Full wwPDB X-ray Structure Validation Report (i)

Oct 27, 2023 – 05:13 AM EDT

PDB ID	:	3NX1
Title	:	Crystal structure of Enterobacter sp. Px6-4 Ferulic Acid Decarboxylase
Authors	:	Gu, W.; Yang, J.K.; Lou, Z.Y.; Meng, Z.H.; Zhang, K.-Q.
Deposited on	:	2010-07-12
Resolution	:	2.40 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 (i) 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 (1)) were used in the production of this report:

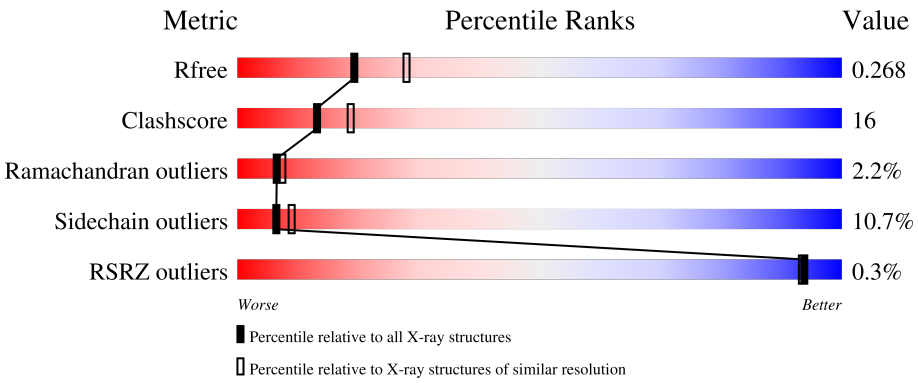
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.40 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}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 <=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	168	61% 25% 9% • •
1	B	168	% 57% 30% 8% • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 2724 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 Ferulic acid decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	162	1301	830	216	250	5	0	0	0
1	B	165	1327	849	219	254	5	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	26	ASN	GLU	SEE REMARK 999	UNP C6F3U5
A	34	ASP	GLU	SEE REMARK 999	UNP C6F3U5
A	37	ILE	LEU	SEE REMARK 999	UNP C6F3U5
A	55	GLU	GLN	SEE REMARK 999	UNP C6F3U5
A	105	PRO	LYS	SEE REMARK 999	UNP C6F3U5
A	122	GLU	ASP	SEE REMARK 999	UNP C6F3U5
B	26	ASN	GLU	SEE REMARK 999	UNP C6F3U5
B	34	ASP	GLU	SEE REMARK 999	UNP C6F3U5
B	37	ILE	LEU	SEE REMARK 999	UNP C6F3U5
B	55	GLU	GLN	SEE REMARK 999	UNP C6F3U5
B	105	PRO	LYS	SEE REMARK 999	UNP C6F3U5
B	122	GLU	ASP	SEE REMARK 999	UNP C6F3U5

- Molecule 2 is water.

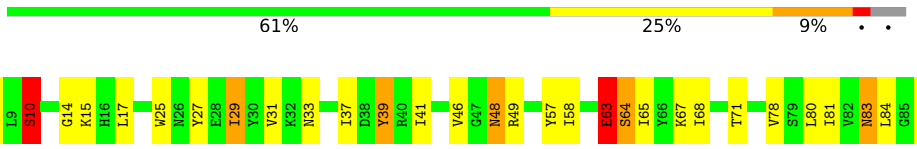
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	42	42	42	0	0
2	B	54	54	54	0	0

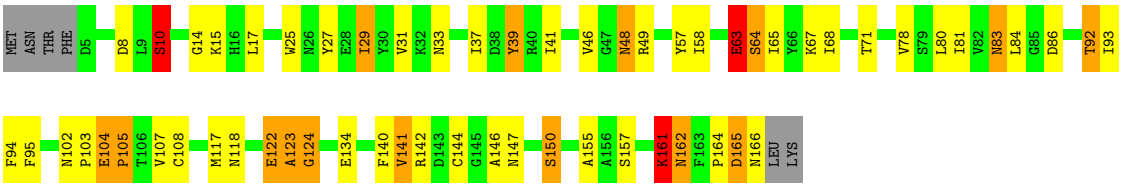
3 Residue-property plots (i)

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: Ferulic acid decarboxylase

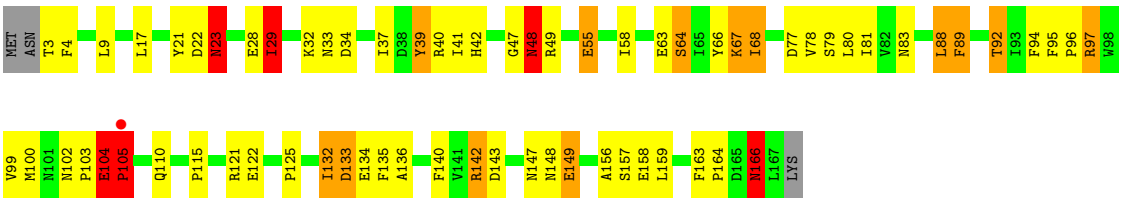
Chain A:





• Molecule 1: Ferulic acid decarboxylase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	41.48Å 88.72Å 49.26Å 90.00° 102.25° 90.00°	Depositor
Resolution (Å)	36.87 – 2.40 36.87 – 2.40	Depositor EDS
% Data completeness (in resolution range)	93.2 (36.87-2.40) 93.2 (36.87-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	6.40 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.162 , 0.262 0.168 , 0.268	Depositor DCC
R_{free} test set	625 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å^2)	35.1	Xtriage
Anisotropy	0.104	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 42.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	2724	wwPDB-VP
Average B, all atoms (Å^2)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.83% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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	1.73	13/1339 (1.0%)	1.40	12/1829 (0.7%)
1	B	1.67	17/1366 (1.2%)	1.43	13/1866 (0.7%)
All	All	1.70	30/2705 (1.1%)	1.41	25/3695 (0.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	2
All	All	0	4

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	64	SER	N-CA	8.21	1.62	1.46
1	A	104	GLU	CG-CD	8.17	1.64	1.51
1	B	89	PHE	CE1-CZ	7.92	1.52	1.37
1	B	55	GLU	CG-CD	7.39	1.63	1.51
1	B	39	TYR	CE2-CZ	6.51	1.47	1.38
1	A	78	VAL	CB-CG1	-6.40	1.39	1.52
1	B	104	GLU	CG-CD	6.27	1.61	1.51
1	A	25	TRP	CZ3-CH2	6.19	1.50	1.40
1	A	31	VAL	CB-CG2	6.10	1.65	1.52
1	B	105	PRO	CG-CD	6.02	1.70	1.50
1	A	164	PRO	CA-C	5.92	1.64	1.52
1	B	104	GLU	CB-CG	5.82	1.63	1.52
1	A	155	ALA	CA-CB	5.65	1.64	1.52
1	B	42	HIS	C-O	-5.65	1.12	1.23
1	A	39	TYR	CB-CG	-5.55	1.43	1.51
1	B	99	VAL	CB-CG2	5.51	1.64	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	146	ALA	CA-CB	5.46	1.64	1.52
1	A	104	GLU	CB-CG	5.44	1.62	1.52
1	B	66	TYR	CG-CD1	5.41	1.46	1.39
1	A	150	SER	CA-CB	5.36	1.60	1.52
1	B	143	ASP	CB-CG	5.30	1.62	1.51
1	B	29	ILE	CA-CB	5.25	1.67	1.54
1	B	94	PHE	CE1-CZ	5.24	1.47	1.37
1	A	10	SER	CB-OG	-5.16	1.35	1.42
1	B	134	GLU	CB-CG	-5.16	1.42	1.52
1	B	55	GLU	CB-CG	5.12	1.61	1.52
1	B	41	ILE	C-O	5.11	1.33	1.23
1	B	149	GLU	CG-CD	5.09	1.59	1.51
1	A	63	GLU	CB-CG	-5.05	1.42	1.52
1	B	115	PRO	CA-C	5.03	1.62	1.52

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	142	ARG	NE-CZ-NH1	-10.25	115.18	120.30
1	A	164	PRO	N-CA-C	8.45	134.08	112.10
1	B	166	ASN	C-N-CA	-8.21	101.17	121.70
1	B	97	ARG	NE-CZ-NH1	-7.63	116.48	120.30
1	A	124	GLY	N-CA-C	-7.63	94.03	113.10
1	B	133	ASP	CB-CG-OD1	-7.58	111.48	118.30
1	B	142	ARG	NE-CZ-NH2	7.12	123.86	120.30
1	A	162	ASN	N-CA-C	7.09	130.15	111.00
1	B	40	ARG	NE-CZ-NH2	-7.04	116.78	120.30
1	B	40	ARG	NE-CZ-NH1	6.99	123.79	120.30
1	A	165	ASP	CB-CG-OD2	6.60	124.24	118.30
1	A	64	SER	N-CA-C	6.53	128.62	111.00
1	B	121	ARG	NE-CZ-NH2	6.29	123.44	120.30
1	A	80	LEU	CB-CG-CD1	-6.22	100.43	111.00
1	A	164	PRO	C-N-CA	5.77	136.13	121.70
1	B	67	LYS	CD-CE-NZ	5.66	124.73	111.70
1	A	150	SER	CB-CA-C	5.63	120.79	110.10
1	A	63	GLU	N-CA-CB	-5.57	100.57	110.60
1	B	88	LEU	CB-CG-CD1	-5.41	101.80	111.00
1	B	64	SER	N-CA-C	5.41	125.59	111.00
1	B	121	ARG	NE-CZ-NH1	-5.40	117.60	120.30
1	A	63	GLU	C-N-CA	-5.35	108.33	121.70
1	B	166	ASN	O-C-N	-5.34	114.16	122.70
1	A	117	MET	CG-SD-CE	5.08	108.33	100.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	165	ASP	N-CA-C	-5.04	97.38	111.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	161	LYS	Peptide
1	A	165	ASP	Peptide
1	B	166	ASN	Peptide
1	B	22	ASP	Peptide

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	1301	0	1221	43	1
1	B	1327	0	1248	45	1
2	A	42	0	0	0	0
2	B	54	0	0	1	0
All	All	2724	0	2469	83	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 16.

All (83) 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:A:8:ASP:OD1	1:A:10:SER:HB2	1.62	0.97
1:B:166:ASN:ND2	1:B:166:ASN:H	1.61	0.93
1:A:102:ASN:O	1:A:105:PRO:HD2	1.67	0.92
1:A:102:ASN:O	1:A:105:PRO:CD	2.21	0.89
1:A:33:ASN:HD22	1:A:147:ASN:HD21	1.23	0.85
1:B:166:ASN:ND2	1:B:166:ASN:N	2.29	0.81
1:A:17:LEU:HD23	1:A:141:VAL:HG13	1.64	0.78
1:A:123:ALA:O	1:A:124:GLY:O	2.08	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:PHE:CB	1:B:100:MET:CE	2.69	0.71
1:B:58:ILE:HG12	1:B:68:ILE:HG23	1.73	0.69
1:A:27:TYR:CE2	1:A:41:ILE:HG12	2.27	0.69
1:A:123:ALA:C	1:A:124:GLY:O	2.28	0.68
1:B:95:PHE:CD2	1:B:100:MET:CE	2.77	0.67
1:B:166:ASN:N	1:B:166:ASN:HD22	1.91	0.67
1:B:95:PHE:CB	1:B:100:MET:HE2	2.25	0.66
1:A:103:PRO:C	1:A:105:PRO:HD2	2.15	0.66
1:B:95:PHE:CD2	1:B:100:MET:HE3	2.31	0.65
1:B:95:PHE:HB3	1:B:100:MET:HE2	1.78	0.65
1:A:33:ASN:HD22	1:A:147:ASN:ND2	1.94	0.63
1:B:3:THR:N	2:B:192:HOH:O	2.32	0.62
1:A:92:THR:HG23	1:B:67:LYS:HZ1	1.65	0.62
1:B:88:LEU:HD21	1:B:135:PHE:HE1	1.67	0.59
1:A:29:ILE:HD11	1:A:37:ILE:CD1	2.33	0.59
1:B:95:PHE:CG	1:B:100:MET:CE	2.87	0.58
1:B:102:ASN:O	1:B:105:PRO:HD2	2.04	0.57
1:A:17:LEU:CD2	1:A:141:VAL:HG13	2.32	0.57
1:A:83:ASN:ND2	1:A:86:ASP:H	2.03	0.56
1:B:33:ASN:HD22	1:B:147:ASN:HD21	1.54	0.56
1:B:95:PHE:HB2	1:B:100:MET:CE	2.36	0.56
1:B:3:THR:HG22	1:B:4:PHE:HD1	1.70	0.55
1:A:29:ILE:HD11	1:A:37:ILE:HD12	1.88	0.54
1:A:58:ILE:HG12	1:A:68:ILE:HG23	1.89	0.54
1:A:140:PHE:CE2	1:A:142:ARG:HD3	2.43	0.53
1:B:140:PHE:CE2	1:B:142:ARG:HG2	2.43	0.53
1:B:63:GLU:O	1:B:64:SER:HB2	2.09	0.53
1:A:118:ASN:O	1:A:122:GLU:HG2	2.09	0.52
1:A:48:ASN:HD21	1:A:157:SER:H	1.57	0.52
1:A:41:ILE:HD13	1:A:46:VAL:HG23	1.91	0.52
1:B:95:PHE:CD2	1:B:100:MET:HE1	2.45	0.51
1:B:21:TYR:HB3	1:B:23:ASN:HB2	1.92	0.51
1:B:3:THR:HG22	1:B:4:PHE:CD1	2.45	0.51
1:B:49:ARG:NH2	1:B:110:GLN:OE1	2.40	0.50
1:A:83:ASN:C	1:A:83:ASN:HD22	2.15	0.50
1:B:97:ARG:HB3	1:B:125:PRO:O	2.12	0.49
1:A:140:PHE:CZ	1:A:142:ARG:HD3	2.48	0.48
1:B:95:PHE:HB2	1:B:100:MET:HE3	1.95	0.48
1:A:63:GLU:O	1:A:63:GLU:HG3	2.12	0.48
1:B:92:THR:HB	1:B:133:ASP:OD2	2.14	0.48
1:A:107:VAL:O	1:A:108:CYS:HB3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:29:ILE:HD12	1:B:39:TYR:HB2	1.96	0.47
1:B:48:ASN:ND2	1:B:156:ALA:HB3	2.29	0.47
1:B:63:GLU:O	1:B:64:SER:CB	2.62	0.47
1:A:67:LYS:NZ	1:B:92:THR:HG22	2.30	0.47
1:A:37:ILE:HG21	1:A:37:ILE:HD13	1.51	0.46
1:B:32:LYS:NZ	1:B:149:GLU:OE1	2.47	0.46
1:A:104:GLU:N	1:A:105:PRO:CD	2.79	0.46
1:A:48:ASN:HD21	1:A:157:SER:N	2.13	0.46
1:B:78:VAL:CG1	1:B:80:LEU:HD21	2.46	0.46
1:A:94:PHE:CD1	1:A:94:PHE:N	2.83	0.45
1:A:39:TYR:CD1	1:A:39:TYR:C	2.90	0.45
1:A:93:ILE:HG21	1:A:95:PHE:CZ	2.52	0.45
1:A:15:LYS:HE2	1:A:84:LEU:HD13	1.98	0.44
1:A:57:TYR:N	1:A:57:TYR:CD2	2.85	0.44
1:B:163:PHE:HA	1:B:164:PRO:HA	1.70	0.44
1:B:48:ASN:HD21	1:B:157:SER:N	2.16	0.44
1:B:95:PHE:HA	1:B:96:PRO:HD3	1.84	0.44
1:A:94:PHE:CE2	1:B:77:ASP:HB3	2.54	0.43
1:B:17:LEU:O	1:B:28:GLU:HA	2.19	0.42
1:B:103:PRO:HD2	1:B:104:GLU:OE1	2.19	0.42
1:B:83:ASN:HB3	1:B:88:LEU:HB3	2.00	0.42
1:B:88:LEU:HD21	1:B:135:PHE:CE1	2.51	0.42
1:A:161:LYS:HB3	1:A:161:LYS:HE2	1.77	0.42
1:A:104:GLU:N	1:A:105:PRO:HD2	2.34	0.41
1:B:89:PHE:CD2	1:B:136:ALA:HB3	2.54	0.41
1:A:81:ILE:HD11	1:B:81:ILE:HD11	2.03	0.41
1:A:14:GLY:O	1:A:144:CYS:N	2.47	0.41
1:A:81:ILE:CD1	1:B:81:ILE:HD11	2.51	0.41
1:B:132:ILE:O	1:B:132:ILE:HG22	2.19	0.41
1:B:156:ALA:HA	1:B:159:LEU:HD12	2.03	0.41
1:A:27:TYR:CD2	1:A:41:ILE:HG12	2.56	0.41
1:A:41:ILE:HD12	1:A:49:ARG:HB3	2.02	0.41
1:A:29:ILE:CD1	1:A:37:ILE:HG21	2.51	0.40
1:A:81:ILE:HD12	1:A:81:ILE:HG23	1.79	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:ASN:ND2	1:B:148:ASN:ND2[2_556]	2.11	0.09

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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	160/168 (95%)	146 (91%)	12 (8%)	2 (1%)	12 17
1	B	163/168 (97%)	145 (89%)	13 (8%)	5 (3%)	4 3
All	All	323/336 (96%)	291 (90%)	25 (8%)	7 (2%)	6 7

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	162	ASN
1	B	23	ASN
1	B	47	GLY
1	B	166	ASN
1	A	123	ALA
1	B	48	ASN
1	B	105	PRO

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	143/149 (96%)	128 (90%)	15 (10%)	7 9
1	B	146/149 (98%)	130 (89%)	16 (11%)	6 8
All	All	289/298 (97%)	258 (89%)	31 (11%)	6 9

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	SER
1	A	29	ILE
1	A	48	ASN
1	A	63	GLU
1	A	64	SER
1	A	65	ILE
1	A	71	THR
1	A	83	ASN
1	A	92	THR
1	A	105	PRO
1	A	122	GLU
1	A	134	GLU
1	A	141	VAL
1	A	150	SER
1	A	161	LYS
1	B	9	LEU
1	B	23	ASN
1	B	29	ILE
1	B	34	ASP
1	B	37	ILE
1	B	48	ASN
1	B	55	GLU
1	B	68	ILE
1	B	79	SER
1	B	92	THR
1	B	104	GLU
1	B	105	PRO
1	B	122	GLU
1	B	132	ILE
1	B	158	GLU
1	B	166	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	ASN
1	A	83	ASN
1	A	113	HIS
1	A	147	ASN
1	B	23	ASN
1	B	48	ASN
1	B	147	ASN
1	B	166	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 monosaccharides 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	162/168 (96%)	-0.45	0 100 100	12, 26, 38, 52	2 (1%)
1	B	165/168 (98%)	-0.39	1 (0%) 89 88	15, 26, 38, 52	2 (1%)
All	All	327/336 (97%)	-0.42	1 (0%) 94 93	12, 26, 38, 52	4 (1%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	105	PRO	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 monosaccharides 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.