



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:49 AM GMT

PDB ID : 2OBA
Title : Pseudomonas aeruginosa 6-pyruvoyl tetrahydrobiopterin synthase
Authors : McGrath, T.E.; Kisselman, G.; Battaile, K.; Romanov, V.; Wu-Brown, J.; Guthrie, J.; Virag, C.; Mansoury, K.; Edwards, A.M.; Pai, E.F.; Chirgadze, N.Y.
Deposited on : 2006-12-18
Resolution : 2.33 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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

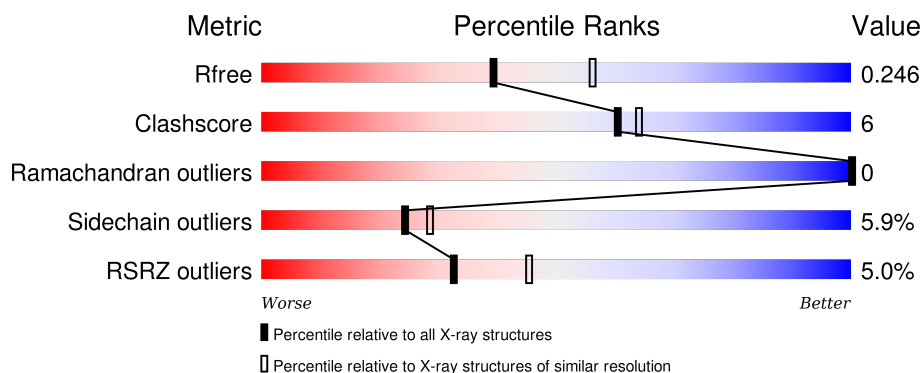
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.33 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}	91344	1406 (2.36-2.32)
Clashscore	102246	1509 (2.36-2.32)
Ramachandran outliers	100387	1490 (2.36-2.32)
Sidechain outliers	100360	1491 (2.36-2.32)
RSRZ outliers	91569	1412 (2.36-2.32)

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. 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	138	<div> <div>5%</div> <div> <div></div> <div>72%</div> <div>14%</div> <div>•</div> <div>13%</div> </div> </div>
1	B	138	<div> <div>4%</div> <div> <div></div> <div>61%</div> <div>25%</div> <div>•</div> <div>12%</div> </div> </div>
1	C	138	<div> <div>3%</div> <div> <div></div> <div>75%</div> <div>13%</div> <div>•</div> <div>12%</div> </div> </div>
1	D	138	<div> <div>4%</div> <div> <div></div> <div>76%</div> <div>10%</div> <div>•</div> <div>13%</div> </div> </div>
1	E	138	<div> <div>6%</div> <div> <div></div> <div>72%</div> <div>14%</div> <div>•</div> <div>12%</div> </div> </div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	138	 A horizontal bar chart showing the quality of chain F. The bar is divided into four segments: a small red segment at the beginning labeled '4%', followed by a long green segment labeled '70%', then a yellow segment labeled '17%', and finally a small grey segment at the end labeled '12%'.

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6228 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 Probable 6-pyruvoyl tetrahydrobiopterin synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	120	Total	C	N	O	S	Se	0	1	0
			994	635	177	177	4	1			
1	B	121	Total	C	N	O	S	Se	0	0	0
			993	632	178	178	4	1			
1	C	121	Total	C	N	O	S	Se	0	0	0
			993	632	178	178	4	1			
1	D	120	Total	C	N	O	S	Se	0	0	0
			989	630	177	177	4	1			
1	E	121	Total	C	N	O	S	Se	0	1	0
			1001	637	181	178	4	1			
1	F	121	Total	C	N	O	S	Se	0	0	0
			993	632	178	178	4	1			

There are 126 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MSE	-	SEE REMARK 999	UNP Q9I0H2
A	-18	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
A	-17	SER	-	CLONING ARTIFACT	UNP Q9I0H2
A	-16	SER	-	CLONING ARTIFACT	UNP Q9I0H2
A	-15	HIS	-	EXPRESSION TAG	UNP Q9I0H2
A	-14	HIS	-	EXPRESSION TAG	UNP Q9I0H2
A	-13	HIS	-	EXPRESSION TAG	UNP Q9I0H2
A	-12	HIS	-	EXPRESSION TAG	UNP Q9I0H2
A	-11	HIS	-	EXPRESSION TAG	UNP Q9I0H2
A	-10	HIS	-	EXPRESSION TAG	UNP Q9I0H2
A	-9	SER	-	CLONING ARTIFACT	UNP Q9I0H2
A	-8	SER	-	CLONING ARTIFACT	UNP Q9I0H2
A	-7	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
A	-6	LEU	-	CLONING ARTIFACT	UNP Q9I0H2
A	-5	VAL	-	CLONING ARTIFACT	UNP Q9I0H2
A	-4	PRO	-	CLONING ARTIFACT	UNP Q9I0H2
A	-3	ARG	-	CLONING ARTIFACT	UNP Q9I0H2

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
A	-1	SER	-	CLONING ARTIFACT	UNP Q9I0H2
A	0	HIS	-	CLONING ARTIFACT	UNP Q9I0H2
A	1	MSE	-	CLONING ARTIFACT	UNP Q9I0H2
B	-19	MSE	-	SEE REMARK 999	UNP Q9I0H2
B	-18	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
B	-17	SER	-	CLONING ARTIFACT	UNP Q9I0H2
B	-16	SER	-	CLONING ARTIFACT	UNP Q9I0H2
B	-15	HIS	-	EXPRESSION TAG	UNP Q9I0H2
B	-14	HIS	-	EXPRESSION TAG	UNP Q9I0H2
B	-13	HIS	-	EXPRESSION TAG	UNP Q9I0H2
B	-12	HIS	-	EXPRESSION TAG	UNP Q9I0H2
B	-11	HIS	-	EXPRESSION TAG	UNP Q9I0H2
B	-10	HIS	-	EXPRESSION TAG	UNP Q9I0H2
B	-9	SER	-	CLONING ARTIFACT	UNP Q9I0H2
B	-8	SER	-	CLONING ARTIFACT	UNP Q9I0H2
B	-7	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
B	-6	LEU	-	CLONING ARTIFACT	UNP Q9I0H2
B	-5	VAL	-	CLONING ARTIFACT	UNP Q9I0H2
B	-4	PRO	-	CLONING ARTIFACT	UNP Q9I0H2
B	-3	ARG	-	CLONING ARTIFACT	UNP Q9I0H2
B	-2	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
B	-1	SER	-	CLONING ARTIFACT	UNP Q9I0H2
B	0	HIS	-	CLONING ARTIFACT	UNP Q9I0H2
B	1	MSE	-	CLONING ARTIFACT	UNP Q9I0H2
C	-19	MSE	-	SEE REMARK 999	UNP Q9I0H2
C	-18	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
C	-17	SER	-	CLONING ARTIFACT	UNP Q9I0H2
C	-16	SER	-	CLONING ARTIFACT	UNP Q9I0H2
C	-15	HIS	-	EXPRESSION TAG	UNP Q9I0H2
C	-14	HIS	-	EXPRESSION TAG	UNP Q9I0H2
C	-13	HIS	-	EXPRESSION TAG	UNP Q9I0H2
C	-12	HIS	-	EXPRESSION TAG	UNP Q9I0H2
C	-11	HIS	-	EXPRESSION TAG	UNP Q9I0H2
C	-10	HIS	-	EXPRESSION TAG	UNP Q9I0H2
C	-9	SER	-	CLONING ARTIFACT	UNP Q9I0H2
C	-8	SER	-	CLONING ARTIFACT	UNP Q9I0H2
C	-7	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
C	-6	LEU	-	CLONING ARTIFACT	UNP Q9I0H2
C	-5	VAL	-	CLONING ARTIFACT	UNP Q9I0H2
C	-4	PRO	-	CLONING ARTIFACT	UNP Q9I0H2
C	-3	ARG	-	CLONING ARTIFACT	UNP Q9I0H2

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
C	-1	SER	-	CLONING ARTIFACT	UNP Q9I0H2
C	0	HIS	-	CLONING ARTIFACT	UNP Q9I0H2
C	1	MSE	-	CLONING ARTIFACT	UNP Q9I0H2
D	-19	MSE	-	SEE REMARK 999	UNP Q9I0H2
D	-18	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
D	-17	SER	-	CLONING ARTIFACT	UNP Q9I0H2
D	-16	SER	-	CLONING ARTIFACT	UNP Q9I0H2
D	-15	HIS	-	EXPRESSION TAG	UNP Q9I0H2
D	-14	HIS	-	EXPRESSION TAG	UNP Q9I0H2
D	-13	HIS	-	EXPRESSION TAG	UNP Q9I0H2
D	-12	HIS	-	EXPRESSION TAG	UNP Q9I0H2
D	-11	HIS	-	EXPRESSION TAG	UNP Q9I0H2
D	-10	HIS	-	EXPRESSION TAG	UNP Q9I0H2
D	-9	SER	-	CLONING ARTIFACT	UNP Q9I0H2
D	-8	SER	-	CLONING ARTIFACT	UNP Q9I0H2
D	-7	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
D	-6	LEU	-	CLONING ARTIFACT	UNP Q9I0H2
D	-5	VAL	-	CLONING ARTIFACT	UNP Q9I0H2
D	-4	PRO	-	CLONING ARTIFACT	UNP Q9I0H2
D	-3	ARG	-	CLONING ARTIFACT	UNP Q9I0H2
D	-2	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
D	-1	SER	-	CLONING ARTIFACT	UNP Q9I0H2
D	0	HIS	-	CLONING ARTIFACT	UNP Q9I0H2
D	1	MSE	-	CLONING ARTIFACT	UNP Q9I0H2
E	-19	MSE	-	SEE REMARK 999	UNP Q9I0H2
E	-18	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
E	-17	SER	-	CLONING ARTIFACT	UNP Q9I0H2
E	-16	SER	-	CLONING ARTIFACT	UNP Q9I0H2
E	-15	HIS	-	EXPRESSION TAG	UNP Q9I0H2
E	-14	HIS	-	EXPRESSION TAG	UNP Q9I0H2
E	-13	HIS	-	EXPRESSION TAG	UNP Q9I0H2
E	-12	HIS	-	EXPRESSION TAG	UNP Q9I0H2
E	-11	HIS	-	EXPRESSION TAG	UNP Q9I0H2
E	-10	HIS	-	EXPRESSION TAG	UNP Q9I0H2
E	-9	SER	-	CLONING ARTIFACT	UNP Q9I0H2
E	-8	SER	-	CLONING ARTIFACT	UNP Q9I0H2
E	-7	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
E	-6	LEU	-	CLONING ARTIFACT	UNP Q9I0H2
E	-5	VAL	-	CLONING ARTIFACT	UNP Q9I0H2
E	-4	PRO	-	CLONING ARTIFACT	UNP Q9I0H2
E	-3	ARG	-	CLONING ARTIFACT	UNP Q9I0H2

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
E	-1	SER	-	CLONING ARTIFACT	UNP Q9I0H2
E	0	HIS	-	CLONING ARTIFACT	UNP Q9I0H2
E	1	MSE	-	CLONING ARTIFACT	UNP Q9I0H2
F	-19	MSE	-	SEE REMARK 999	UNP Q9I0H2
F	-18	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
F	-17	SER	-	CLONING ARTIFACT	UNP Q9I0H2
F	-16	SER	-	CLONING ARTIFACT	UNP Q9I0H2
F	-15	HIS	-	EXPRESSION TAG	UNP Q9I0H2
F	-14	HIS	-	EXPRESSION TAG	UNP Q9I0H2
F	-13	HIS	-	EXPRESSION TAG	UNP Q9I0H2
F	-12	HIS	-	EXPRESSION TAG	UNP Q9I0H2
F	-11	HIS	-	EXPRESSION TAG	UNP Q9I0H2
F	-10	HIS	-	EXPRESSION TAG	UNP Q9I0H2
F	-9	SER	-	CLONING ARTIFACT	UNP Q9I0H2
F	-8	SER	-	CLONING ARTIFACT	UNP Q9I0H2
F	-7	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
F	-6	LEU	-	CLONING ARTIFACT	UNP Q9I0H2
F	-5	VAL	-	CLONING ARTIFACT	UNP Q9I0H2
F	-4	PRO	-	CLONING ARTIFACT	UNP Q9I0H2
F	-3	ARG	-	CLONING ARTIFACT	UNP Q9I0H2
F	-2	GLY	-	CLONING ARTIFACT	UNP Q9I0H2
F	-1	SER	-	CLONING ARTIFACT	UNP Q9I0H2
F	0	HIS	-	CLONING ARTIFACT	UNP Q9I0H2
F	1	MSE	-	CLONING ARTIFACT	UNP Q9I0H2

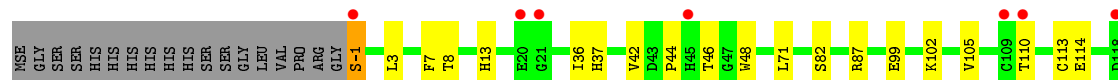
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	1	Total Zn 1 1	0	0
2	E	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	A	1	Total Zn 1 1	0	0
2	F	1	Total Zn 1 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	63	Total 63	O 63	0	0
3	B	32	Total 32	O 32	0	0
3	C	45	Total 45	O 45	0	0
3	D	52	Total 52	O 52	0	0
3	E	31	Total 31	O 31	0	0
3	F	36	Total 36	O 36	0	0

- Molecule 1: Probable 6-pyruvoyl tetrahydrobiopterin synthase



- | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|----|----|----|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|
| L71 | MSE | GLY | SER | SER | HIS | HIS | HIS | HIS | SER | SER | GLY | LEU | VAL | PRO | ARG | G-2 | H-1 | M1 | E2 | I3 | F4 | F9 | E10 | S11 | A12 | H13 | R14 | L15 | E20 | C24 | R33 | V34 | A35 | I36 | H37 | I38 | V42 | W48 | L49 | R50 | I55 | K60 | P61 | I62 | V63 | E64 | G65 | L66 |
|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|----|----|----|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|

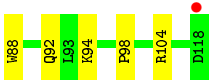
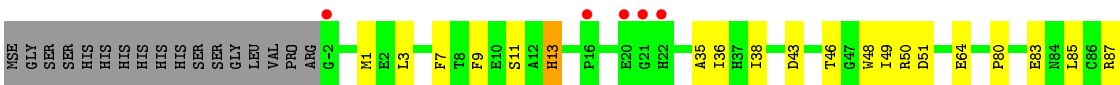
- | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|----|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|------|------|------|------|------|------|
| MSE | GLY | SER | SER | HIS | HIS | HIS | HIS | HIS | HIS | SER | SER | GLY | LEU | VAL | PRO | ARG | G-2 | S-1 | L3 | F4 | K5 | H13 | R33 | H37 | I38 | I48 | I49 | R50 | D51 | F52 | A53 | E54 | I55 | L71 | C86 | I93 | K102 | C109 | T110 | S111 | G112 | D113 |
|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|----|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|------|------|------|------|------|------|

- [illegible]

- | Category | Item | Value | Color |
|------------|--------|--------|--------|
| Nucleotide | MSE | 0.0000 | Grey |
| | GLY | 0.0000 | Grey |
| | SER | 0.0000 | Grey |
| | SER | 0.0000 | Grey |
| | HIS | 0.0000 | Grey |
| | HIS | 0.0000 | Grey |
| | HIS | 0.0000 | Grey |
| | HIS | 0.0000 | Grey |
| | HIS | 0.0000 | Grey |
| | HIS | 0.0000 | Grey |
| Amino Acid | SER | 0.0000 | Grey |
| | SER | 0.0000 | Grey |
| | GLY | 0.0000 | Grey |
| | VAL | 0.0000 | Grey |
| | LEU | 0.0000 | Grey |
| | ARG | 0.0000 | Grey |
| | ARG | 0.0000 | Grey |
| | G-2 | 0.0000 | Red |
| | S-1 | 0.0000 | Red |
| | H-0 | 0.0000 | Red |
| Other | F7 | 0.0000 | Yellow |
| | S11 | 0.0000 | Yellow |
| | A12 | 0.0000 | Yellow |
| | H13 | 0.0000 | Yellow |
| | R14 | 0.0000 | Yellow |
| | E20 | 0.0000 | Red |
| | G21 | 0.0000 | Red |
| | H30 | 0.0000 | Yellow |
| | V34 | 0.0000 | Yellow |
| | H37 | 0.0000 | Yellow |
| Other | P44 | 0.0000 | Red |
| | H45 | 0.0000 | Red |
| | W48 | 0.0000 | Orange |
| | K56 | 0.0000 | Yellow |
| | L71 | 0.0000 | Orange |
| | S82 | 0.0000 | Yellow |
| | E83 | 0.0000 | Yellow |
| | W88 | 0.0000 | Yellow |
| | Q92 | 0.0000 | Yellow |
| | K102 | 0.0000 | Yellow |
| V103 | 0.0000 | Yellow | |
| R104 | 0.0000 | Yellow | |
| Other | E107 | 0.0000 | Yellow |
| | T108 | 0.0000 | Green |



● Molecule 1: Probable 6-pyruvoyl tetrahydrobiopterin synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.08 Å 87.52 Å 124.67 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.22 – 2.33 42.23 – 2.33	Depositor EDS
% Data completeness (in resolution range)	97.7 (42.22-2.33) 97.7 (42.23-2.33)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.79 (at 2.32 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.186 , 0.256 0.182 , 0.246	Depositor DCC
R_{free} test set	1767 reflections (5.23%)	DCC
Wilson B-factor (Å ²)	44.8	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 35575 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6228	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.37% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: ZN

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	0.88	2/1028 (0.2%)	0.76	0/1392
1	B	0.68	0/1024	0.75	0/1386
1	C	0.70	1/1024 (0.1%)	0.72	0/1386
1	D	0.81	0/1020	0.77	0/1381
1	E	0.58	0/1035	0.67	0/1400
1	F	0.64	0/1024	0.73	4/1386 (0.3%)
All	All	0.72	3/6155 (0.0%)	0.73	4/8331 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	46	THR	C-O	9.57	1.41	1.23
1	A	44	PRO	C-N	6.03	1.48	1.34
1	C	86	CYS	CB-SG	-5.16	1.73	1.81

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	50	ARG	NE-CZ-NH2	-5.43	117.59	120.30
1	F	51	ASP	CB-CG-OD1	5.26	123.03	118.30
1	F	85	LEU	CA-CB-CG	5.13	127.11	115.30
1	F	85	LEU	CB-CG-CD2	-5.07	102.37	111.00

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	994	0	957	12	0
1	B	993	0	949	20	0
1	C	993	0	949	11	0
1	D	989	0	946	7	0
1	E	1001	0	962	11	0
1	F	993	0	949	13	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	63	0	0	3	0
3	B	32	0	0	0	0
3	C	45	0	0	1	0
3	D	52	0	0	1	0
3	E	31	0	0	1	0
3	F	36	0	0	0	0
All	All	6228	0	5712	65	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (65) 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:109:CYS:SG	3:E:208:HOH:O	2.39	0.80
1:C:3:LEU:HD12	1:C:49:ILE:HD12	1.64	0.79
1:D:104:ARG:HG3	1:D:114:GLU:HG3	1.69	0.73
1:D:114:GLU:OE2	1:D:116:ARG:NH2	2.22	0.72
1:D:107:GLU:HB3	3:D:250:HOH:O	1.89	0.71
1:C:3:LEU:CD1	1:C:49:ILE:HD12	2.20	0.71
1:B:102:LYS:HD2	1:B:114:GLU:CD	2.19	0.63
1:F:3:LEU:HG	1:F:49:ILE:HD12	1.81	0.63
1:F:7:PHE:CE2	1:F:36:ILE:HD12	2.34	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:34:VAL:HA	1:B:104:ARG:O	2.00	0.60
1:A:8:THR:HG21	3:A:232:HOH:O	1.99	0.60
1:F:94:LYS:NZ	1:F:98:PRO:O	2.30	0.60
1:D:37:HIS:HB2	1:D:102:LYS:HB3	1.83	0.59
1:A:-1:SER:O	1:B:118:ASP:O	2.20	0.59
1:C:38:ILE:HD13	1:C:55:ILE:HD11	1.84	0.58
1:B:60:LYS:HG2	1:B:64:GLU:OE2	2.04	0.58
1:B:37:HIS:HB2	1:B:102:LYS:HB3	1.86	0.58
1:B:62:ILE:CD1	1:B:93:LEU:HD13	2.33	0.57
1:A:37:HIS:HB2	1:A:102:LYS:HB3	1.86	0.57
1:E:88:TRP:O	1:E:92:GLN:HG2	2.06	0.56
1:E:104:ARG:HG3	1:E:114:GLU:HG3	1.88	0.55
1:E:30:HIS:CD2	1:E:107:GLU:HG2	2.43	0.54
1:F:43:ASP:HB3	1:F:46:THR:OG1	2.08	0.53
1:B:38:ILE:HD13	1:B:55:ILE:HD11	1.89	0.53
1:A:87:ARG:HH22	1:C:-1:SER:HB2	1.73	0.53
1:E:48:TRP:HB2	1:F:83:GLU:OE2	2.09	0.53
1:E:11:SER:HB3	1:E:71:LEU:HD13	1.91	0.52
1:E:7:PHE:HE1	1:E:56:LYS:HG3	1.75	0.52
1:B:33:ARG:NH2	1:B:109:CYS:HA	2.24	0.52
1:C:37:HIS:HB2	1:C:102:LYS:HB3	1.92	0.51
1:F:35:ALA:HB3	1:F:104:ARG:HB3	1.94	0.50
1:A:-1:SER:HB2	1:B:87:ARG:NH1	2.26	0.50
1:F:7:PHE:HE2	1:F:36:ILE:HD12	1.77	0.49
1:A:110:THR:O	1:C:5:LYS:HD2	2.13	0.48
1:F:13:HIS:HB2	1:F:80:PRO:HG2	1.95	0.48
1:B:11:SER:HB3	1:B:71:LEU:HD13	1.96	0.48
1:C:49:ILE:O	3:C:213:HOH:O	2.20	0.47
1:B:60:LYS:HB3	1:B:61:PRO:HD3	1.96	0.47
1:B:94:LYS:HA	1:B:94:LYS:HD2	1.64	0.47
1:A:87:ARG:NE	3:A:242:HOH:O	2.35	0.47
1:B:84:ASN:OD1	1:B:87:ARG:NH2	2.47	0.46
1:D:11:SER:HB3	1:D:71:LEU:HD13	1.96	0.46
1:D:102:LYS:HD2	1:D:114:GLU:OE2	2.16	0.46
1:F:3:LEU:HD12	1:F:38:ILE:HD11	1.98	0.46
1:E:-1:SER:OG	1:F:87:ARG:HD3	2.15	0.46
1:A:113:CYS:HB2	1:C:3:LEU:HD23	1.98	0.46
1:B:9:PHE:HD2	1:B:11:SER:OG	2.01	0.44
1:B:50:ARG:HD2	1:B:99:GLU:OE2	2.18	0.44
1:F:9:PHE:HD2	1:F:11:SER:OG	2.00	0.44
1:F:1:MSE:HB2	1:F:49:ILE:HD11	1.99	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:4:PHE:CE1	1:C:112:GLY:HA3	2.54	0.43
1:A:102:LYS:HD2	1:A:114:GLU:CD	2.39	0.43
1:A:87:ARG:NH2	3:A:242:HOH:O	2.44	0.42
1:B:4:PHE:HA	1:B:36:ILE:O	2.19	0.42
1:E:34:VAL:HA	1:E:104:ARG:O	2.18	0.42
1:B:15:LEU:HD12	1:B:24:CYS:HB3	2.02	0.42
1:D:42:VAL:HG22	1:E:83:GLU:HG2	2.02	0.42
1:C:51:ASP:HB3	1:C:54:GLU:HG3	2.02	0.42
1:A:7:PHE:HE2	1:A:36[A]:ILE:HD12	1.85	0.41
1:C:33:ARG:NH2	1:C:109:CYS:HA	2.35	0.41
1:E:37:HIS:HB2	1:E:102:LYS:HB3	2.02	0.41
1:A:102:LYS:HD2	1:A:114:GLU:CG	2.50	0.41
1:F:88:TRP:O	1:F:92:GLN:HG2	2.20	0.41
1:B:3:LEU:HD13	1:B:49:ILE:HD12	2.03	0.40
1:B:1:MSE:HB2	1:B:49:ILE:HD11	2.03	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	119/138 (86%)	116 (98%)	3 (2%)	0	100	100
1	B	119/138 (86%)	119 (100%)	0	0	100	100
1	C	119/138 (86%)	117 (98%)	2 (2%)	0	100	100
1	D	118/138 (86%)	116 (98%)	2 (2%)	0	100	100
1	E	120/138 (87%)	117 (98%)	3 (2%)	0	100	100
1	F	119/138 (86%)	117 (98%)	2 (2%)	0	100	100
All	All	714/828 (86%)	702 (98%)	12 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	109/121 (90%)	100 (92%)	9 (8%)	14	14
1	B	108/121 (89%)	99 (92%)	9 (8%)	14	14
1	C	108/121 (89%)	104 (96%)	4 (4%)	41	53
1	D	108/121 (89%)	101 (94%)	7 (6%)	21	24
1	E	109/121 (90%)	103 (94%)	6 (6%)	27	32
1	F	108/121 (89%)	105 (97%)	3 (3%)	51	64
All	All	650/726 (90%)	612 (94%)	38 (6%)	24	29

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-1	SER
1	A	3	LEU
1	A	13	HIS
1	A	42	VAL
1	A	48	TRP
1	A	71	LEU
1	A	82	SER
1	A	99	GLU
1	A	105	VAL
1	B	3	LEU
1	B	13	HIS
1	B	42	VAL
1	B	48	TRP
1	B	66	LEU
1	B	71	LEU
1	B	82	SER
1	B	85	LEU
1	B	92	GLN
1	C	13	HIS
1	C	48	TRP
1	C	71	LEU
1	C	93	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	3	LEU
1	D	13	HIS
1	D	20	GLU
1	D	48	TRP
1	D	71	LEU
1	D	93	LEU
1	D	105	VAL
1	E	13	HIS
1	E	14	ARG
1	E	20	GLU
1	E	48	TRP
1	E	71	LEU
1	E	82	SER
1	F	13	HIS
1	F	48	TRP
1	F	64	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

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.

5.7 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	119/138 (86%)	0.33	7 (5%)	26 37	39, 45, 54, 58	0
1	B	120/138 (86%)	0.18	5 (4%)	40 52	39, 45, 53, 57	0
1	C	120/138 (86%)	0.17	4 (3%)	50 61	39, 45, 52, 58	0
1	D	119/138 (86%)	0.40	6 (5%)	32 45	39, 45, 53, 60	0
1	E	120/138 (86%)	0.37	8 (6%)	21 31	40, 45, 54, 61	0
1	F	120/138 (86%)	0.20	6 (5%)	32 45	39, 45, 54, 62	0
All	All	718/828 (86%)	0.27	36 (5%)	32 45	39, 45, 54, 62	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	-2	GLY	5.1
1	A	45	HIS	4.8
1	E	-2	GLY	4.3
1	F	-2	GLY	4.0
1	E	0	HIS	3.8
1	E	-1	SER	3.6
1	E	21	GLY	3.5
1	B	-1	SER	3.4
1	C	118	ASP	3.3
1	A	20	GLU	3.1
1	F	20	GLU	3.1
1	B	20	GLU	3.0
1	A	109	CYS	3.0
1	E	45	HIS	2.9
1	F	16	PRO	2.9
1	E	118	ASP	2.9
1	F	118	ASP	2.9
1	D	118	ASP	2.7
1	D	-1	SER	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	109	CYS	2.6
1	F	21	GLY	2.6
1	D	45	HIS	2.6
1	A	-1	SER	2.5
1	C	109	CYS	2.4
1	B	118	ASP	2.3
1	C	110	THR	2.3
1	A	110	THR	2.2
1	A	21	GLY	2.2
1	F	22	HIS	2.2
1	D	34	VAL	2.2
1	E	44	PRO	2.2
1	D	0	HIS	2.2
1	E	20	GLU	2.1
1	C	52	PHE	2.1
1	B	87	ARG	2.1
1	A	118	ASP	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	ZN	F	200	1/1	0.98	0.03	-3.33	56,56,56,56	0
2	ZN	D	200	1/1	0.99	0.02	-4.05	42,42,42,42	0
2	ZN	C	200	1/1	0.99	0.02	-4.99	50,50,50,50	0
2	ZN	E	200	1/1	0.97	0.04	-6.84	53,53,53,53	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	ZN	B	200	1/1	0.99	0.02	-8.29	54,54,54,54	0
2	ZN	A	200	1/1	1.00	0.02	-	37,37,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.