



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:43 AM UTC

PDB ID : 4KWV / pdb_00004kwv
Title : Crystal Structure of human apo-QPRT
Authors : Malik, S.S.; Patterson, D.N.; Ncube, Z.; Toth, E.A.
Deposited on : 2013-05-24
Resolution : 2.80 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

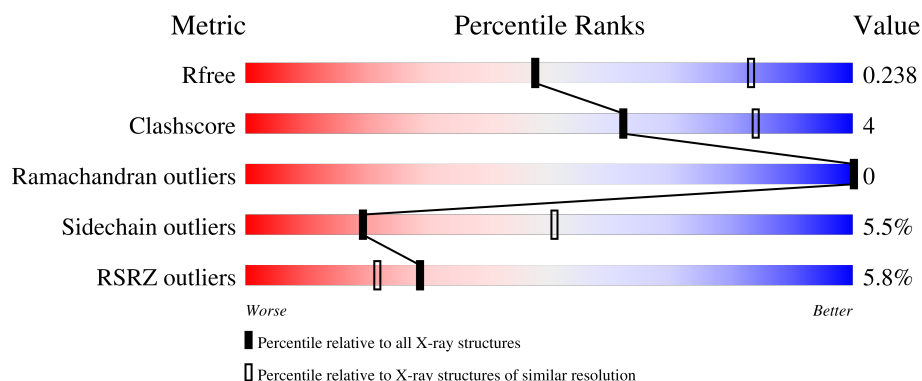
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	301	5% 84% 9% 5%
1	B	301	4% 86% 9% 5%
1	C	301	11% 85% 9% 5%
1	D	301	4% 84% 10% 5%
1	E	301	4% 83% 12% 5%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	301	5% 85% 9% ...

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12001 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 Nicotinate-nucleotide pyrophosphorylase [carboxylating].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	0	0	0
			1995	1281	337	368	9			
1	B	286	Total	C	N	O	S	0	0	0
			1976	1268	330	369	9			
1	C	285	Total	C	N	O	S	0	0	0
			1951	1243	331	368	9			
1	D	285	Total	C	N	O	S	0	0	0
			2005	1285	341	370	9			
1	E	288	Total	C	N	O	S	0	0	0
			2023	1295	345	374	9			
1	F	291	Total	C	N	O	S	0	0	0
			2036	1301	345	381	9			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP Q15274
A	-2	PRO	-	expression tag	UNP Q15274
A	-1	GLY	-	expression tag	UNP Q15274
A	0	SER	-	expression tag	UNP Q15274
B	-3	GLY	-	expression tag	UNP Q15274
B	-2	PRO	-	expression tag	UNP Q15274
B	-1	GLY	-	expression tag	UNP Q15274
B	0	SER	-	expression tag	UNP Q15274
C	-3	GLY	-	expression tag	UNP Q15274
C	-2	PRO	-	expression tag	UNP Q15274
C	-1	GLY	-	expression tag	UNP Q15274
C	0	SER	-	expression tag	UNP Q15274
D	-3	GLY	-	expression tag	UNP Q15274
D	-2	PRO	-	expression tag	UNP Q15274
D	-1	GLY	-	expression tag	UNP Q15274
D	0	SER	-	expression tag	UNP Q15274
E	-3	GLY	-	expression tag	UNP Q15274

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	PRO	-	expression tag	UNP Q15274
E	-1	GLY	-	expression tag	UNP Q15274
E	0	SER	-	expression tag	UNP Q15274
F	-3	GLY	-	expression tag	UNP Q15274
F	-2	PRO	-	expression tag	UNP Q15274
F	-1	GLY	-	expression tag	UNP Q15274
F	0	SER	-	expression tag	UNP Q15274

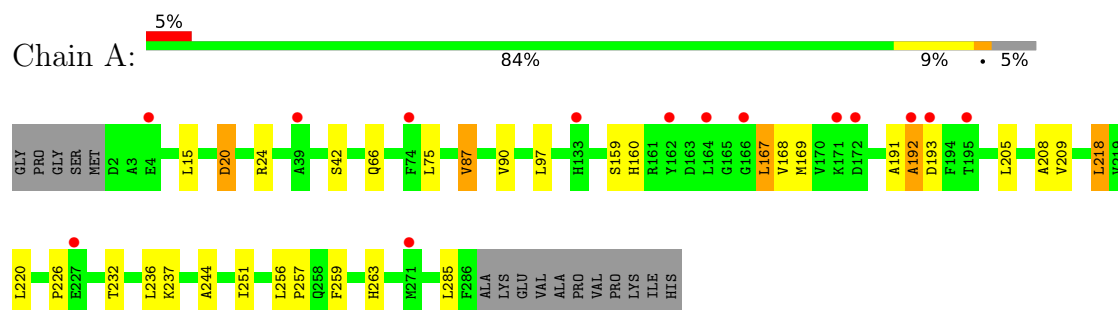
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	4	Total O 4 4	0	0
2	B	1	Total O 1 1	0	0
2	C	1	Total O 1 1	0	0
2	D	3	Total O 3 3	0	0
2	E	1	Total O 1 1	0	0
2	F	5	Total O 5 5	0	0

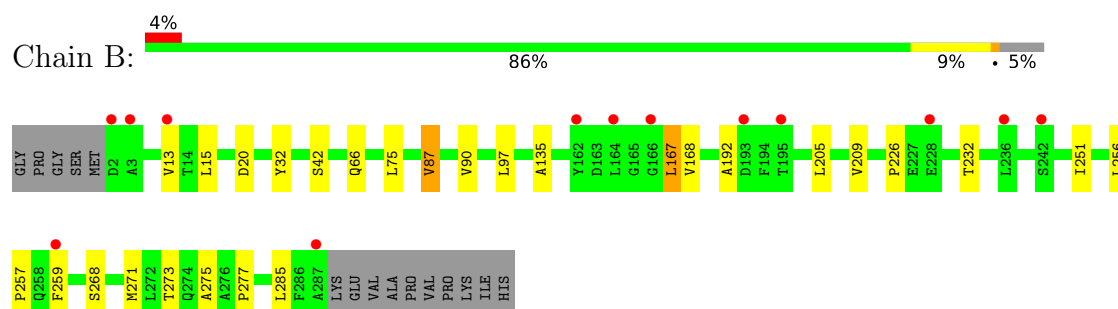
3 Residue-property plots [i](#)

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.

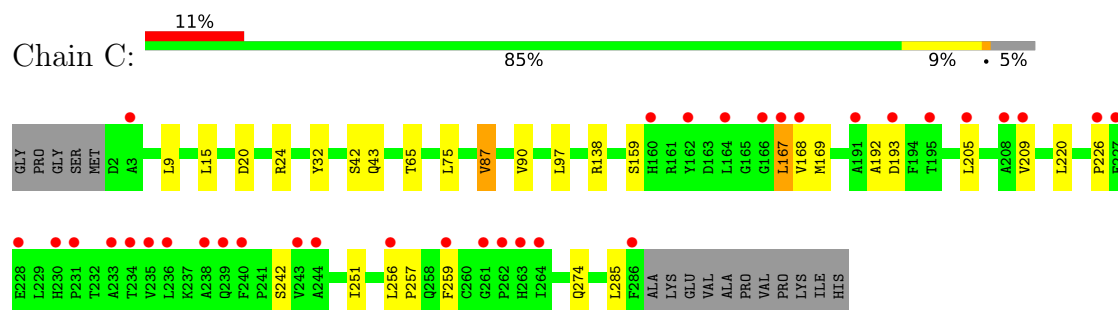
- Molecule 1: Nicotinate-nucleotide pyrophosphorylase [carboxylating]



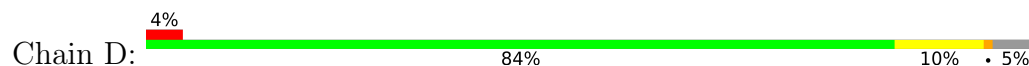
- Molecule 1: Nicotinate-nucleotide pyrophosphorylase [carboxylating]

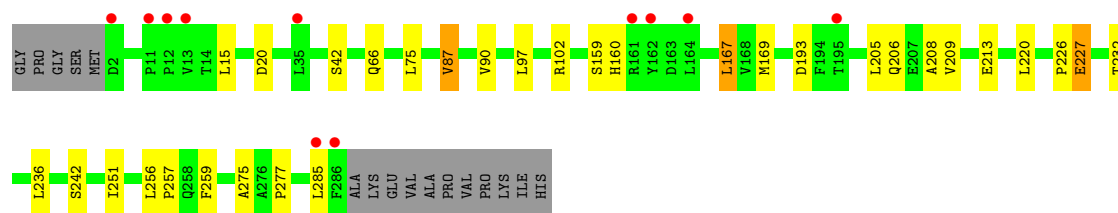


- Molecule 1: Nicotinate-nucleotide pyrophosphorylase [carboxylating]

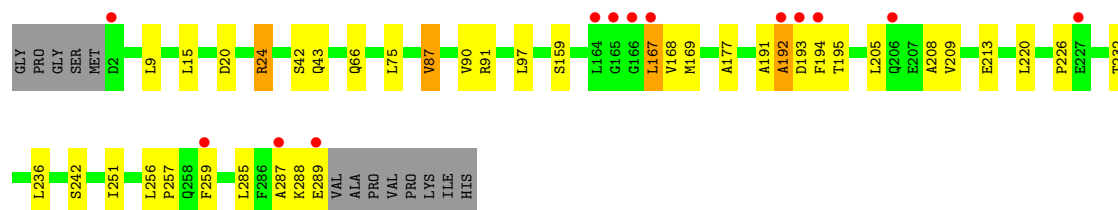
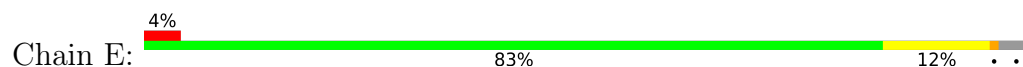


- Molecule 1: Nicotinate-nucleotide pyrophosphorylase [carboxylating]

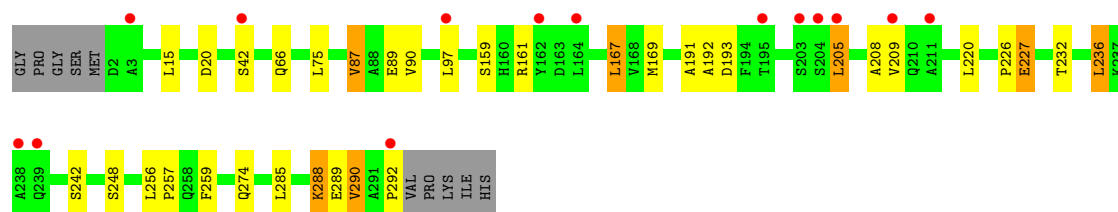
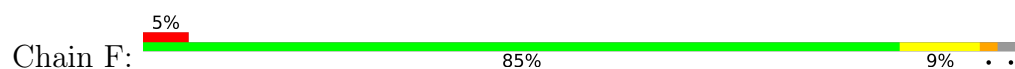




● Molecule 1: Nicotinate-nucleotide pyrophosphorylase [carboxylating]



● Molecule 1: Nicotinate-nucleotide pyrophosphorylase [carboxylating]



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	178.56Å 178.56Å 121.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.93 – 2.80 42.93 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (42.93-2.80) 99.8 (42.93-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.212 , 0.237 0.212 , 0.238	Depositor DCC
R_{free} test set	2812 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	54.4	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 40.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.021 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12001	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 2.71% 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	0.50	0/2038	0.76	2/2796 (0.1%)
1	B	0.55	0/2018	0.77	0/2773
1	C	0.55	0/1990	0.78	1/2731 (0.0%)
1	D	0.57	0/2048	0.75	0/2806
1	E	0.51	0/2065	0.78	2/2829 (0.1%)
1	F	0.72	2/2078 (0.1%)	0.81	3/2849 (0.1%)
All	All	0.57	2/12237 (0.0%)	0.77	8/16784 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	89	GLU	C-O	-5.25	1.17	1.23
1	F	289	GLU	C-O	-5.08	1.17	1.23

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	292	PRO	N-CA-CB	5.95	109.54	103.00
1	A	192	ALA	CB-CA-C	5.72	118.91	109.70
1	E	192	ALA	CB-CA-C	5.61	117.51	109.71
1	F	274	GLN	CB-CA-C	5.35	120.49	110.70
1	C	274	GLN	CB-CA-C	5.33	120.46	110.70
1	F	290	VAL	CB-CA-C	-5.31	103.81	111.34
1	E	287	ALA	N-CA-C	5.27	116.81	108.96
1	A	192	ALA	N-CA-C	-5.18	102.94	110.24

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	1995	0	1941	16	0
1	B	1976	0	1899	16	0
1	C	1951	0	1861	13	0
1	D	2005	0	1962	17	0
1	E	2023	0	1982	27	0
1	F	2036	0	1986	14	0
2	A	4	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	3	0	0	0	0
2	E	1	0	0	0	0
2	F	5	0	0	0	0
All	All	12001	0	11631	97	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 4.

All (97) 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:168:VAL:HG21	1:E:192:ALA:HB2	1.31	1.13
1:D:275:ALA:O	1:D:277:PRO:HD3	1.78	0.83
1:B:275:ALA:O	1:B:277:PRO:HD3	1.79	0.82
1:E:168:VAL:CG2	1:E:192:ALA:HB2	2.11	0.80
1:E:168:VAL:HG21	1:E:192:ALA:CB	2.10	0.79
1:F:20:ASP:OD1	1:F:66:GLN:NE2	2.21	0.73
1:D:20:ASP:OD1	1:D:66:GLN:NE2	2.22	0.72
1:A:208:ALA:HB1	1:A:236:LEU:HD11	1.71	0.72
1:E:20:ASP:OD1	1:E:66:GLN:NE2	2.22	0.71
1:E:43:GLN:HG2	1:E:91:ARG:HG2	1.72	0.71
1:B:20:ASP:OD1	1:B:66:GLN:NE2	2.24	0.70
1:E:208:ALA:HB1	1:E:236:LEU:HD11	1.75	0.66
1:A:20:ASP:OD1	1:A:66:GLN:NE2	2.29	0.65
1:B:168:VAL:HG21	1:B:192:ALA:HB2	1.79	0.64
1:E:193:ASP:OD1	1:E:195:THR:HG23	1.96	0.64

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:167:LEU:HD23	1:E:195:THR:HB	1.83	0.61
1:D:208:ALA:HB1	1:D:236:LEU:HD11	1.86	0.57
1:E:167:LEU:C	1:E:167:LEU:HD12	2.31	0.56
1:C:168:VAL:HG21	1:C:192:ALA:HB2	1.88	0.56
1:A:167:LEU:C	1:A:167:LEU:HD12	2.32	0.55
1:A:191:ALA:O	1:A:192:ALA:C	2.49	0.55
1:E:191:ALA:O	1:E:192:ALA:C	2.50	0.55
1:A:168:VAL:HG21	1:A:192:ALA:HB2	1.88	0.55
1:D:167:LEU:C	1:D:167:LEU:HD12	2.33	0.54
1:E:177:ALA:HA	1:F:288:LYS:HG3	1.90	0.53
1:B:167:LEU:C	1:B:167:LEU:HD12	2.34	0.53
1:F:167:LEU:HD12	1:F:167:LEU:C	2.33	0.53
1:C:205:LEU:O	1:C:209:VAL:HG23	2.09	0.53
1:E:194:PHE:CE2	1:E:195:THR:HG22	2.44	0.52
1:C:167:LEU:HD12	1:C:167:LEU:C	2.34	0.52
1:C:90:VAL:HG12	1:C:97:LEU:HD21	1.93	0.51
1:A:205:LEU:O	1:A:209:VAL:HG23	2.10	0.51
1:F:90:VAL:HG12	1:F:97:LEU:HD21	1.93	0.50
1:E:90:VAL:HG12	1:E:97:LEU:HD21	1.93	0.50
1:B:90:VAL:HG12	1:B:97:LEU:HD21	1.93	0.50
1:D:90:VAL:HG12	1:D:97:LEU:HD21	1.94	0.50
1:B:32:TYR:HA	1:C:9:LEU:CD2	2.42	0.50
1:E:205:LEU:O	1:E:209:VAL:HG23	2.12	0.49
1:B:75:LEU:HD12	1:B:87:VAL:HG13	1.94	0.49
1:F:205:LEU:HB2	1:F:232:THR:HG23	1.93	0.49
1:A:90:VAL:HG12	1:A:97:LEU:HD21	1.94	0.49
1:A:218:LEU:HG	1:A:244:ALA:HB3	1.95	0.48
1:F:75:LEU:HD12	1:F:87:VAL:HG13	1.95	0.48
1:D:205:LEU:O	1:D:209:VAL:HG23	2.13	0.48
1:D:75:LEU:HD12	1:D:87:VAL:HG13	1.96	0.48
1:F:191:ALA:O	1:F:192:ALA:C	2.55	0.47
1:A:20:ASP:OD2	1:A:24:ARG:NH2	2.48	0.47
1:B:205:LEU:O	1:B:209:VAL:HG23	2.13	0.47
1:F:226:PRO:HG3	1:F:259:PHE:CE1	2.50	0.47
1:F:208:ALA:HB1	1:F:236:LEU:HD21	1.96	0.47
1:A:75:LEU:HD12	1:A:87:VAL:HG13	1.96	0.47
1:D:227:GLU:H	1:D:227:GLU:CD	2.21	0.47
1:D:226:PRO:HG3	1:D:259:PHE:CE1	2.51	0.46
1:C:75:LEU:HD12	1:C:87:VAL:HG13	1.97	0.46
1:C:226:PRO:HG3	1:C:259:PHE:CE1	2.51	0.46
1:B:226:PRO:HG3	1:B:259:PHE:CE1	2.50	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:226:PRO:HG3	1:E:259:PHE:CE1	2.51	0.46
1:E:75:LEU:HD12	1:E:87:VAL:HG13	1.98	0.46
1:C:138:ARG:CB	1:D:102:ARG:HD3	2.45	0.45
1:E:168:VAL:CG2	1:E:192:ALA:CB	2.85	0.45
1:A:226:PRO:HG3	1:A:259:PHE:CE1	2.52	0.45
1:D:205:LEU:HB2	1:D:232:THR:HG23	1.98	0.45
1:B:13:VAL:HG12	1:E:24:ARG:HD2	1.97	0.45
1:F:169:MET:HE1	1:F:220:LEU:HD23	1.98	0.45
1:C:32:TYR:HA	1:E:9:LEU:CD2	2.47	0.45
1:E:288:LYS:O	1:E:289:GLU:C	2.60	0.45
1:B:205:LEU:HB2	1:B:232:THR:HG23	1.99	0.44
1:D:169:MET:HE1	1:D:220:LEU:HD23	2.00	0.44
1:E:193:ASP:OD1	1:E:194:PHE:CD1	2.70	0.44
1:F:220:LEU:HD11	1:F:248:SER:HB3	2.00	0.43
1:A:205:LEU:HB2	1:A:232:THR:HG23	1.99	0.42
1:F:205:LEU:O	1:F:209:VAL:HG23	2.19	0.42
1:B:251:ILE:HG12	1:B:259:PHE:CE2	2.54	0.42
1:E:193:ASP:OD1	1:E:194:PHE:CE1	2.72	0.42
1:C:251:ILE:HG12	1:C:259:PHE:CE2	2.55	0.42
1:B:135:ALA:O	1:B:268:SER:HA	2.20	0.42
1:F:256:LEU:N	1:F:257:PRO:CD	2.82	0.42
1:F:227:GLU:H	1:F:227:GLU:CD	2.28	0.42
1:B:256:LEU:N	1:B:257:PRO:CD	2.83	0.41
1:A:237:LYS:HG3	1:A:263:HIS:HB3	2.02	0.41
1:D:251:ILE:HG12	1:D:259:PHE:CE2	2.55	0.41
1:E:256:LEU:N	1:E:257:PRO:CD	2.83	0.41
1:C:169:MET:HE1	1:C:220:LEU:HD23	2.01	0.41
1:E:169:MET:HE1	1:E:220:LEU:HD23	2.03	0.41
1:C:256:LEU:N	1:C:257:PRO:CD	2.84	0.41
1:E:205:LEU:HB2	1:E:232:THR:HG23	2.01	0.41
1:E:251:ILE:HG12	1:E:259:PHE:CE2	2.54	0.41
1:A:256:LEU:N	1:A:257:PRO:CD	2.84	0.41
1:B:271:MET:C	1:B:273:THR:N	2.77	0.41
1:A:251:ILE:HG12	1:A:259:PHE:CE2	2.55	0.41
1:B:251:ILE:HG12	1:B:259:PHE:CD2	2.56	0.40
1:D:205:LEU:HD13	1:D:232:THR:HG23	2.03	0.40
1:C:251:ILE:HG12	1:C:259:PHE:CD2	2.56	0.40
1:D:256:LEU:N	1:D:257:PRO:CD	2.84	0.40
1:E:251:ILE:HG12	1:E:259:PHE:CD2	2.56	0.40
1:A:169:MET:HE1	1:A:220:LEU:HD23	2.03	0.40
1:D:251:ILE:HG12	1:D:259:PHE:CD2	2.56	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	283/301 (94%)	282 (100%)	1 (0%)	0	100	100
1	B	284/301 (94%)	282 (99%)	2 (1%)	0	100	100
1	C	283/301 (94%)	281 (99%)	2 (1%)	0	100	100
1	D	283/301 (94%)	282 (100%)	1 (0%)	0	100	100
1	E	286/301 (95%)	284 (99%)	2 (1%)	0	100	100
1	F	289/301 (96%)	286 (99%)	3 (1%)	0	100	100
All	All	1708/1806 (95%)	1697 (99%)	11 (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	188/222 (85%)	178 (95%)	10 (5%)	20	52
1	B	183/222 (82%)	178 (97%)	5 (3%)	39	74
1	C	179/222 (81%)	167 (93%)	12 (7%)	15	42
1	D	191/222 (86%)	179 (94%)	12 (6%)	16	45
1	E	192/222 (86%)	183 (95%)	9 (5%)	23	57
1	F	195/222 (88%)	181 (93%)	14 (7%)	13	39

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1128/1332 (85%)	1066 (94%)	62 (6%)	19 51

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	LEU
1	A	20	ASP
1	A	42	SER
1	A	87	VAL
1	A	159	SER
1	A	160	HIS
1	A	167	LEU
1	A	193	ASP
1	A	218	LEU
1	A	285	LEU
1	B	15	LEU
1	B	42	SER
1	B	87	VAL
1	B	167	LEU
1	B	285	LEU
1	C	15	LEU
1	C	20	ASP
1	C	24	ARG
1	C	42	SER
1	C	43	GLN
1	C	65	THR
1	C	87	VAL
1	C	159	SER
1	C	167	LEU
1	C	193	ASP
1	C	242	SER
1	C	285	LEU
1	D	15	LEU
1	D	42	SER
1	D	87	VAL
1	D	159	SER
1	D	160	HIS
1	D	167	LEU
1	D	193	ASP
1	D	206	GLN
1	D	213	GLU
1	D	227	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	242	SER
1	D	285	LEU
1	E	15	LEU
1	E	24	ARG
1	E	42	SER
1	E	87	VAL
1	E	159	SER
1	E	167	LEU
1	E	213	GLU
1	E	242	SER
1	E	285	LEU
1	F	15	LEU
1	F	42	SER
1	F	87	VAL
1	F	159	SER
1	F	161	ARG
1	F	167	LEU
1	F	193	ASP
1	F	205	LEU
1	F	227	GLU
1	F	236	LEU
1	F	242	SER
1	F	285	LEU
1	F	288	LYS
1	F	290	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	174	HIS
1	A	210	GLN
1	B	66	GLN
1	B	95	HIS
1	B	239	GLN
1	B	263	HIS
1	C	43	GLN
1	C	66	GLN
1	C	70	GLN
1	C	95	HIS
1	D	66	GLN
1	D	95	HIS
1	E	66	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	E	239	GLN
1	F	66	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](#)

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

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	285/301 (94%)	0.51	14 (4%)	35 27	39, 63, 83, 116	0
1	B	286/301 (95%)	0.49	13 (4%)	38 30	40, 63, 89, 107	0
1	C	285/301 (94%)	0.64	34 (11%)	9 7	33, 63, 104, 120	0
1	D	285/301 (94%)	0.32	11 (3%)	43 34	36, 52, 76, 100	0
1	E	288/301 (95%)	0.25	13 (4%)	38 30	30, 51, 83, 106	0
1	F	291/301 (96%)	0.17	14 (4%)	35 28	27, 44, 70, 90	0
All	All	1720/1806 (95%)	0.40	99 (5%)	29 22	27, 56, 90, 120	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	162	TYR	6.2
1	C	263	HIS	5.6
1	E	193	ASP	4.9
1	D	2	ASP	4.5
1	C	166	GLY	4.2
1	A	193	ASP	4.0
1	C	205	LEU	4.0
1	B	287	ALA	3.8
1	A	39	ALA	3.8
1	E	289	GLU	3.7
1	A	4	GLU	3.6
1	C	233	ALA	3.6
1	C	230	HIS	3.5
1	E	192	ALA	3.4
1	E	2	ASP	3.3
1	C	160	HIS	3.3
1	D	35	LEU	3.3
1	F	162	TYR	3.3
1	F	292	PRO	3.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	F	205	LEU	3.2
1	B	162	TYR	3.2
1	F	195	THR	3.1
1	C	240	PHE	3.1
1	B	193	ASP	3.0
1	F	3	ALA	3.0
1	C	261	GLY	3.0
1	C	167	LEU	3.0
1	A	162	TYR	3.0
1	D	164	LEU	2.9
1	C	262	PRO	2.8
1	A	195	THR	2.8
1	A	227	GLU	2.8
1	C	239	GLN	2.8
1	D	12	PRO	2.7
1	C	286	PHE	2.7
1	D	162	TYR	2.7
1	E	164	LEU	2.7
1	B	228	GLU	2.7
1	D	13	VAL	2.6
1	D	286	PHE	2.6
1	A	74	PHE	2.6
1	C	195	THR	2.6
1	C	234	THR	2.6
1	E	227	GLU	2.6
1	B	166	GLY	2.6
1	A	192	ALA	2.5
1	F	238	ALA	2.5
1	C	264	ILE	2.5
1	F	203	SER	2.5
1	A	172	ASP	2.5
1	F	209	VAL	2.5
1	A	171	LYS	2.4
1	B	236	LEU	2.4
1	C	243	VAL	2.4
1	A	164	LEU	2.4
1	C	164	LEU	2.4
1	C	256	LEU	2.4
1	C	193	ASP	2.4
1	C	235	VAL	2.4
1	C	3	ALA	2.4
1	B	242	SER	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	161	ARG	2.3
1	B	164	LEU	2.3
1	F	204	SER	2.3
1	C	168	VAL	2.3
1	E	165	GLY	2.3
1	B	195	THR	2.3
1	D	11	PRO	2.3
1	E	206	GLN	2.2
1	F	239	GLN	2.2
1	E	194	PHE	2.2
1	C	191	ALA	2.2
1	C	244	ALA	2.2
1	E	287	ALA	2.2
1	A	166	GLY	2.2
1	F	164	LEU	2.2
1	B	259	PHE	2.2
1	C	238	ALA	2.2
1	F	211	ALA	2.2
1	B	2	ASP	2.2
1	D	195	THR	2.1
1	C	231	PRO	2.1
1	A	271	MET	2.1
1	F	42	SER	2.1
1	C	208	ALA	2.1
1	C	228	GLU	2.1
1	F	97	LEU	2.1
1	E	259	PHE	2.1
1	C	226	PRO	2.1
1	E	167	LEU	2.1
1	C	259	PHE	2.1
1	C	236	LEU	2.1
1	E	166	GLY	2.1
1	B	13	VAL	2.1
1	B	3	ALA	2.0
1	D	285	LEU	2.0
1	A	133	HIS	2.0
1	C	209	VAL	2.0
1	C	227	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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.