



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 12:05 PM UTC

PDB ID : 8OSY / pdb_00008osy
Title : Trimeric catalytic domain of the E. coli Dihydrolipoamide Acetyltransferase (E2) of the pyruvate dehydrogenase complex
Authors : Meinhold, S.; Zdanowicz, R.; Glockshuber, R.
Deposited on : 2023-04-20
Resolution : 1.89 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

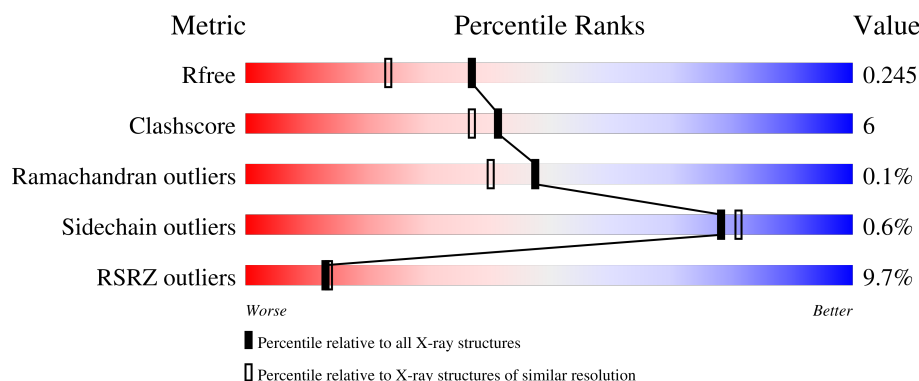
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	247	4% 86% 13% .
1	B	247	2% 88% 11% .
1	C	247	3% 91% 8% .
1	D	247	30% 77% 22% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8096 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 Dihydrolipoyllysine-residue acetyltransferase component of pyruvate dehydrogenase complex.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	245	Total	C	N	O	S	0	0	0
			1901	1214	331	347	9			
1	B	245	Total	C	N	O	S	0	0	0
			1901	1214	331	347	9			
1	C	245	Total	C	N	O	S	0	1	0
			1910	1219	333	349	9			
1	D	245	Total	C	N	O	S	0	0	0
			1901	1214	331	347	9			

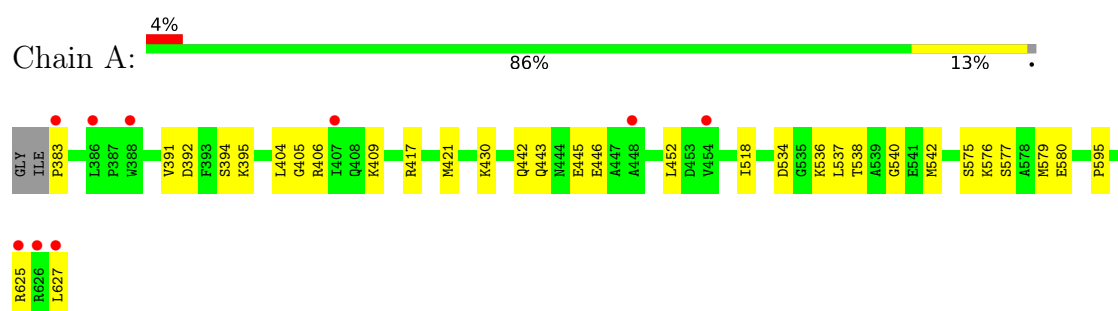
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	142	Total	O	0	0
			142	142		
2	B	180	Total	O	0	0
			180	180		
2	C	150	Total	O	0	0
			150	150		
2	D	11	Total	O	0	0
			11	11		

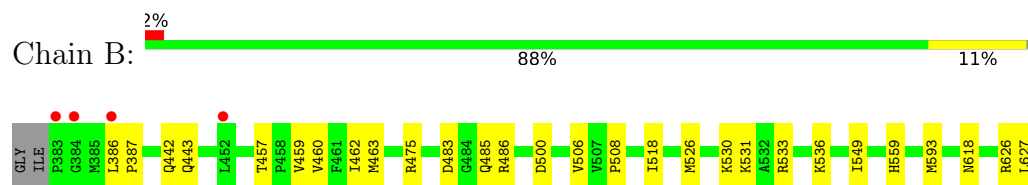
3 Residue-property plots [i](#)

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.

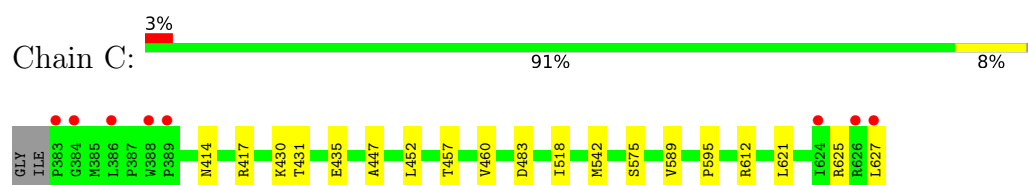
- Molecule 1: Dihydrolipoyllysine-residue acetyltransferase component of pyruvate dehydrogenase complex



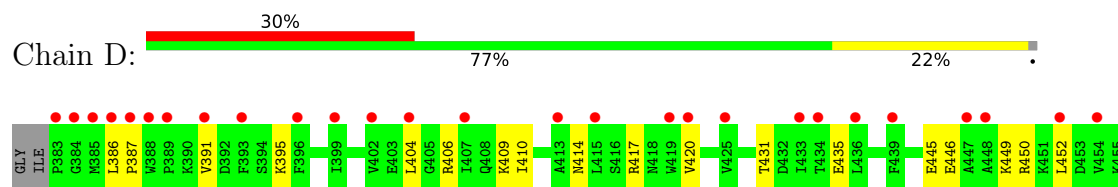
- Molecule 1: Dihydrolipoyllysine-residue acetyltransferase component of pyruvate dehydrogenase complex

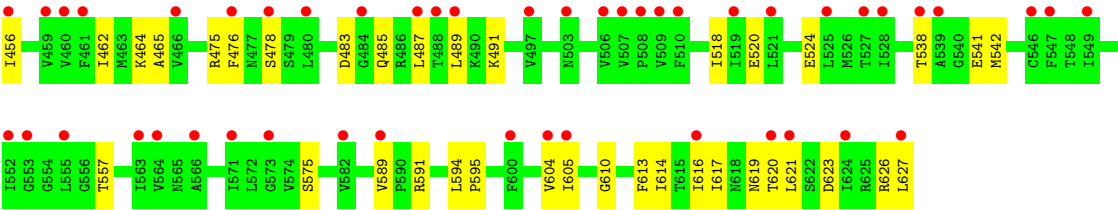


- Molecule 1: Dihydrolipoyllysine-residue acetyltransferase component of pyruvate dehydrogenase complex



- Molecule 1: Dihydrolipoyllysine-residue acetyltransferase component of pyruvate dehydrogenase complex





4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	148.15Å 148.15Å 119.95Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.96 – 1.89 44.96 – 1.89	Depositor EDS
% Data completeness (in resolution range)	99.7 (44.96-1.89) 100.0 (44.96-1.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.15 (at 1.88Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.206 , 0.244 (Not available) , 0.245	Depositor DCC
R_{free} test set	3951 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	40.8	Xtriage
Anisotropy	0.118	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 40.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.016 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8096	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.24	0/1936	0.46	0/2613
1	B	0.23	0/1936	0.45	0/2613
1	C	0.23	0/1945	0.46	0/2625
1	D	0.17	0/1936	0.41	0/2613
All	All	0.22	0/7753	0.45	0/10464

There are no bond length outliers.

There are no bond angle outliers.

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	1901	0	1967	25	0
1	B	1901	0	1967	23	0
1	C	1910	0	1974	14	0
1	D	1901	0	1967	43	0
2	A	142	0	0	0	0
2	B	180	0	0	2	0
2	C	150	0	0	1	0
2	D	11	0	0	0	0
All	All	8096	0	7875	99	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 (99) 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:D:414:ASN:CG	1:D:417:ARG:HH12	1.40	1.27
1:D:414:ASN:OD1	1:D:417:ARG:NH1	1.70	1.23
1:D:414:ASN:HA	1:D:417:ARG:NH1	1.67	1.10
1:D:414:ASN:HA	1:D:417:ARG:HH11	1.16	1.10
1:D:414:ASN:CG	1:D:417:ARG:NH1	2.19	0.96
1:D:414:ASN:CA	1:D:417:ARG:NH1	2.42	0.82
1:B:518:ILE:H	1:B:627:LEU:HD23	1.48	0.79
1:B:526:MET:O	1:B:530:LYS:HD3	1.83	0.79
1:D:518:ILE:H	1:D:627:LEU:HD22	1.49	0.76
1:A:383:PRO:HD3	1:C:483:ASP:HA	1.67	0.75
1:B:386:LEU:HD12	1:B:387:PRO:HD2	1.72	0.71
1:D:478:SER:HB2	1:D:487:LEU:HD11	1.75	0.69
1:D:414:ASN:CB	1:D:417:ARG:NH1	2.59	0.66
1:A:430:LYS:HE2	1:A:580:GLU:OE2	1.96	0.64
1:A:576:LYS:HD3	1:A:577:SER:N	2.11	0.64
1:C:414:ASN:HD22	1:C:417:ARG:NH2	1.95	0.64
1:C:414:ASN:HD22	1:C:417:ARG:HH22	1.47	0.62
1:B:530:LYS:HD2	1:B:533:ARG:HH22	1.66	0.61
1:A:406:ARG:HB2	1:D:623:ASP:HB3	1.83	0.60
1:D:404:LEU:O	1:D:409:LYS:NZ	2.35	0.60
1:C:518:ILE:HB	1:C:627:LEU:HB3	1.84	0.59
1:D:478:SER:HB3	1:D:489:LEU:HD23	1.83	0.58
1:D:626:ARG:HB3	1:D:627:LEU:HD12	1.86	0.58
1:D:616:ILE:O	1:D:620:THR:HG22	2.04	0.58
1:D:475:ARG:O	1:D:478:SER:OG	2.19	0.57
1:B:526:MET:HE3	1:B:530:LYS:HZ1	1.72	0.55
1:A:442:GLN:NE2	1:A:443:GLN:OE1	2.35	0.55
1:D:476:PHE:CE1	1:D:605:ILE:HD12	2.42	0.54
1:A:625:ARG:HH11	1:A:625:ARG:HG3	1.73	0.54
1:D:610:GLY:O	1:D:614:ILE:HG22	2.08	0.54
1:B:486:ARG:NH2	2:B:701:HOH:O	2.23	0.53
1:A:392:ASP:OD1	1:A:394:SER:OG	2.25	0.53
1:B:457:THR:O	1:B:460:VAL:HG12	2.09	0.53
1:D:417:ARG:HA	1:D:420:VAL:HG22	1.92	0.52
1:C:447:ALA:HA	1:C:452:LEU:HD12	1.90	0.52
1:D:538:THR:OG1	1:D:541:GLU:HG3	2.09	0.52
1:A:538:THR:HG22	1:A:540:GLY:N	2.26	0.51
1:D:386:LEU:HD12	1:D:387:PRO:HD2	1.92	0.51
1:C:435:GLU:OE1	1:C:625:ARG:NH2	2.43	0.51
1:A:392:ASP:O	1:A:395:LYS:HG3	2.11	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:442:GLN:HA	1:A:445:GLU:OE1	2.11	0.51
1:B:475:ARG:HG3	1:B:475:ARG:HH11	1.76	0.50
1:B:626:ARG:O	1:B:627:LEU:HB2	2.12	0.50
1:A:518:ILE:H	1:A:627:LEU:HD23	1.78	0.49
1:B:475:ARG:HG3	1:B:475:ARG:NH1	2.28	0.48
1:B:500:ASP:OD2	1:C:612:ARG:NH2	2.40	0.48
1:D:445:GLU:O	1:D:449:LYS:HG3	2.13	0.48
1:A:409:LYS:HD3	1:B:485:GLN:OE1	2.14	0.47
1:A:405:GLY:O	1:A:409:LYS:HG3	2.15	0.47
1:A:417:ARG:O	1:A:421:MET:HG3	2.15	0.46
1:D:604:VAL:HG12	1:D:605:ILE:HG12	1.98	0.46
1:D:386:LEU:HD12	1:D:386:LEU:HA	1.73	0.46
1:D:542:MET:HE2	1:D:542:MET:HB2	1.77	0.46
1:B:442:GLN:OE1	1:B:443:GLN:NE2	2.47	0.46
1:D:575:SER:HB2	1:D:595:PRO:HG2	1.97	0.46
1:D:491:LYS:N	1:D:491:LYS:HD2	2.31	0.46
1:A:575:SER:HB2	1:A:595:PRO:HG2	1.98	0.45
1:D:450:ARG:HB2	1:D:452:LEU:HG	1.97	0.45
1:A:579:MET:HB2	1:A:579:MET:HE2	1.72	0.45
1:D:395:LYS:HE2	1:D:395:LYS:HA	1.98	0.45
1:A:404:LEU:HD12	1:B:485:GLN:C	2.42	0.45
1:A:625:ARG:HG3	1:A:625:ARG:NH1	2.32	0.45
1:D:483:ASP:CG	1:D:485:GLN:HG2	2.42	0.45
1:A:538:THR:HG22	1:A:540:GLY:H	1.82	0.45
1:D:456:ILE:H	1:D:456:ILE:HD12	1.82	0.45
1:D:619:ASN:HD22	1:D:619:ASN:N	2.14	0.45
1:A:406:ARG:HH11	1:D:623:ASP:CG	2.24	0.44
1:D:520:GLU:O	1:D:524:GLU:HG3	2.17	0.44
1:C:575:SER:HB2	1:C:595:PRO:HG2	2.00	0.44
1:C:542:MET:HE3	1:C:542:MET:HB3	1.86	0.44
1:B:593:MET:HE3	1:B:593:MET:HB3	1.76	0.44
1:B:459:VAL:O	1:B:463:MET:HG3	2.18	0.43
1:C:460:VAL:HG23	1:C:518:ILE:HD12	1.99	0.43
1:D:464:LYS:HE3	1:D:464:LYS:HB3	1.65	0.43
1:D:465:ALA:HB2	1:D:621:LEU:HG	2.00	0.43
1:D:446:GLU:O	1:D:450:ARG:HG3	2.17	0.43
1:D:431:THR:HG23	1:D:594:LEU:HB3	2.00	0.43
1:D:589:VAL:HG23	1:D:591:ARG:HG2	2.01	0.43
1:B:462:ILE:HG22	1:B:549:ILE:HD13	2.01	0.43
1:A:534:ASP:HB3	1:A:536:LYS:HE3	2.00	0.43
1:A:421:MET:HE2	1:A:421:MET:HB3	1.89	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:506:VAL:HG23	1:B:508:PRO:HD3	2.01	0.42
1:B:483:ASP:O	1:B:485:GLN:HG2	2.20	0.41
1:C:414:ASN:ND2	1:C:417:ARG:HH22	2.16	0.41
1:C:457:THR:O	1:C:460:VAL:HG22	2.20	0.41
1:D:462:ILE:HG12	1:D:621:LEU:HD11	2.02	0.41
1:D:613:PHE:CE2	1:D:617:ILE:HD11	2.54	0.41
1:A:537:LEU:HD12	1:A:537:LEU:HA	1.88	0.41
1:B:626:ARG:NH1	2:B:706:HOH:O	2.45	0.41
1:A:452:LEU:O	1:A:452:LEU:HD23	2.20	0.41
1:B:530:LYS:HD2	1:B:533:ARG:NH2	2.33	0.41
1:C:430:LYS:HE3	2:C:810:HOH:O	2.20	0.41
1:C:431:THR:HG21	1:C:621:LEU:HD23	2.03	0.41
1:D:406:ARG:HA	1:D:409:LYS:HD3	2.01	0.41
1:D:435:GLU:OE1	1:D:591:ARG:NH2	2.54	0.41
1:B:531:LYS:HG2	1:B:536:LYS:HB2	2.03	0.40
1:A:542:MET:HE3	1:A:542:MET:HB2	1.86	0.40
1:B:618:ASN:OD1	1:B:618:ASN:C	2.65	0.40
1:D:406:ARG:O	1:D:410:ILE:HD12	2.21	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	243/247 (98%)	235 (97%)	8 (3%)	0	100	100
1	B	243/247 (98%)	239 (98%)	4 (2%)	0	100	100
1	C	244/247 (99%)	239 (98%)	5 (2%)	0	100	100
1	D	243/247 (98%)	234 (96%)	8 (3%)	1 (0%)	30	22
All	All	973/988 (98%)	947 (97%)	25 (3%)	1 (0%)	48	40

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	557	THR

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	209/210 (100%)	207 (99%)	2 (1%)	68	70
1	B	209/210 (100%)	208 (100%)	1 (0%)	81	84
1	C	210/210 (100%)	209 (100%)	1 (0%)	81	84
1	D	209/210 (100%)	208 (100%)	1 (0%)	81	84
All	All	837/840 (100%)	832 (99%)	5 (1%)	78	81

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	391	VAL
1	A	446	GLU
1	B	559	HIS
1	C	589	VAL
1	D	391	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	472	GLN
1	A	559	HIS
1	C	414	ASN
1	C	485	GLN
1	D	619	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	245/247 (99%)	0.23	9 (3%) 45 48	27, 46, 80, 94	0
1	B	245/247 (99%)	0.04	4 (1%) 70 74	26, 45, 73, 102	0
1	C	245/247 (99%)	0.12	8 (3%) 49 52	26, 47, 91, 141	1 (0%)
1	D	245/247 (99%)	1.64	74 (30%) 1 1	55, 87, 118, 157	0
All	All	980/988 (99%)	0.51	95 (9%) 13 14	26, 53, 106, 157	1 (0%)

All (95) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	448	ALA	4.9
1	D	383	PRO	4.4
1	D	627	LEU	4.2
1	B	383	PRO	4.1
1	C	627	LEU	4.0
1	D	386	LEU	3.9
1	A	627	LEU	3.9
1	D	391	VAL	3.9
1	D	624	ILE	3.9
1	D	452	LEU	3.7
1	D	484	GLY	3.7
1	C	383	PRO	3.6
1	A	386	LEU	3.6
1	D	387	PRO	3.5
1	D	547	PHE	3.5
1	A	388	TRP	3.4
1	D	407	ILE	3.4
1	B	452	LEU	3.4
1	D	456	ILE	3.3
1	D	497	VAL	3.3
1	D	388	TRP	3.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	393	PHE	3.2
1	D	510	PHE	3.2
1	D	480	LEU	3.1
1	D	389	PRO	3.1
1	D	461	PHE	3.0
1	A	383	PRO	3.0
1	C	389	PRO	3.0
1	D	553	GLY	3.0
1	A	625	ARG	3.0
1	D	506	VAL	3.0
1	C	386	LEU	2.9
1	D	454	VAL	2.9
1	D	399	ILE	2.9
1	D	415	LEU	2.9
1	D	488	THR	2.8
1	D	564	VAL	2.8
1	D	434	THR	2.8
1	D	604	VAL	2.8
1	C	388	TRP	2.8
1	D	487	LEU	2.7
1	C	626	ARG	2.7
1	D	447	ALA	2.7
1	C	384	GLY	2.6
1	D	546	CYS	2.6
1	D	616	ILE	2.6
1	D	402	VAL	2.6
1	D	413	ALA	2.6
1	D	566	ALA	2.5
1	D	555	LEU	2.5
1	D	433	ILE	2.5
1	A	448	ALA	2.5
1	D	459	VAL	2.5
1	D	525	LEU	2.5
1	D	621	LEU	2.5
1	D	489	LEU	2.4
1	D	563	ILE	2.4
1	D	419	TRP	2.4
1	D	549	ILE	2.4
1	D	552	ILE	2.4
1	D	527	THR	2.4
1	A	626	ARG	2.4
1	D	420	VAL	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	507	VAL	2.3
1	D	404	LEU	2.3
1	D	439	PHE	2.3
1	D	605	ILE	2.3
1	D	460	VAL	2.3
1	D	425	VAL	2.2
1	D	582	VAL	2.2
1	D	385	MET	2.2
1	B	386	LEU	2.2
1	C	624	ILE	2.2
1	B	384	GLY	2.2
1	D	384	GLY	2.2
1	D	478	SER	2.2
1	D	396	PHE	2.2
1	A	454	VAL	2.2
1	D	620	THR	2.2
1	D	508	PRO	2.2
1	D	519	ILE	2.1
1	D	509	VAL	2.1
1	D	503	ASN	2.1
1	D	539	ALA	2.1
1	D	571	ILE	2.1
1	D	521	LEU	2.1
1	D	528	ILE	2.1
1	D	538	THR	2.1
1	D	466	VAL	2.1
1	D	436	LEU	2.1
1	A	407	ILE	2.0
1	D	573	GLY	2.0
1	D	476	PHE	2.0
1	D	600	PHE	2.0
1	D	589	VAL	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.