



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:35 PM UTC

PDB ID : 4ITV / pdb_00004itv
Title : Structure of a 16 nm protein cage designed by fusing symmetric oligomeric domains, triple mutant, P212121 form
Authors : Lai, Y.-T.; Sawaya, M.R.; Yeates, T.O.
Deposited on : 2013-01-18
Resolution : 3.60 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

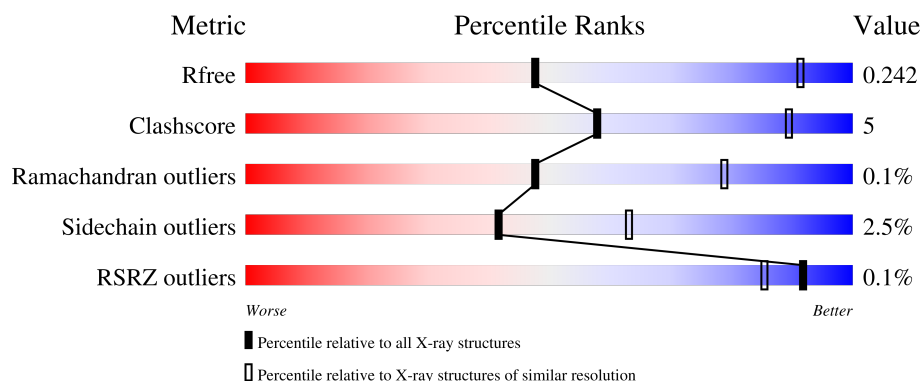
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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	1747 (3.70-3.50)
Clashscore	190562	1827 (3.70-3.50)
Ramachandran outliers	187476	1773 (3.70-3.50)
Sidechain outliers	187428	1772 (3.70-3.50)
RSRZ outliers	180081	1745 (3.70-3.50)

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	456	 81% 15% .
1	B	456	 83% 12% . .
1	C	456	 85% 11% . .
1	D	456	 84% 11% . .
1	E	456	 84% 12% . .

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	456	 84%12%.
1	G	456	 85%11%.
1	H	456	 83%12%. .
1	I	456	 85%11%.
1	J	456	 84%12%.
1	K	456	 84%12%.
1	L	456	 83%13%..

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 40776 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 Non-haem bromoperoxidase BPO-A2, Matrix protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	0	0	0
			3398	2166	572	652	8			
1	B	440	Total	C	N	O	S	0	0	0
			3398	2166	572	652	8			
1	C	440	Total	C	N	O	S	0	0	0
			3398	2166	572	652	8			
1	D	440	Total	C	N	O	S	0	0	0
			3398	2166	572	652	8			
1	E	440	Total	C	N	O	S	0	0	0
			3398	2166	572	652	8			
1	F	440	Total	C	N	O	S	0	0	0
			3398	2166	572	652	8			
1	G	440	Total	C	N	O	S	0	0	0
			3398	2166	572	652	8			
1	H	440	Total	C	N	O	S	0	0	0
			3398	2166	572	652	8			
1	I	440	Total	C	N	O	S	0	0	0
			3398	2166	572	652	8			
1	J	440	Total	C	N	O	S	0	0	0
			3398	2166	572	652	8			
1	K	440	Total	C	N	O	S	0	0	0
			3398	2166	572	652	8			
1	L	440	Total	C	N	O	S	0	0	0
			3398	2166	572	652	8			

There are 216 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	THR	GLN	engineered mutation	UNP P29715
A	118	ALA	LYS	engineered mutation	UNP P29715
A	278	ALA	-	linker	UNP P03485
A	279	GLN	-	linker	UNP P03485
A	280	GLU	-	linker	UNP P03485

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	281	ALA	-	linker	UNP P03485
A	282	GLN	-	linker	UNP P03485
A	283	LYS	-	linker	UNP P03485
A	284	GLN	-	linker	UNP P03485
A	285	LYS	-	linker	UNP P03485
A	448	LEU	-	expression tag	UNP P03485
A	449	GLU	-	expression tag	UNP P03485
A	450	HIS	-	expression tag	UNP P03485
A	451	HIS	-	expression tag	UNP P03485
A	452	HIS	-	expression tag	UNP P03485
A	453	HIS	-	expression tag	UNP P03485
A	454	HIS	-	expression tag	UNP P03485
A	455	HIS	-	expression tag	UNP P03485
B	24	THR	GLN	engineered mutation	UNP P29715
B	118	ALA	LYS	engineered mutation	UNP P29715
B	278	ALA	-	linker	UNP P03485
B	279	GLN	-	linker	UNP P03485
B	280	GLU	-	linker	UNP P03485
B	281	ALA	-	linker	UNP P03485
B	282	GLN	-	linker	UNP P03485
B	283	LYS	-	linker	UNP P03485
B	284	GLN	-	linker	UNP P03485
B	285	LYS	-	linker	UNP P03485
B	448	LEU	-	expression tag	UNP P03485
B	449	GLU	-	expression tag	UNP P03485
B	450	HIS	-	expression tag	UNP P03485
B	451	HIS	-	expression tag	UNP P03485
B	452	HIS	-	expression tag	UNP P03485
B	453	HIS	-	expression tag	UNP P03485
B	454	HIS	-	expression tag	UNP P03485
B	455	HIS	-	expression tag	UNP P03485
C	24	THR	GLN	engineered mutation	UNP P29715
C	118	ALA	LYS	engineered mutation	UNP P29715
C	278	ALA	-	linker	UNP P03485
C	279	GLN	-	linker	UNP P03485
C	280	GLU	-	linker	UNP P03485
C	281	ALA	-	linker	UNP P03485
C	282	GLN	-	linker	UNP P03485
C	283	LYS	-	linker	UNP P03485
C	284	GLN	-	linker	UNP P03485
C	285	LYS	-	linker	UNP P03485
C	448	LEU	-	expression tag	UNP P03485

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	449	GLU	-	expression tag	UNP P03485
C	450	HIS	-	expression tag	UNP P03485
C	451	HIS	-	expression tag	UNP P03485
C	452	HIS	-	expression tag	UNP P03485
C	453	HIS	-	expression tag	UNP P03485
C	454	HIS	-	expression tag	UNP P03485
C	455	HIS	-	expression tag	UNP P03485
D	24	THR	GLN	engineered mutation	UNP P29715
D	118	ALA	LYS	engineered mutation	UNP P29715
D	278	ALA	-	linker	UNP P03485
D	279	GLN	-	linker	UNP P03485
D	280	GLU	-	linker	UNP P03485
D	281	ALA	-	linker	UNP P03485
D	282	GLN	-	linker	UNP P03485
D	283	LYS	-	linker	UNP P03485
D	284	GLN	-	linker	UNP P03485
D	285	LYS	-	linker	UNP P03485
D	448	LEU	-	expression tag	UNP P03485
D	449	GLU	-	expression tag	UNP P03485
D	450	HIS	-	expression tag	UNP P03485
D	451	HIS	-	expression tag	UNP P03485
D	452	HIS	-	expression tag	UNP P03485
D	453	HIS	-	expression tag	UNP P03485
D	454	HIS	-	expression tag	UNP P03485
D	455	HIS	-	expression tag	UNP P03485
E	24	THR	GLN	engineered mutation	UNP P29715
E	118	ALA	LYS	engineered mutation	UNP P29715
E	278	ALA	-	linker	UNP P03485
E	279	GLN	-	linker	UNP P03485
E	280	GLU	-	linker	UNP P03485
E	281	ALA	-	linker	UNP P03485
E	282	GLN	-	linker	UNP P03485
E	283	LYS	-	linker	UNP P03485
E	284	GLN	-	linker	UNP P03485
E	285	LYS	-	linker	UNP P03485
E	448	LEU	-	expression tag	UNP P03485
E	449	GLU	-	expression tag	UNP P03485
E	450	HIS	-	expression tag	UNP P03485
E	451	HIS	-	expression tag	UNP P03485
E	452	HIS	-	expression tag	UNP P03485
E	453	HIS	-	expression tag	UNP P03485
E	454	HIS	-	expression tag	UNP P03485

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	455	HIS	-	expression tag	UNP P03485
F	24	THR	GLN	engineered mutation	UNP P29715
F	118	ALA	LYS	engineered mutation	UNP P29715
F	278	ALA	-	linker	UNP P03485
F	279	GLN	-	linker	UNP P03485
F	280	GLU	-	linker	UNP P03485
F	281	ALA	-	linker	UNP P03485
F	282	GLN	-	linker	UNP P03485
F	283	LYS	-	linker	UNP P03485
F	284	GLN	-	linker	UNP P03485
F	285	LYS	-	linker	UNP P03485
F	448	LEU	-	expression tag	UNP P03485
F	449	GLU	-	expression tag	UNP P03485
F	450	HIS	-	expression tag	UNP P03485
F	451	HIS	-	expression tag	UNP P03485
F	452	HIS	-	expression tag	UNP P03485
F	453	HIS	-	expression tag	UNP P03485
F	454	HIS	-	expression tag	UNP P03485
F	455	HIS	-	expression tag	UNP P03485
G	24	THR	GLN	engineered mutation	UNP P29715
G	118	ALA	LYS	engineered mutation	UNP P29715
G	278	ALA	-	linker	UNP P03485
G	279	GLN	-	linker	UNP P03485
G	280	GLU	-	linker	UNP P03485
G	281	ALA	-	linker	UNP P03485
G	282	GLN	-	linker	UNP P03485
G	283	LYS	-	linker	UNP P03485
G	284	GLN	-	linker	UNP P03485
G	285	LYS	-	linker	UNP P03485
G	448	LEU	-	expression tag	UNP P03485
G	449	GLU	-	expression tag	UNP P03485
G	450	HIS	-	expression tag	UNP P03485
G	451	HIS	-	expression tag	UNP P03485
G	452	HIS	-	expression tag	UNP P03485
G	453	HIS	-	expression tag	UNP P03485
G	454	HIS	-	expression tag	UNP P03485
G	455	HIS	-	expression tag	UNP P03485
H	24	THR	GLN	engineered mutation	UNP P29715
H	118	ALA	LYS	engineered mutation	UNP P29715
H	278	ALA	-	linker	UNP P03485
H	279	GLN	-	linker	UNP P03485
H	280	GLU	-	linker	UNP P03485

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
H	281	ALA	-	linker	UNP P03485
H	282	GLN	-	linker	UNP P03485
H	283	LYS	-	linker	UNP P03485
H	284	GLN	-	linker	UNP P03485
H	285	LYS	-	linker	UNP P03485
H	448	LEU	-	expression tag	UNP P03485
H	449	GLU	-	expression tag	UNP P03485
H	450	HIS	-	expression tag	UNP P03485
H	451	HIS	-	expression tag	UNP P03485
H	452	HIS	-	expression tag	UNP P03485
H	453	HIS	-	expression tag	UNP P03485
H	454	HIS	-	expression tag	UNP P03485
H	455	HIS	-	expression tag	UNP P03485
I	24	THR	GLN	engineered mutation	UNP P29715
I	118	ALA	LYS	engineered mutation	UNP P29715
I	278	ALA	-	linker	UNP P03485
I	279	GLN	-	linker	UNP P03485
I	280	GLU	-	linker	UNP P03485
I	281	ALA	-	linker	UNP P03485
I	282	GLN	-	linker	UNP P03485
I	283	LYS	-	linker	UNP P03485
I	284	GLN	-	linker	UNP P03485
I	285	LYS	-	linker	UNP P03485
I	448	LEU	-	expression tag	UNP P03485
I	449	GLU	-	expression tag	UNP P03485
I	450	HIS	-	expression tag	UNP P03485
I	451	HIS	-	expression tag	UNP P03485
I	452	HIS	-	expression tag	UNP P03485
I	453	HIS	-	expression tag	UNP P03485
I	454	HIS	-	expression tag	UNP P03485
I	455	HIS	-	expression tag	UNP P03485
J	24	THR	GLN	engineered mutation	UNP P29715
J	118	ALA	LYS	engineered mutation	UNP P29715
J	278	ALA	-	linker	UNP P03485
J	279	GLN	-	linker	UNP P03485
J	280	GLU	-	linker	UNP P03485
J	281	ALA	-	linker	UNP P03485
J	282	GLN	-	linker	UNP P03485
J	283	LYS	-	linker	UNP P03485
J	284	GLN	-	linker	UNP P03485
J	285	LYS	-	linker	UNP P03485
J	448	LEU	-	expression tag	UNP P03485

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
J	449	GLU	-	expression tag	UNP P03485
J	450	HIS	-	expression tag	UNP P03485
J	451	HIS	-	expression tag	UNP P03485
J	452	HIS	-	expression tag	UNP P03485
J	453	HIS	-	expression tag	UNP P03485
J	454	HIS	-	expression tag	UNP P03485
J	455	HIS	-	expression tag	UNP P03485
K	24	THR	GLN	engineered mutation	UNP P29715
K	118	ALA	LYS	engineered mutation	UNP P29715
K	278	ALA	-	linker	UNP P03485
K	279	GLN	-	linker	UNP P03485
K	280	GLU	-	linker	UNP P03485
K	281	ALA	-	linker	UNP P03485
K	282	GLN	-	linker	UNP P03485
K	283	LYS	-	linker	UNP P03485
K	284	GLN	-	linker	UNP P03485
K	285	LYS	-	linker	UNP P03485
K	448	LEU	-	expression tag	UNP P03485
K	449	GLU	-	expression tag	UNP P03485
K	450	HIS	-	expression tag	UNP P03485
K	451	HIS	-	expression tag	UNP P03485
K	452	HIS	-	expression tag	UNP P03485
K	453	HIS	-	expression tag	UNP P03485
K	454	HIS	-	expression tag	UNP P03485
K	455	HIS	-	expression tag	UNP P03485
L	24	THR	GLN	engineered mutation	UNP P29715
L	118	ALA	LYS	engineered mutation	UNP P29715
L	278	ALA	-	linker	UNP P03485
L	279	GLN	-	linker	UNP P03485
L	280	GLU	-	linker	UNP P03485
L	281	ALA	-	linker	UNP P03485
L	282	GLN	-	linker	UNP P03485
L	283	LYS	-	linker	UNP P03485
L	284	GLN	-	linker	UNP P03485
L	285	LYS	-	linker	UNP P03485
L	448	LEU	-	expression tag	UNP P03485
L	449	GLU	-	expression tag	UNP P03485
L	450	HIS	-	expression tag	UNP P03485
L	451	HIS	-	expression tag	UNP P03485
L	452	HIS	-	expression tag	UNP P03485
L	453	HIS	-	expression tag	UNP P03485
L	454	HIS	-	expression tag	UNP P03485

Continued on next page...

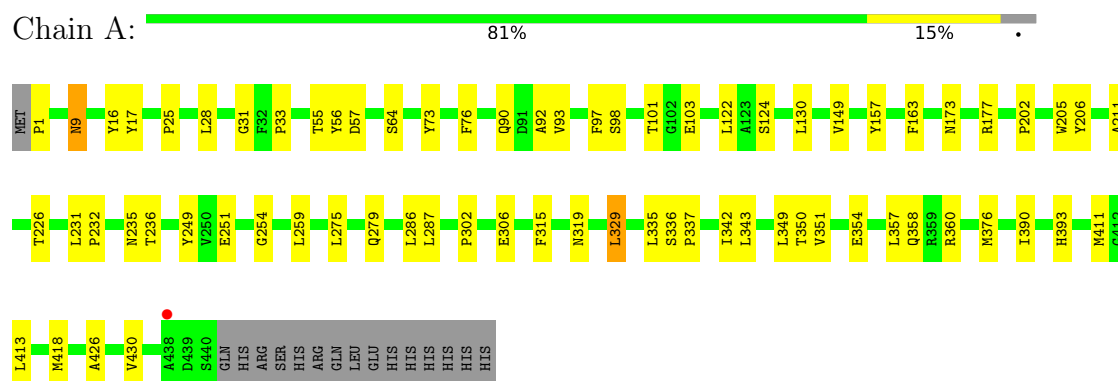
Continued from previous page...

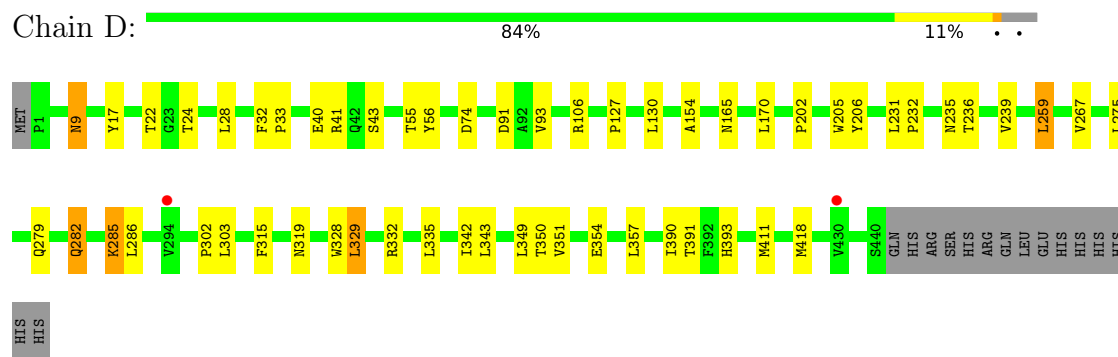
Chain	Residue	Modelled	Actual	Comment	Reference
L	455	HIS	-	expression tag	UNP P03485

3 Residue-property plots

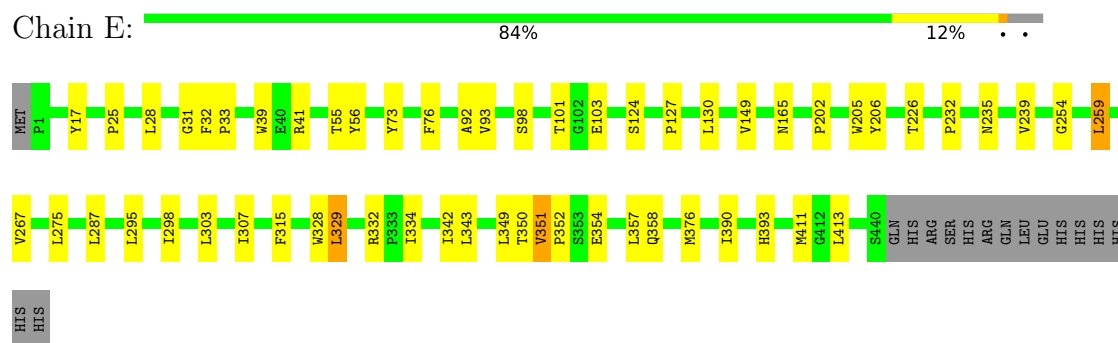
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Non-haem bromoperoxidase BPO-A2, Matrix protein 1

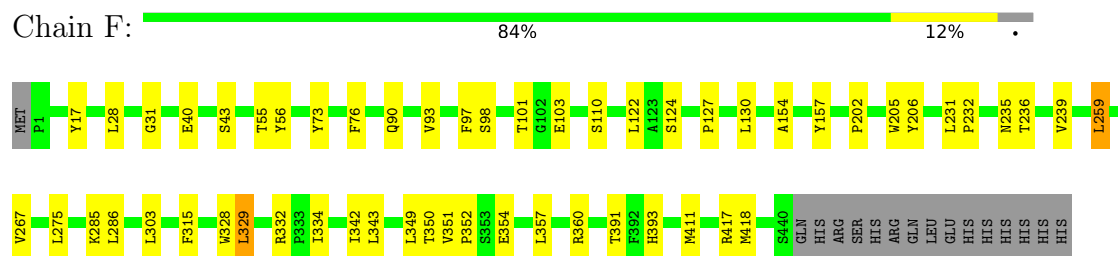




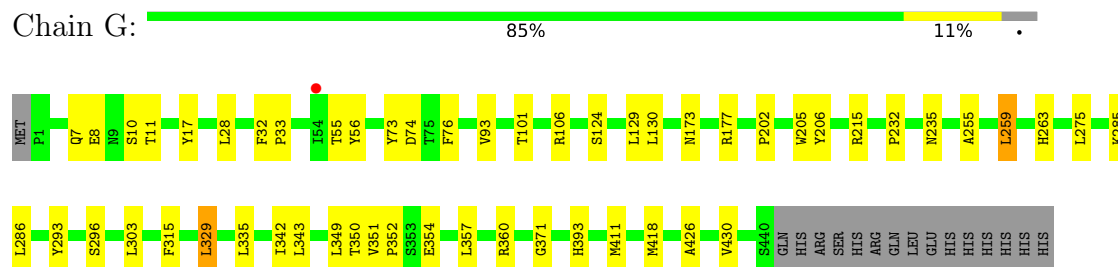
- Molecule 1: Non-haem bromoperoxidase BPO-A2, Matrix protein 1



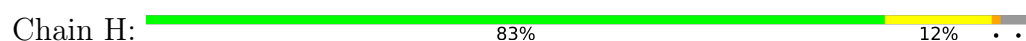
- Molecule 1: Non-haem bromoperoxidase BPO-A2, Matrix protein 1



- Molecule 1: Non-haem bromoperoxidase BPO-A2, Matrix protein 1



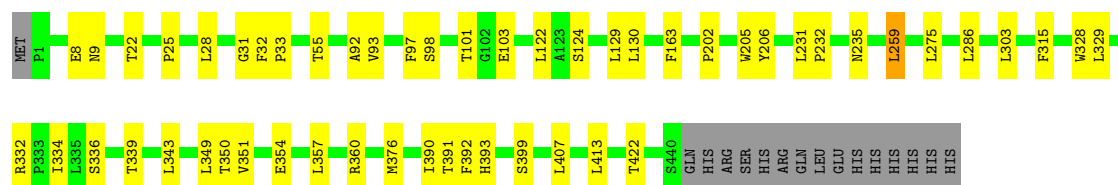
- Molecule 1: Non-haem bromoperoxidase BPO-A2, Matrix protein 1





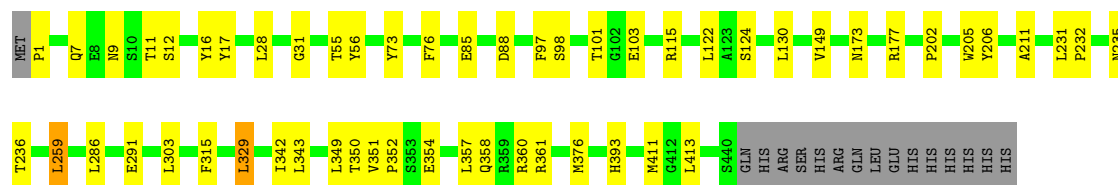
- Molecule 1: Non-haem bromoperoxidase BPO-A2, Matrix protein 1

Chain I: .



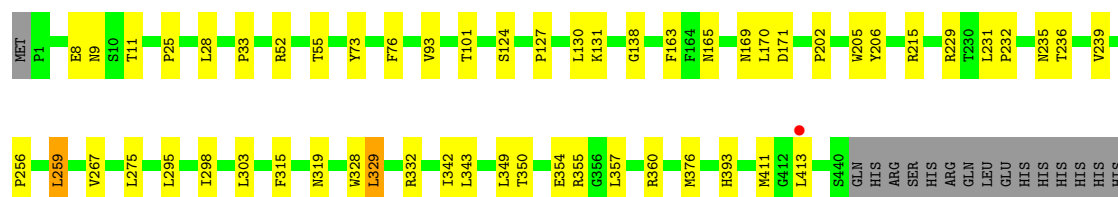
- Molecule 1: Non-haem bromoperoxidase BPO-A2, Matrix protein 1

Chain J: .



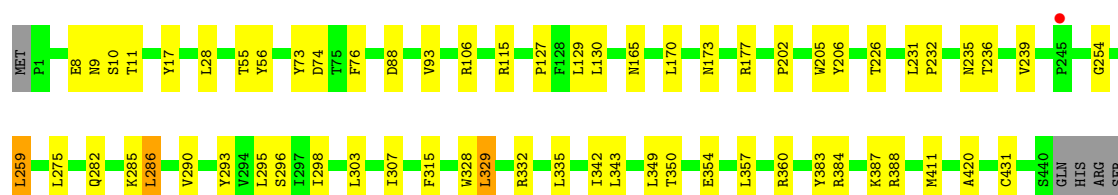
- Molecule 1: Non-haem bromoperoxidase BPO-A2, Matrix protein 1

Chain K: .



- Molecule 1: Non-haem bromoperoxidase BPO-A2, Matrix protein 1

Chain L: .



HIS
ARG
GLN
LEU
GLU
HIS
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	131.62Å 156.61Å 321.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	96.14 – 3.60 96.14 – 3.60	Depositor EDS
% Data completeness (in resolution range)	98.3 (96.14-3.60) 98.3 (96.14-3.60)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 3.58Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, R_{free}	0.206 , 0.240 0.208 , 0.242	Depositor DCC
R_{free} test set	3824 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	114.8	Xtriage
Anisotropy	0.164	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 99.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	40776	wwPDB-VP
Average B, all atoms (Å ²)	151.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.33% 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	0.31	0/3475	0.73	2/4729 (0.0%)
1	B	0.30	0/3475	0.74	3/4729 (0.1%)
1	C	0.30	0/3475	0.73	0/4729
1	D	0.29	0/3475	0.72	2/4729 (0.0%)
1	E	0.30	0/3475	0.71	2/4729 (0.0%)
1	F	0.29	0/3475	0.73	0/4729
1	G	0.30	0/3475	0.73	0/4729
1	H	0.29	0/3475	0.73	2/4729 (0.0%)
1	I	0.29	0/3475	0.72	2/4729 (0.0%)
1	J	0.30	0/3475	0.73	0/4729
1	K	0.30	0/3475	0.73	0/4729
1	L	0.29	0/3475	0.73	0/4729
All	All	0.30	0/41700	0.73	13/56748 (0.0%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	9	ASN	CB-CA-C	-6.03	109.60	116.54
1	H	301	GLY	CA-C-N	5.80	125.42	119.56
1	H	301	GLY	C-N-CA	5.80	125.42	119.56
1	B	351	VAL	CA-C-N	5.65	125.76	119.32
1	B	351	VAL	C-N-CA	5.65	125.76	119.32
1	A	351	VAL	CA-C-N	5.44	124.63	118.97
1	A	351	VAL	C-N-CA	5.44	124.63	118.97
1	D	351	VAL	CA-C-N	5.27	124.90	119.05
1	D	351	VAL	C-N-CA	5.27	124.90	119.05
1	E	351	VAL	CA-C-N	5.22	125.27	119.32
1	E	351	VAL	C-N-CA	5.22	125.27	119.32
1	I	351	VAL	CA-C-N	5.17	125.22	119.32
1	I	351	VAL	C-N-CA	5.17	125.22	119.32

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	3398	0	3326	42	0
1	B	3398	0	3326	38	0
1	C	3398	0	3326	29	0
1	D	3398	0	3326	34	0
1	E	3398	0	3326	33	0
1	F	3398	0	3326	29	0
1	G	3398	0	3326	33	0
1	H	3398	0	3326	34	0
1	I	3398	0	3326	29	0
1	J	3398	0	3326	29	0
1	K	3398	0	3326	31	0
1	L	3398	0	3326	40	0
All	All	40776	0	39912	365	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 (365) 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:130:LEU:HB2	1:E:206:TYR:HB2	1.70	0.73
1:A:90:GLN:HB2	1:A:390:ILE:HD13	1.71	0.73
1:L:387:LYS:NZ	1:L:420:ALA:O	2.22	0.72
1:B:130:LEU:HB2	1:B:206:TYR:HB2	1.71	0.72
1:B:1:PRO:HD2	1:B:16:TYR:HE1	1.57	0.69
1:D:9:ASN:HA	1:E:390:ILE:HD11	1.73	0.69
1:C:23:GLY:HA2	1:C:390:ILE:HD13	1.74	0.69
1:F:417:ARG:HD2	1:L:387:LYS:HE2	1.74	0.68
1:K:130:LEU:HB2	1:K:206:TYR:HB2	1.75	0.68
1:H:130:LEU:HB2	1:H:206:TYR:HB2	1.74	0.68
1:F:28:LEU:HB2	1:F:55:THR:HG22	1.76	0.68
1:G:130:LEU:HB2	1:G:206:TYR:HB2	1.75	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:130:LEU:HB2	1:C:206:TYR:HB2	1.77	0.66
1:A:360:ARG:HH21	1:J:358:GLN:HG3	1.60	0.66
1:B:1:PRO:HD2	1:B:16:TYR:CE1	2.29	0.66
1:K:28:LEU:HB2	1:K:55:THR:HG22	1.77	0.66
1:C:315:PHE:HE1	1:C:350:THR:HG21	1.60	0.66
1:B:329:LEU:HD11	1:B:343:LEU:HB3	1.79	0.65
1:J:130:LEU:HB2	1:J:206:TYR:HB2	1.78	0.65
1:F:418:MET:HE3	1:L:388:ARG:NH1	2.10	0.65
1:A:28:LEU:HB2	1:A:55:THR:HG22	1.78	0.65
1:L:130:LEU:HB2	1:L:206:TYR:HB2	1.79	0.64
1:L:8:GLU:HG2	1:L:9:ASN:H	1.62	0.64
1:D:302:PRO:HD2	1:G:215:ARG:HH22	1.63	0.64
1:K:315:PHE:HE1	1:K:350:THR:HG21	1.63	0.63
1:D:22:THR:HA	1:D:390:ILE:HD11	1.79	0.63
1:I:130:LEU:HB2	1:I:206:TYR:HB2	1.81	0.63
1:D:130:LEU:HB2	1:D:206:TYR:HB2	1.80	0.63
1:K:8:GLU:HG2	1:K:9:ASN:H	1.64	0.63
1:D:91:ASP:OD2	1:D:285:LYS:NZ	2.33	0.62
1:C:232:PRO:HG2	1:C:235:ASN:HB2	1.82	0.61
1:I:8:GLU:HG2	1:I:9:ASN:H	1.65	0.61
1:K:319:ASN:OD1	1:K:355:ARG:NH2	2.33	0.61
1:A:376:MET:HE3	1:J:413:LEU:HD11	1.82	0.61
1:I:28:LEU:HB2	1:I:55:THR:HG22	1.82	0.61
1:D:231:LEU:HB3	1:D:236:THR:HG21	1.81	0.61
1:K:171:ASP:OD1	1:K:229:ARG:NH1	2.34	0.61
1:J:232:PRO:HG2	1:J:235:ASN:HB2	1.82	0.61
1:G:315:PHE:HE1	1:G:350:THR:HG21	1.66	0.61
1:L:28:LEU:HB2	1:L:55:THR:HG22	1.83	0.60
1:D:232:PRO:HG2	1:D:235:ASN:HB2	1.83	0.60
1:K:232:PRO:HG2	1:K:235:ASN:HB2	1.84	0.60
1:L:74:ASP:OD1	1:L:106:ARG:NH1	2.34	0.60
1:D:418:MET:HE3	1:H:417:ARG:HD3	1.84	0.60
1:E:28:LEU:HB2	1:E:55:THR:HG22	1.84	0.60
1:L:232:PRO:HG2	1:L:235:ASN:HB2	1.83	0.60
1:C:342:ILE:HA	1:C:411:MET:HE1	1.82	0.59
1:L:315:PHE:HE1	1:L:350:THR:HG21	1.67	0.59
1:A:358:GLN:HG3	1:J:360:ARG:HH21	1.67	0.58
1:B:93:VAL:HG11	1:B:275:LEU:HD21	1.85	0.58
1:I:232:PRO:HG2	1:I:235:ASN:HB2	1.85	0.58
1:K:93:VAL:HG11	1:K:275:LEU:HD21	1.85	0.58
1:B:232:PRO:HG2	1:B:235:ASN:HB2	1.85	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:LEU:HB2	1:A:206:TYR:HB2	1.87	0.57
1:B:101:THR:HG21	1:B:124:SER:HA	1.87	0.57
1:E:93:VAL:HG11	1:E:275:LEU:HD21	1.86	0.57
1:D:28:LEU:HB2	1:D:55:THR:HG22	1.85	0.57
1:E:315:PHE:HE1	1:E:350:THR:HG21	1.69	0.56
1:G:232:PRO:HG2	1:G:235:ASN:HB2	1.87	0.56
1:F:232:PRO:HG2	1:F:235:ASN:HB2	1.87	0.56
1:L:226:THR:HG1	1:L:254:GLY:H	1.54	0.56
1:J:7:GLN:HA	1:J:12:SER:HA	1.86	0.56
1:B:376:MET:HE3	1:E:413:LEU:HD11	1.86	0.56
1:B:74:ASP:OD1	1:B:106:ARG:NH1	2.39	0.56
1:F:334:ILE:HD13	1:K:215:ARG:HH21	1.71	0.56
1:L:342:ILE:HA	1:L:411:MET:HE1	1.87	0.56
1:H:93:VAL:HG11	1:H:275:LEU:HD21	1.88	0.56
1:J:231:LEU:HB3	1:J:236:THR:HG21	1.86	0.56
1:K:165:ASN:HA	1:K:170:LEU:HD12	1.87	0.56
1:L:329:LEU:HD11	1:L:343:LEU:HB3	1.88	0.56
1:E:329:LEU:HD11	1:E:343:LEU:HB3	1.87	0.55
1:F:130:LEU:HB2	1:F:206:TYR:HB2	1.88	0.55
1:H:28:LEU:HB2	1:H:55:THR:HG22	1.89	0.55
1:G:28:LEU:HB2	1:G:55:THR:HG22	1.89	0.55
1:K:329:LEU:HD11	1:K:343:LEU:HB3	1.86	0.55
1:A:93:VAL:HG11	1:A:275:LEU:HD21	1.88	0.55
1:E:232:PRO:HG2	1:E:235:ASN:HB2	1.87	0.55
1:A:413:LEU:HD11	1:J:376:MET:HE3	1.87	0.55
1:A:211:ALA:HA	1:E:334:ILE:HA	1.90	0.55
1:E:226:THR:HG1	1:E:254:GLY:H	1.54	0.54
1:B:7:GLN:HB2	1:G:371:GLY:HA3	1.89	0.54
1:E:101:THR:HG21	1:E:124:SER:HA	1.90	0.54
1:E:73:TYR:HA	1:E:76:PHE:HB2	1.89	0.53
1:A:9:ASN:HA	1:B:390:ILE:HD11	1.90	0.53
1:C:384:ARG:HG3	1:G:418:MET:HG2	1.91	0.53
1:G:173:ASN:HB3	1:G:177:ARG:HB2	1.91	0.53
1:G:259:LEU:HD13	1:G:259:LEU:H	1.74	0.53
1:F:315:PHE:HE1	1:F:350:THR:HG21	1.73	0.53
1:H:101:THR:HG21	1:H:124:SER:HA	1.90	0.53
1:B:28:LEU:HB2	1:B:55:THR:HG22	1.89	0.53
1:D:202:PRO:HA	1:D:205:TRP:CE2	2.45	0.52
1:B:413:LEU:HD11	1:E:376:MET:HE3	1.91	0.52
1:L:259:LEU:H	1:L:259:LEU:HD13	1.73	0.52
1:D:329:LEU:HD11	1:D:343:LEU:HB3	1.91	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:232:PRO:HG2	1:H:235:ASN:HB2	1.90	0.52
1:I:413:LEU:HD11	1:K:376:MET:HE3	1.90	0.52
1:G:342:ILE:HA	1:G:411:MET:HE1	1.92	0.52
1:F:231:LEU:HB3	1:F:236:THR:HG21	1.92	0.52
1:C:28:LEU:HB2	1:C:55:THR:HG22	1.92	0.51
1:B:173:ASN:HB3	1:B:177:ARG:HB2	1.93	0.51
1:J:329:LEU:HD11	1:J:343:LEU:HB3	1.92	0.51
1:F:259:LEU:HD13	1:F:259:LEU:H	1.75	0.51
1:K:259:LEU:HD13	1:K:259:LEU:H	1.75	0.51
1:F:93:VAL:HG11	1:F:275:LEU:HD21	1.92	0.51
1:K:342:ILE:HA	1:K:411:MET:HE1	1.93	0.50
1:A:232:PRO:HG2	1:A:235:ASN:HB2	1.92	0.50
1:C:286:LEU:HD11	1:C:392:PHE:CD1	2.47	0.50
1:I:93:VAL:HG11	1:I:275:LEU:HD21	1.93	0.50
1:D:9:ASN:HA	1:E:390:ILE:CD1	2.41	0.50
1:A:31:GLY:HA3	1:A:98:SER:HB3	1.94	0.50
1:H:259:LEU:HD13	1:H:259:LEU:H	1.76	0.50
1:A:173:ASN:HB3	1:A:177:ARG:HB2	1.92	0.50
1:C:384:ARG:NE	1:G:418:MET:HA	2.26	0.50
1:D:40:GLU:OE2	1:F:157:TYR:OH	2.28	0.50
1:F:110:SER:O	1:H:365:ASN:HB3	2.12	0.50
1:L:73:TYR:HA	1:L:76:PHE:HB2	1.94	0.49
1:A:226:THR:HG1	1:A:254:GLY:H	1.57	0.49
1:G:74:ASP:OD1	1:G:106:ARG:NH1	2.45	0.49
1:B:157:TYR:OH	1:C:40:GLU:OE2	2.21	0.49
1:B:165:ASN:HA	1:B:170:LEU:HD12	1.94	0.49
1:J:17:TYR:HA	1:J:56:TYR:HA	1.95	0.49
1:B:73:TYR:HA	1:B:76:PHE:HB2	1.94	0.49
1:D:154:ALA:O	1:E:41:ARG:NH2	2.41	0.49
1:I:259:LEU:HD13	1:I:259:LEU:H	1.78	0.49
1:A:231:LEU:HB3	1:A:236:THR:HG21	1.93	0.48
1:G:329:LEU:HD11	1:G:343:LEU:HB3	1.95	0.48
1:L:93:VAL:HG11	1:L:275:LEU:HD21	1.95	0.48
1:B:342:ILE:HA	1:B:411:MET:HE1	1.94	0.48
1:C:101:THR:HG21	1:C:124:SER:HA	1.96	0.48
1:E:17:TYR:HA	1:E:56:TYR:HA	1.95	0.48
1:I:286:LEU:HD21	1:I:392:PHE:CG	2.48	0.48
1:J:28:LEU:HB2	1:J:55:THR:HG22	1.94	0.48
1:L:9:ASN:C	1:L:11:THR:H	2.21	0.48
1:I:315:PHE:HE1	1:I:350:THR:HG21	1.79	0.48
1:J:259:LEU:HD13	1:J:259:LEU:H	1.79	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:315:PHE:HE1	1:B:350:THR:HG21	1.78	0.48
1:F:417:ARG:CD	1:L:387:LYS:HE2	2.44	0.48
1:A:315:PHE:HE1	1:A:350:THR:HG21	1.78	0.48
1:J:85:GLU:OE1	1:J:115:ARG:NH2	2.44	0.48
1:J:315:PHE:HE1	1:J:350:THR:HG21	1.79	0.48
1:F:328:TRP:O	1:F:332:ARG:HG2	2.13	0.48
1:K:342:ILE:HG12	1:K:411:MET:HE1	1.96	0.48
1:A:342:ILE:HA	1:A:411:MET:HE1	1.95	0.48
1:C:259:LEU:HD13	1:C:259:LEU:H	1.78	0.48
1:D:259:LEU:HD13	1:D:259:LEU:H	1.78	0.48
1:I:202:PRO:HA	1:I:205:TRP:CE2	2.48	0.48
1:K:231:LEU:HB3	1:K:236:THR:HG21	1.96	0.48
1:D:390:ILE:HG23	1:D:391:THR:HG23	1.96	0.48
1:E:31:GLY:HA3	1:E:98:SER:HB3	1.95	0.48
1:G:342:ILE:HG12	1:G:411:MET:HE1	1.96	0.47
1:K:259:LEU:HD23	1:K:267:VAL:HG21	1.96	0.47
1:C:31:GLY:HA3	1:C:98:SER:HB3	1.95	0.47
1:F:342:ILE:HA	1:F:411:MET:HE1	1.95	0.47
1:I:25:PRO:HG2	1:I:92:ALA:HA	1.95	0.47
1:L:295:LEU:HD23	1:L:298:ILE:HD12	1.96	0.47
1:D:302:PRO:HD2	1:G:215:ARG:NH2	2.28	0.47
1:H:342:ILE:HA	1:H:411:MET:HE1	1.96	0.47
1:L:231:LEU:HB3	1:L:236:THR:HG21	1.97	0.47
1:J:173:ASN:HB3	1:J:177:ARG:HB2	1.96	0.47
1:G:202:PRO:HA	1:G:205:TRP:CD2	2.50	0.47
1:H:202:PRO:HA	1:H:205:TRP:CD2	2.50	0.47
1:B:1:PRO:HB2	1:B:17:TYR:O	2.15	0.47
1:B:17:TYR:HA	1:B:56:TYR:HA	1.96	0.47
1:H:202:PRO:HA	1:H:205:TRP:CE2	2.50	0.47
1:C:381:LYS:HG2	1:C:384:ARG:NH2	2.30	0.47
1:F:329:LEU:HD11	1:F:343:LEU:HB3	1.96	0.47
1:L:202:PRO:HA	1:L:205:TRP:CD2	2.50	0.47
1:G:32:PHE:HA	1:G:33:PRO:HA	1.75	0.46
1:L:290:VAL:HG22	1:L:431:CYS:SG	2.55	0.46
1:A:17:TYR:HA	1:A:56:TYR:HA	1.96	0.46
1:A:9:ASN:HA	1:B:390:ILE:CD1	2.45	0.46
1:A:202:PRO:HA	1:A:205:TRP:CE2	2.51	0.46
1:E:342:ILE:HG12	1:E:411:MET:HE1	1.97	0.46
1:K:73:TYR:HA	1:K:76:PHE:HB2	1.96	0.46
1:A:73:TYR:HA	1:A:76:PHE:HB2	1.97	0.46
1:J:149:VAL:HG22	1:J:202:PRO:HB2	1.97	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:73:TYR:HA	1:G:76:PHE:HB2	1.97	0.46
1:J:11:THR:HG22	1:J:12:SER:N	2.31	0.46
1:J:31:GLY:HA3	1:J:98:SER:HB3	1.97	0.46
1:C:93:VAL:HG11	1:C:275:LEU:HD21	1.98	0.46
1:K:202:PRO:HA	1:K:205:TRP:CE2	2.51	0.46
1:B:360:ARG:HE	1:B:360:ARG:HA	1.81	0.46
1:E:25:PRO:HG2	1:E:92:ALA:HA	1.96	0.46
1:K:328:TRP:O	1:K:332:ARG:HG2	2.16	0.46
1:L:88:ASP:HA	1:L:115:ARG:HH12	1.81	0.46
1:D:17:TYR:HA	1:D:56:TYR:HA	1.98	0.45
1:D:342:ILE:HA	1:D:411:MET:HE1	1.97	0.45
1:J:202:PRO:HA	1:J:205:TRP:CE2	2.50	0.45
1:L:165:ASN:HA	1:L:170:LEU:HD12	1.98	0.45
1:D:74:ASP:OD1	1:D:106:ARG:NH1	2.47	0.45
1:D:328:TRP:O	1:D:332:ARG:HG2	2.17	0.45
1:B:259:LEU:HD13	1:B:259:LEU:H	1.81	0.45
1:B:328:TRP:O	1:B:332:ARG:HG2	2.16	0.45
1:F:73:TYR:HA	1:F:76:PHE:HB2	1.98	0.45
1:K:101:THR:HG21	1:K:124:SER:HA	1.97	0.45
1:D:165:ASN:HA	1:D:170:LEU:HD12	1.98	0.45
1:H:329:LEU:HD11	1:H:343:LEU:HB3	1.97	0.45
1:A:202:PRO:HA	1:A:205:TRP:CD2	2.51	0.45
1:H:303:LEU:HD13	1:H:335:LEU:HG	1.99	0.45
1:G:129:LEU:HB2	1:G:206:TYR:HA	1.99	0.45
1:H:17:TYR:HA	1:H:56:TYR:HA	1.97	0.45
1:I:329:LEU:HD11	1:I:343:LEU:HB3	1.98	0.45
1:D:259:LEU:HD23	1:D:267:VAL:HG21	1.99	0.45
1:K:33:PRO:HD3	1:K:163:PHE:CE2	2.52	0.45
1:B:231:LEU:HB3	1:B:236:THR:HG21	1.98	0.45
1:D:32:PHE:HA	1:D:33:PRO:HA	1.81	0.45
1:D:41:ARG:NH2	1:F:154:ALA:O	2.48	0.45
1:H:73:TYR:HA	1:H:76:PHE:HB2	1.97	0.45
1:C:17:TYR:HA	1:C:56:TYR:HA	1.98	0.45
1:I:22:THR:HG21	1:I:422:THR:HG21	1.98	0.45
1:I:376:MET:HE3	1:K:413:LEU:HD11	1.99	0.45
1:E:287:LEU:HD13	1:E:315:PHE:CD1	2.52	0.44
1:C:32:PHE:HA	1:C:33:PRO:HA	1.77	0.44
1:G:360:ARG:HE	1:G:360:ARG:HA	1.83	0.44
1:E:28:LEU:HB3	1:E:39:TRP:CE2	2.53	0.44
1:C:329:LEU:HD11	1:C:343:LEU:HB3	1.98	0.44
1:H:9:ASN:OD1	1:I:390:ILE:HD12	2.18	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:286:LEU:HD22	1:L:286:LEU:HA	1.86	0.44
1:D:202:PRO:HA	1:D:205:TRP:CD2	2.53	0.44
1:E:202:PRO:HA	1:E:205:TRP:CE2	2.51	0.44
1:G:11:THR:HB	1:H:1:PRO:HG3	1.98	0.44
1:L:9:ASN:O	1:L:10:SER:OG	2.30	0.44
1:L:295:LEU:HA	1:L:298:ILE:HD12	2.00	0.44
1:H:315:PHE:HE1	1:H:350:THR:HG21	1.83	0.44
1:I:129:LEU:HB2	1:I:206:TYR:HA	1.99	0.44
1:L:17:TYR:HA	1:L:56:TYR:HA	1.99	0.44
1:C:165:ASN:HA	1:C:170:LEU:HD12	1.99	0.44
1:H:131:LYS:HA	1:H:138:GLY:HA3	2.00	0.44
1:H:173:ASN:HB3	1:H:177:ARG:HB2	1.99	0.44
1:L:328:TRP:O	1:L:332:ARG:HG2	2.18	0.44
1:D:24:THR:OG1	1:D:282:GLN:OE1	2.33	0.43
1:L:202:PRO:HA	1:L:205:TRP:CE2	2.52	0.43
1:A:101:THR:HG21	1:A:124:SER:HA	1.99	0.43
1:C:25:PRO:HG2	1:C:92:ALA:HA	2.00	0.43
1:F:90:GLN:HB3	1:F:391:THR:CG2	2.48	0.43
1:I:32:PHE:HA	1:I:33:PRO:HA	1.79	0.43
1:C:91:ASP:OD1	1:C:393:HIS:HB2	2.17	0.43
1:F:127:PRO:HD3	1:F:239:VAL:HG13	2.01	0.43
1:G:93:VAL:HG11	1:G:275:LEU:HD21	1.99	0.43
1:I:31:GLY:HA3	1:I:98:SER:HB3	2.00	0.43
1:A:57:ASP:O	1:A:64:SER:OG	2.36	0.43
1:F:97:PHE:HA	1:F:122:LEU:O	2.18	0.43
1:F:101:THR:HG21	1:F:124:SER:HA	2.00	0.43
1:G:17:TYR:HA	1:G:56:TYR:HA	2.00	0.43
1:G:101:THR:HG21	1:G:124:SER:HA	2.01	0.43
1:K:169:ASN:HB2	1:K:256:PRO:HB3	1.99	0.43
1:L:383:TYR:CZ	1:L:387:LYS:HD2	2.53	0.43
1:B:99:MET:HE2	1:B:99:MET:HB3	1.93	0.43
1:J:342:ILE:HA	1:J:411:MET:HE1	2.00	0.43
1:C:328:TRP:O	1:C:332:ARG:HG2	2.19	0.43
1:C:351:VAL:HA	1:C:352:PRO:HD3	1.91	0.43
1:E:328:TRP:O	1:E:332:ARG:HG2	2.19	0.43
1:H:28:LEU:HB3	1:H:39:TRP:CE2	2.53	0.43
1:I:101:THR:HG21	1:I:124:SER:HA	2.00	0.43
1:L:360:ARG:HE	1:L:360:ARG:HA	1.84	0.43
1:E:149:VAL:HG22	1:E:202:PRO:HB2	2.00	0.43
1:E:342:ILE:HA	1:E:411:MET:HE1	2.01	0.43
1:H:293:TYR:O	1:H:296:SER:OG	2.31	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:25:PRO:HA	1:K:52:ARG:HB3	2.01	0.43
1:C:73:TYR:HA	1:C:76:PHE:HB2	2.00	0.42
1:G:10:SER:O	1:H:17:TYR:OH	2.37	0.42
1:K:170:LEU:HD23	1:K:170:LEU:HA	1.88	0.42
1:B:342:ILE:HG12	1:B:411:MET:HE1	2.00	0.42
1:B:360:ARG:HH21	1:E:358:GLN:HG3	1.84	0.42
1:F:202:PRO:HA	1:F:205:TRP:CD2	2.53	0.42
1:H:19:ASP:OD2	1:H:52:ARG:NH2	2.44	0.42
1:I:202:PRO:HA	1:I:205:TRP:CD2	2.54	0.42
1:A:90:GLN:CB	1:A:390:ILE:HD13	2.45	0.42
1:G:7:GLN:HG3	1:G:8:GLU:N	2.34	0.42
1:H:259:LEU:HD23	1:H:267:VAL:HG21	2.01	0.42
1:L:282:GLN:HA	1:L:285:LYS:HB3	2.00	0.42
1:D:282:GLN:HE21	1:D:282:GLN:HB2	1.68	0.42
1:G:293:TYR:O	1:G:296:SER:OG	2.33	0.42
1:L:129:LEU:HB2	1:L:206:TYR:HA	2.00	0.42
1:L:173:ASN:HB3	1:L:177:ARG:HB2	2.01	0.42
1:L:286:LEU:O	1:L:290:VAL:HG23	2.19	0.42
1:L:307:ILE:HD11	1:L:335:LEU:HD11	2.02	0.42
1:A:418:MET:HE2	1:A:418:MET:HB3	1.93	0.42
1:L:293:TYR:O	1:L:296:SER:OG	2.35	0.42
1:A:329:LEU:HD11	1:A:343:LEU:HB3	2.01	0.42
1:A:1:PRO:HD2	1:A:16:TYR:HE1	1.85	0.42
1:A:33:PRO:HD3	1:A:163:PHE:CE2	2.55	0.42
1:A:249:TYR:CE2	1:A:251:GLU:HG3	2.54	0.42
1:D:127:PRO:HD3	1:D:239:VAL:HG13	2.01	0.42
1:I:399:SER:O	1:I:407:LEU:HD21	2.19	0.42
1:J:291:GLU:HB2	1:J:315:PHE:HE2	1.85	0.42
1:K:127:PRO:HD3	1:K:239:VAL:HG13	2.01	0.42
1:A:97:PHE:HA	1:A:122:LEU:O	2.20	0.42
1:A:157:TYR:OH	1:B:40:GLU:OE2	2.27	0.42
1:E:32:PHE:HA	1:E:33:PRO:HA	1.80	0.42
1:J:97:PHE:HA	1:J:122:LEU:O	2.19	0.42
1:A:287:LEU:HD13	1:A:315:PHE:CD1	2.55	0.42
1:H:360:ARG:HE	1:H:360:ARG:HA	1.84	0.42
1:I:33:PRO:HD3	1:I:163:PHE:CE2	2.54	0.42
1:A:1:PRO:HD2	1:A:16:TYR:CE1	2.55	0.42
1:D:315:PHE:HE1	1:D:350:THR:HG21	1.84	0.42
1:L:282:GLN:O	1:L:286:LEU:N	2.52	0.42
1:A:302:PRO:O	1:A:306:GLU:HG3	2.20	0.41
1:I:97:PHE:HA	1:I:122:LEU:O	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:40:GLU:HA	1:D:43:SER:HB2	2.01	0.41
1:H:231:LEU:HB3	1:H:236:THR:HG21	2.01	0.41
1:G:351:VAL:HA	1:G:352:PRO:HD3	1.91	0.41
1:I:334:ILE:HA	1:J:211:ALA:HA	2.02	0.41
1:A:25:PRO:HG2	1:A:92:ALA:HA	2.02	0.41
1:A:149:VAL:HG22	1:A:202:PRO:HB2	2.02	0.41
1:A:358:GLN:OE1	1:J:361:ARG:N	2.54	0.41
1:B:25:PRO:HG2	1:B:92:ALA:HA	2.01	0.41
1:G:10:SER:OG	1:H:19:ASP:HB3	2.21	0.41
1:H:282:GLN:HA	1:H:285:LYS:HG2	2.02	0.41
1:A:231:LEU:HA	1:A:232:PRO:HD2	1.88	0.41
1:A:335:LEU:HD23	1:A:335:LEU:HA	1.90	0.41
1:D:93:VAL:HG11	1:D:275:LEU:HD21	2.03	0.41
1:F:259:LEU:HD23	1:F:267:VAL:HG21	2.02	0.41
1:K:295:LEU:HD23	1:K:298:ILE:HD12	2.02	0.41
1:L:127:PRO:HD3	1:L:239:VAL:HG13	2.02	0.41
1:B:202:PRO:HA	1:B:205:TRP:CE2	2.55	0.41
1:E:307:ILE:HD13	1:E:343:LEU:HD11	2.02	0.41
1:F:351:VAL:HA	1:F:352:PRO:HD3	1.90	0.41
1:H:1:PRO:HD2	1:H:16:TYR:HE1	1.85	0.41
1:K:360:ARG:HE	1:K:360:ARG:HA	1.85	0.41
1:B:32:PHE:HA	1:B:33:PRO:HA	1.76	0.41
1:B:40:GLU:HA	1:B:43:SER:HB2	2.03	0.41
1:C:129:LEU:HD23	1:C:129:LEU:HA	1.88	0.41
1:F:17:TYR:HA	1:F:56:TYR:HA	2.02	0.41
1:G:202:PRO:HA	1:G:205:TRP:CE2	2.56	0.41
1:G:255:ALA:HA	1:G:263:HIS:CE1	2.56	0.41
1:H:74:ASP:OD1	1:H:106:ARG:NH1	2.48	0.41
1:H:97:PHE:HA	1:H:122:LEU:O	2.21	0.41
1:H:129:LEU:HD23	1:H:129:LEU:HA	1.95	0.41
1:L:231:LEU:HA	1:L:232:PRO:HD2	1.89	0.41
1:A:426:ALA:O	1:A:430:VAL:HG23	2.21	0.41
1:B:351:VAL:HA	1:B:352:PRO:HD3	1.86	0.41
1:C:279:GLN:O	1:C:282:GLN:HG3	2.21	0.41
1:D:335:LEU:HD23	1:D:335:LEU:HA	1.97	0.41
1:E:351:VAL:HA	1:E:352:PRO:HD3	1.87	0.41
1:F:31:GLY:HA3	1:F:98:SER:HB3	2.02	0.41
1:I:231:LEU:HA	1:I:232:PRO:HD2	1.86	0.41
1:I:336:SER:HB3	1:I:339:THR:OG1	2.21	0.41
1:I:390:ILE:HG23	1:I:391:THR:HG23	2.02	0.41
1:J:351:VAL:HA	1:J:352:PRO:HD3	1.92	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:234:GLU:H	1:B:234:GLU:CD	2.29	0.40
1:C:307:ILE:HD11	1:C:335:LEU:HD11	2.04	0.40
1:E:202:PRO:HA	1:E:205:TRP:CD2	2.57	0.40
1:E:295:LEU:HD23	1:E:298:ILE:HD12	2.03	0.40
1:J:101:THR:HG21	1:J:124:SER:HA	2.03	0.40
1:K:131:LYS:HA	1:K:138:GLY:HA3	2.03	0.40
1:A:336:SER:HA	1:A:337:PRO:HD3	1.98	0.40
1:B:202:PRO:HA	1:B:205:TRP:CD2	2.57	0.40
1:C:97:PHE:HA	1:C:122:LEU:O	2.20	0.40
1:E:259:LEU:HD23	1:E:267:VAL:HG21	2.03	0.40
1:G:426:ALA:O	1:G:430:VAL:HG23	2.21	0.40
1:J:88:ASP:HA	1:J:115:ARG:HH12	1.87	0.40
1:B:407:LEU:HD12	1:B:436:GLN:HG3	2.03	0.40
1:C:315:PHE:CE1	1:C:350:THR:HG21	2.49	0.40
1:E:127:PRO:HD3	1:E:239:VAL:HG13	2.03	0.40
1:G:335:LEU:HD23	1:G:335:LEU:HA	1.97	0.40
1:I:328:TRP:O	1:I:332:ARG:HG2	2.21	0.40
1:I:360:ARG:HE	1:I:360:ARG:HA	1.86	0.40
1:J:1:PRO:HD2	1:J:16:TYR:HE1	1.86	0.40
1:D:275:LEU:O	1:D:279:GLN:N	2.47	0.40
1:F:40:GLU:HA	1:F:43:SER:HB2	2.03	0.40
1:J:73:TYR:HA	1:J:76:PHE:HB2	2.02	0.40
1:F:360:ARG:HE	1:F:360:ARG:HA	1.86	0.40
1:H:27:VAL:HG21	1:H:84:LEU:HD21	2.04	0.40
1:K:9:ASN:C	1:K:11:THR:H	2.30	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	438/456 (96%)	414 (94%)	23 (5%)	1 (0%)	43 72

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	438/456 (96%)	416 (95%)	22 (5%)	0	100	100
1	C	438/456 (96%)	415 (95%)	23 (5%)	0	100	100
1	D	438/456 (96%)	414 (94%)	22 (5%)	2 (0%)	24	57
1	E	438/456 (96%)	413 (94%)	25 (6%)	0	100	100
1	F	438/456 (96%)	414 (94%)	24 (6%)	0	100	100
1	G	438/456 (96%)	418 (95%)	20 (5%)	0	100	100
1	H	438/456 (96%)	418 (95%)	19 (4%)	1 (0%)	43	72
1	I	438/456 (96%)	412 (94%)	26 (6%)	0	100	100
1	J	438/456 (96%)	414 (94%)	23 (5%)	1 (0%)	43	72
1	K	438/456 (96%)	414 (94%)	24 (6%)	0	100	100
1	L	438/456 (96%)	414 (94%)	24 (6%)	0	100	100
All	All	5256/5472 (96%)	4976 (95%)	275 (5%)	5 (0%)	48	79

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	D	9	ASN
1	H	9	ASN
1	J	9	ASN
1	D	319	ASN

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	354/370 (96%)	344 (97%)	10 (3%)	38	60
1	B	354/370 (96%)	346 (98%)	8 (2%)	44	64
1	C	354/370 (96%)	345 (98%)	9 (2%)	42	63
1	D	354/370 (96%)	344 (97%)	10 (3%)	38	60

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	354/370 (96%)	345 (98%)	9 (2%)	42	63
1	F	354/370 (96%)	344 (97%)	10 (3%)	38	60
1	G	354/370 (96%)	345 (98%)	9 (2%)	42	63
1	H	354/370 (96%)	343 (97%)	11 (3%)	35	59
1	I	354/370 (96%)	347 (98%)	7 (2%)	48	66
1	J	354/370 (96%)	345 (98%)	9 (2%)	42	63
1	K	354/370 (96%)	347 (98%)	7 (2%)	48	66
1	L	354/370 (96%)	346 (98%)	8 (2%)	44	64
All	All	4248/4440 (96%)	4141 (98%)	107 (2%)	42	63

All (107) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	103	GLU
1	A	259	LEU
1	A	279	GLN
1	A	286	LEU
1	A	319	ASN
1	A	329	LEU
1	A	349	LEU
1	A	354	GLU
1	A	357	LEU
1	A	393	HIS
1	B	103	GLU
1	B	259	LEU
1	B	285	LYS
1	B	329	LEU
1	B	338	LEU
1	B	349	LEU
1	B	354	GLU
1	B	357	LEU
1	C	103	GLU
1	C	259	LEU
1	C	286	LEU
1	C	329	LEU
1	C	338	LEU
1	C	349	LEU
1	C	354	GLU
1	C	357	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	384	ARG
1	D	259	LEU
1	D	282	GLN
1	D	285	LYS
1	D	286	LEU
1	D	303	LEU
1	D	329	LEU
1	D	349	LEU
1	D	354	GLU
1	D	357	LEU
1	D	393	HIS
1	E	103	GLU
1	E	165	ASN
1	E	259	LEU
1	E	303	LEU
1	E	329	LEU
1	E	349	LEU
1	E	354	GLU
1	E	357	LEU
1	E	393	HIS
1	F	103	GLU
1	F	259	LEU
1	F	285	LYS
1	F	286	LEU
1	F	303	LEU
1	F	329	LEU
1	F	349	LEU
1	F	354	GLU
1	F	357	LEU
1	F	393	HIS
1	G	259	LEU
1	G	285	LYS
1	G	286	LEU
1	G	303	LEU
1	G	329	LEU
1	G	349	LEU
1	G	354	GLU
1	G	357	LEU
1	G	393	HIS
1	H	103	GLU
1	H	259	LEU
1	H	282	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	H	286	LEU
1	H	303	LEU
1	H	329	LEU
1	H	338	LEU
1	H	349	LEU
1	H	354	GLU
1	H	357	LEU
1	H	393	HIS
1	I	103	GLU
1	I	259	LEU
1	I	303	LEU
1	I	349	LEU
1	I	354	GLU
1	I	357	LEU
1	I	393	HIS
1	J	103	GLU
1	J	259	LEU
1	J	286	LEU
1	J	303	LEU
1	J	329	LEU
1	J	349	LEU
1	J	354	GLU
1	J	357	LEU
1	J	393	HIS
1	K	259	LEU
1	K	303	LEU
1	K	329	LEU
1	K	349	LEU
1	K	354	GLU
1	K	357	LEU
1	K	393	HIS
1	L	259	LEU
1	L	286	LEU
1	L	303	LEU
1	L	329	LEU
1	L	349	LEU
1	L	354	GLU
1	L	357	LEU
1	L	384	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	GLN
1	A	81	ASN
1	A	165	ASN
1	A	365	ASN
1	B	81	ASN
1	B	165	ASN
1	B	365	ASN
1	C	81	ASN
1	C	165	ASN
1	C	364	GLN
1	C	365	ASN
1	D	63	GLN
1	D	81	ASN
1	D	365	ASN
1	E	81	ASN
1	E	365	ASN
1	F	81	ASN
1	F	165	ASN
1	F	257	HIS
1	F	365	ASN
1	F	370	ASN
1	G	268	ASN
1	G	282	GLN
1	G	358	GLN
1	H	81	ASN
1	H	165	ASN
1	H	364	GLN
1	I	9	ASN
1	I	81	ASN
1	I	257	HIS
1	I	365	ASN
1	J	63	GLN
1	J	81	ASN
1	J	165	ASN
1	J	279	GLN
1	J	365	ASN
1	K	9	ASN
1	K	81	ASN
1	K	165	ASN
1	K	365	ASN
1	L	42	GLN
1	L	81	ASN
1	L	165	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	L	364	GLN
1	L	365	ASN

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 [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	440/456 (96%)	-0.56	1 (0%) 91 80	63, 94, 154, 182	0
1	B	440/456 (96%)	-0.55	0 100 100	72, 106, 155, 220	0
1	C	440/456 (96%)	-0.44	0 100 100	68, 122, 214, 270	0
1	D	440/456 (96%)	-0.43	2 (0%) 87 66	84, 135, 228, 266	0
1	E	440/456 (96%)	-0.57	0 100 100	74, 114, 155, 200	0
1	F	440/456 (96%)	-0.47	0 100 100	113, 184, 300, 340	0
1	G	440/456 (96%)	-0.32	1 (0%) 91 80	118, 167, 247, 287	0
1	H	440/456 (96%)	-0.36	0 100 100	87, 129, 332, 377	0
1	I	440/456 (96%)	-0.46	0 100 100	90, 132, 168, 218	0
1	J	440/456 (96%)	-0.56	0 100 100	81, 118, 153, 187	0
1	K	440/456 (96%)	-0.48	1 (0%) 91 80	93, 156, 213, 271	0
1	L	440/456 (96%)	-0.28	1 (0%) 91 80	130, 211, 375, 404	0
All	All	5280/5472 (96%)	-0.46	6 (0%) 92 85	63, 134, 291, 404	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	430	VAL	2.5
1	K	413	LEU	2.4
1	L	245	PRO	2.2
1	A	438	ALA	2.2
1	G	54	ILE	2.1
1	D	294	VAL	2.1

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.