



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:54 AM UTC

PDB ID : 3AX3 / pdb_00003ax3
Title : Crystal structure of rat TOM20-ALDH presequence complex: a complex (form2) between Tom20 and a disulfide-bridged presequence peptide containing D-Cys and L-Cys at the i and i+3 positions.
Authors : Saitoh, T.; Maita, Y.; Kohda, D.
Deposited on : 2011-03-28
Resolution : 2.10 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:

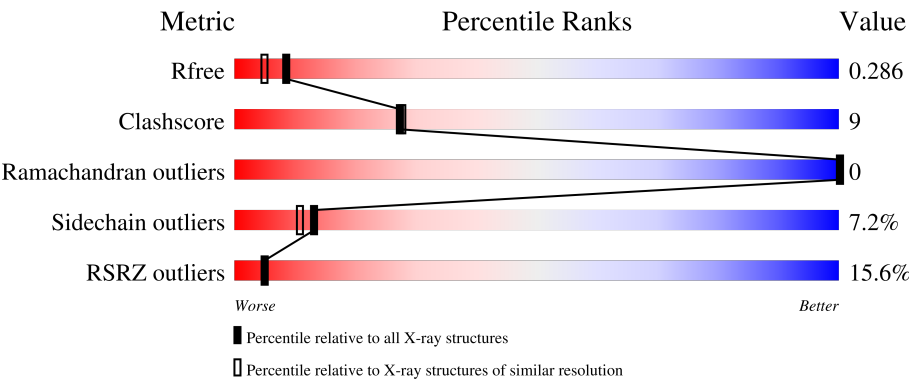
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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.



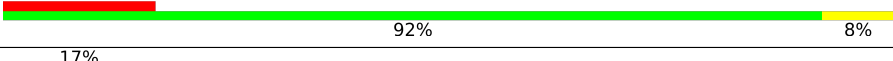
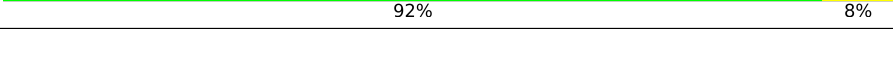
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	73	8% 59% 29% 10%
1	C	73	10% 68% 16% 5% 10%
1	E	73	5% 64% 27% 7%
1	G	73	30% 67% 22% 7%
2	B	12	33% 75% 25%

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Mol	Chain	Length	Quality of chain
2	D	12	
2	F	12	
2	H	12	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2488 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 Mitochondrial import receptor subunit TOM20 homolog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	66	Total	C	N	O	S	0	0	0
			519	334	83	101	1			
1	C	66	Total	C	N	O	S	0	0	0
			519	334	83	101	1			
1	E	68	Total	C	N	O	S	0	0	0
			535	344	85	105	1			
1	G	68	Total	C	N	O	S	0	0	0
			535	344	85	105	1			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	54	GLY	-	expression tag	UNP Q62760
A	55	PRO	-	expression tag	UNP Q62760
A	56	LEU	-	expression tag	UNP Q62760
A	57	GLY	-	expression tag	UNP Q62760
A	58	SER	-	expression tag	UNP Q62760
A	100	SER	CYS	engineered mutation	UNP Q62760
C	54	GLY	-	expression tag	UNP Q62760
C	55	PRO	-	expression tag	UNP Q62760
C	56	LEU	-	expression tag	UNP Q62760
C	57	GLY	-	expression tag	UNP Q62760
C	58	SER	-	expression tag	UNP Q62760
C	100	SER	CYS	engineered mutation	UNP Q62760
E	54	GLY	-	expression tag	UNP Q62760
E	55	PRO	-	expression tag	UNP Q62760
E	56	LEU	-	expression tag	UNP Q62760
E	57	GLY	-	expression tag	UNP Q62760
E	58	SER	-	expression tag	UNP Q62760
E	100	SER	CYS	engineered mutation	UNP Q62760
G	54	GLY	-	expression tag	UNP Q62760
G	55	PRO	-	expression tag	UNP Q62760
G	56	LEU	-	expression tag	UNP Q62760

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Chain	Residue	Modelled	Actual	Comment	Reference
G	57	GLY	-	expression tag	UNP Q62760
G	58	SER	-	expression tag	UNP Q62760
G	100	SER	CYS	engineered mutation	UNP Q62760

- Molecule 2 is a protein called Aldehyde dehydrogenase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	12	Total	C	N	O	S	0	0	1
			86	53	18	13	2			
2	D	12	Total	C	N	O	S	0	0	1
			86	53	18	13	2			
2	F	12	Total	C	N	O	S	0	0	1
			86	53	18	13	2			
2	H	12	Total	C	N	O	S	0	0	1
			86	53	18	13	2			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	13	DCY	PRO	engineered mutation	UNP P11884
B	16	CYS	SER	engineered mutation	UNP P11884
B	21	TYR	-	expression tag	UNP P11884
B	22	ALA	-	expression tag	UNP P11884
B	23	NH2	-	amidation	UNP P11884
D	13	DCY	PRO	engineered mutation	UNP P11884
D	16	CYS	SER	engineered mutation	UNP P11884
D	21	TYR	-	expression tag	UNP P11884
D	22	ALA	-	expression tag	UNP P11884
D	23	NH2	-	amidation	UNP P11884
F	13	DCY	PRO	engineered mutation	UNP P11884
F	16	CYS	SER	engineered mutation	UNP P11884
F	21	TYR	-	expression tag	UNP P11884
F	22	ALA	-	expression tag	UNP P11884
F	23	NH2	-	amidation	UNP P11884
H	13	DCY	PRO	engineered mutation	UNP P11884
H	16	CYS	SER	engineered mutation	UNP P11884
H	21	TYR	-	expression tag	UNP P11884
H	22	ALA	-	expression tag	UNP P11884
H	23	NH2	-	amidation	UNP P11884

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	6	Total 6	O 6	0	0
3	C	10	Total 10	O 10	0	0
3	E	13	Total 13	O 13	0	0
3	F	2	Total 2	O 2	0	0
3	G	3	Total 3	O 3	0	0
3	H	2	Total 2	O 2	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.

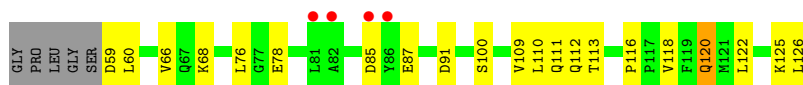
- Molecule 1: Mitochondrial import receptor subunit TOM20 homolog



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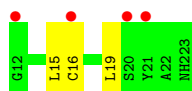
- Molecule 1: Mitochondrial import receptor subunit TOM20 homolog



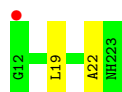
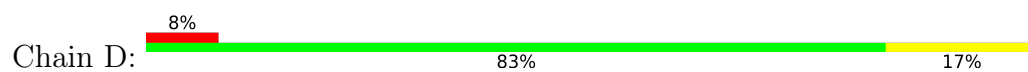
- Molecule 1: Mitochondrial import receptor subunit TOM20 homolog



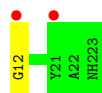
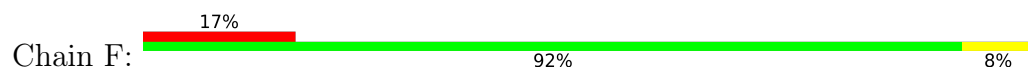
- Molecule 2: Aldehyde dehydrogenase, mitochondrial



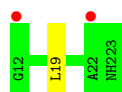
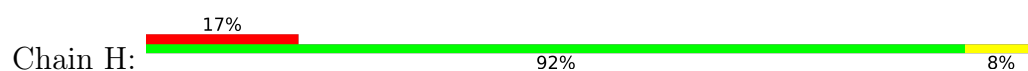
- Molecule 2: Aldehyde dehydrogenase, mitochondrial



- Molecule 2: Aldehyde dehydrogenase, mitochondrial



- Molecule 2: Aldehyde dehydrogenase, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	99.60Å 99.60Å 195.18Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.94 – 2.10 28.94 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.1 (28.94-2.10) 99.0 (28.94-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.78 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.243 , 0.287 0.243 , 0.286	Depositor DCC
R_{free} test set	1127 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	35.0	Xtriage
Anisotropy	0.109	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 35.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2488	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 44.10 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.6029e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: NH2, DCY

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	1.53	3/527 (0.6%)	1.22	1/713 (0.1%)
1	C	1.21	0/527	1.30	1/713 (0.1%)
1	E	1.53	2/543 (0.4%)	1.40	6/735 (0.8%)
1	G	1.15	2/543 (0.4%)	1.27	3/735 (0.4%)
2	B	1.14	0/78	1.10	0/101
2	D	1.23	0/78	1.01	0/101
2	F	1.38	0/78	1.26	0/101
2	H	1.06	0/78	1.17	0/101
All	All	1.35	7/2452 (0.3%)	1.28	11/3300 (0.3%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	118	VAL	CA-CB	-7.85	1.45	1.54
1	G	116	PRO	CA-C	6.74	1.58	1.52
1	E	110	LEU	C-O	-6.06	1.17	1.24
1	A	108	GLN	CA-C	-5.73	1.45	1.52
1	A	99	VAL	CA-CB	5.37	1.61	1.54
1	G	65	ALA	CA-CB	5.20	1.61	1.53
1	A	105	GLN	C-O	5.09	1.30	1.24

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	115	PRO	CA-C-N	-7.81	112.33	120.38
1	G	115	PRO	C-N-CA	-7.81	112.33	120.38
1	E	112	GLN	N-CA-C	-6.36	105.55	113.38
1	C	107	LEU	N-CA-C	6.01	117.83	111.28
1	E	116	PRO	N-CA-C	5.99	118.01	110.70
1	G	116	PRO	N-CA-C	5.81	117.78	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	85	ASP	CA-C-N	5.62	128.27	120.63
1	E	85	ASP	C-N-CA	5.62	128.27	120.63
1	A	71	LEU	N-CA-C	5.62	117.40	111.28
1	E	116	PRO	CA-C-N	-5.29	113.58	119.19
1	E	116	PRO	C-N-CA	-5.29	113.58	119.19

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	519	0	527	12	1
1	C	519	0	527	10	0
1	E	535	0	542	11	0
1	G	535	0	542	13	0
2	B	86	0	87	2	0
2	D	86	0	88	2	0
2	F	86	0	86	1	0
2	H	86	0	87	1	0
3	A	6	0	0	0	0
3	C	10	0	0	1	0
3	E	13	0	0	2	0
3	F	2	0	0	0	0
3	G	3	0	0	0	0
3	H	2	0	0	0	0
All	All	2488	0	2486	45	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 9.

All (45) 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:E:91:ASP:OD1	1:E:125:LYS:HE2	1.27	1.27

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:59:ASP:HB2	3:E:8:HOH:O	1.79	0.82
1:A:80:LEU:HD13	1:A:88:LYS:HG2	1.61	0.82
1:G:86:TYR:O	1:G:90:VAL:HG23	1.82	0.80
1:C:111:GLN:NE2	1:C:112:GLN:HE21	1.79	0.80
1:A:75:GLN:O	1:A:79:GLU:HG3	1.83	0.79
1:E:91:ASP:OD1	1:E:125:LYS:CE	2.22	0.78
1:E:120:GLN:OE1	3:E:29:HOH:O	2.07	0.73
1:A:109:VAL:O	1:A:113:THR:HG23	1.92	0.69
1:A:66:VAL:HG21	1:E:66:VAL:HG21	1.75	0.69
1:E:126:LEU:OXT	1:E:126:LEU:HG	1.96	0.64
1:A:72:GLU:O	1:A:76:LEU:HG	1.98	0.64
1:C:106:LEU:HD12	2:D:22:ALA:HB3	1.82	0.61
1:G:109:VAL:O	1:G:113:THR:HG22	2.01	0.60
1:G:110:LEU:HA	1:G:113:THR:HG23	1.83	0.60
1:C:76:LEU:HD23	1:C:92:HIS:CE1	2.37	0.60
1:C:68:LYS:NZ	1:C:72:GLU:OE2	2.35	0.59
1:C:111:GLN:HE22	1:C:112:GLN:HE21	1.51	0.58
1:G:116:PRO:N	1:G:117:PRO:HD2	2.19	0.58
1:G:77:GLY:HA3	1:G:93:LEU:CD1	2.34	0.57
1:G:64:GLU:O	1:G:68:LYS:HG2	2.05	0.56
2:B:15:LEU:O	2:B:19:LEU:HD13	2.09	0.52
1:G:109:VAL:O	1:G:113:THR:CG2	2.58	0.52
1:A:66:VAL:CG2	1:E:66:VAL:HG21	2.40	0.50
1:G:90:VAL:CG1	1:G:122:LEU:HA	2.41	0.50
1:E:126:LEU:OXT	1:E:126:LEU:CG	2.56	0.49
1:E:109:VAL:O	1:E:113:THR:HG23	2.12	0.49
1:E:78:GLU:OE2	2:F:12:GLY:O	2.32	0.48
1:A:108:GLN:O	1:A:112:GLN:HG2	2.15	0.47
1:E:122:LEU:C	1:E:122:LEU:HD23	2.39	0.47
1:C:102:GLN:HB3	1:C:105:GLN:OE1	2.16	0.46
1:C:110:LEU:HD21	2:D:19:LEU:HD11	1.96	0.46
1:G:115:PRO:C	1:G:117:PRO:HD2	2.40	0.45
1:A:120:GLN:O	1:A:124:THR:HG23	2.16	0.45
1:A:87:GLU:CG	1:A:125:LYS:HE3	2.47	0.45
1:G:76:LEU:O	1:G:77:GLY:C	2.59	0.44
1:C:116:PRO:N	1:C:117:PRO:HD2	2.33	0.43
1:G:90:VAL:HG11	1:G:122:LEU:HA	1.99	0.43
1:C:111:GLN:HG3	3:C:6:HOH:O	2.16	0.43
1:G:116:PRO:O	1:G:120:GLN:NE2	2.52	0.43
1:G:110:LEU:HG	2:H:19:LEU:HD21	2.01	0.43
1:A:80:LEU:CD1	1:A:88:LYS:HG2	2.42	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:VAL:O	1:A:94:THR:HG23	2.19	0.42
1:A:71:LEU:HD22	2:B:16:CYS:HB3	2.02	0.41
1:C:115:PRO:HA	1:C:116:PRO:HD3	1.92	0.41

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:85:ASP:OD2	1:A:125:LYS:NZ[2_565]	2.19	0.01

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	64/73 (88%)	64 (100%)	0	0	100	100
1	C	64/73 (88%)	64 (100%)	0	0	100	100
1	E	66/73 (90%)	66 (100%)	0	0	100	100
1	G	66/73 (90%)	64 (97%)	2 (3%)	0	100	100
2	B	9/12 (75%)	9 (100%)	0	0	100	100
2	D	9/12 (75%)	9 (100%)	0	0	100	100
2	F	9/12 (75%)	9 (100%)	0	0	100	100
2	H	9/12 (75%)	9 (100%)	0	0	100	100
All	All	296/340 (87%)	294 (99%)	2 (1%)	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	57/62 (92%)	55 (96%)	2 (4%)	32	35
1	C	57/62 (92%)	51 (90%)	6 (10%)	6	4
1	E	59/62 (95%)	52 (88%)	7 (12%)	5	3
1	G	59/62 (95%)	55 (93%)	4 (7%)	14	12
2	B	8/8 (100%)	8 (100%)	0	100	100
2	D	8/8 (100%)	8 (100%)	0	100	100
2	F	8/8 (100%)	8 (100%)	0	100	100
2	H	8/8 (100%)	8 (100%)	0	100	100
All	All	264/280 (94%)	245 (93%)	19 (7%)	13	11

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	LYS
1	A	126	LEU
1	C	61	LYS
1	C	64	GLU
1	C	68	LYS
1	C	102	GLN
1	C	106	LEU
1	C	111	GLN
1	E	60	LEU
1	E	68	LYS
1	E	76	LEU
1	E	87	GLU
1	E	100	SER
1	E	111	GLN
1	E	120	GLN
1	G	60	LEU
1	G	80	LEU
1	G	87	GLU
1	G	113	THR

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	105	GLN
1	C	111	GLN
1	E	120	GLN
1	G	120	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](#)

4 non-standard protein/DNA/RNA residues are modelled in this entry.

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.

5.5 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	66/73 (90%)	0.54	6 (9%) 15 15	23, 33, 43, 63	0
1	C	66/73 (90%)	0.79	7 (10%) 11 12	27, 38, 55, 66	0
1	E	68/73 (93%)	0.58	4 (5%) 28 30	22, 33, 53, 59	0
1	G	68/73 (93%)	1.64	22 (32%) 1 1	32, 48, 65, 69	0
2	B	10/12 (83%)	1.57	4 (40%) 1 1	41, 47, 56, 56	0
2	D	10/12 (83%)	1.09	1 (10%) 12 13	33, 38, 46, 47	0
2	F	10/12 (83%)	0.96	2 (20%) 3 3	27, 32, 40, 49	0
2	H	10/12 (83%)	1.22	2 (20%) 3 3	35, 40, 48, 51	0
All	All	308/340 (90%)	0.93	48 (15%) 5 5	22, 38, 59, 69	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	126	LEU	4.8
1	G	123	LEU	4.4
1	G	115	PRO	4.3
1	G	117	PRO	4.3
1	A	63	ALA	4.2
1	C	63	ALA	3.9
2	F	12	GLY	3.9
1	E	82	ALA	3.8
1	C	62	ASP	3.7
1	G	80	LEU	3.6
1	G	114	LEU	3.5
1	G	86	TYR	3.4
1	G	79	GLU	3.3
1	G	124	THR	3.3
1	G	116	PRO	3.0
2	F	21	TYR	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	61	LYS	2.9
1	G	82	ALA	2.9
2	H	22	ALA	2.8
2	H	12	GLY	2.8
1	G	118	VAL	2.8
1	G	106	LEU	2.6
1	A	105	GLN	2.6
1	G	88	LYS	2.6
1	G	83	GLN	2.5
2	B	21	TYR	2.5
1	E	86	TYR	2.4
1	G	109	VAL	2.4
1	C	124	THR	2.4
2	B	12	GLY	2.4
1	C	103	PRO	2.4
1	C	102	GLN	2.3
1	G	77	GLY	2.3
1	G	90	VAL	2.3
1	G	121	MET	2.3
1	E	81	LEU	2.3
1	A	104	GLN	2.2
2	B	20	SER	2.2
1	C	61	LYS	2.2
2	D	12	GLY	2.2
1	G	87	GLU	2.2
2	B	16	CYS	2.2
1	E	85	ASP	2.2
1	C	82	ALA	2.1
1	A	103	PRO	2.1
1	A	82	ALA	2.1
1	G	102	GLN	2.1
1	G	108	GLN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
2	DCY	B	13	6/7	0.83	0.15	49,50,51,51	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	DCY	H	13	6/7	0.84	0.14	43,46,48,51	0
2	DCY	F	13	6/7	0.90	0.15	36,38,40,45	0
2	DCY	D	13	6/7	0.90	0.11	38,42,43,44	0

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.