A:165:HIS:HE1	3:A:1073:HOH:O	2.01	0.43
1:A:39:GLN:HG2	3:A:1002:HOH:O	2.19	0.42
1:A:105:LEU:HB3	1:A:110:LEU:O	2.20	0.42
1:A:167:PRO:O	1:A:169:PRO:HD3	2.20	0.42
1:A:106:LYS:CD	1:A:111:PRO:HB3	2.47	0.41
1:A:163:TYR:HE2	1:A:165:HIS:CE1	2.37	0.41
1:A:227:ASP:O	1:A:231:GLU:HG3	2.21	0.41
1:A:224:VAL:HB	3:A:1090:HOH:O	2.20	0.41
1:A:52:GLN:HG2	1:A:159:PRO:O	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)
1:A:174:ARG:NH2	3:A:1099:HOH:O[2_755]	2.01	0.19

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	250/261 (96%)	246 (98%)	3 (1%)	1 (0%)	30	27

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	258	LYS

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	191/194 (98%)	179 (94%)	12 (6%)	16	13

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	THR
1	A	51	ASP
1	A	52	GLN
1	A	83	LEU
1	A	89	LEU
1	A	129	GLN
1	A	131	ARG
1	A	145	LEU
1	A	243	GLU
1	A	256	LEU

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Mol	Chain	Res	Type
1	A	257	GLN
1	A	258	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	HIS
1	A	137	GLN
1	A	165	HIS
1	A	257	GLN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GOL	A	1001	-	5,5,5	1.39	1 (20%)	5,5,5	0.54	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	1001	-	-	2/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	GOL	O3-C3	2.21	1.51	1.42

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1001	GOL	C1-C2-C3-O3
2	A	1001	GOL	O2-C2-C3-O3

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	129:GLN	C	130:GLY	N	0.92

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	248/261 (95%)	-0.00	8 (3%)	50 49	10, 20, 38, 53	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	243	GLU	2.7
1	A	259	GLU	2.5
1	A	131	ARG	2.5
1	A	129	GLN	2.5
1	A	151	ALA	2.5
1	A	203	LYS	2.1
1	A	17	ARG	2.1
1	A	103	ARG	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](#)

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. 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	B-factors($\text{\AA}^2$)	Q<0.9
2	GOL	A	1001	6/6	0.83	0.14	32,35,38,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.