wwPDB X-ray Structure Validation Summary Report (i)

May 15, 2020 – 08:29 am BST

PDB ID	:	4B3K
Title	:	Family 1 6-phospho-beta-D glycosidase from Streptococcus pyogenes
Authors	:	Stepper, J.; Dabin, J.; Ekloef, J.M.; Thongpoo, P.; Kongsaeree, P.T.; Taylor, E.J.; Turkenburg, J.P.; Brumer, H.; Davies, G.J.
Deposited on	:	2012-07-24
Resolution	:	2.60 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 following versions of software and data (see references (1)) were used in the production of this report:

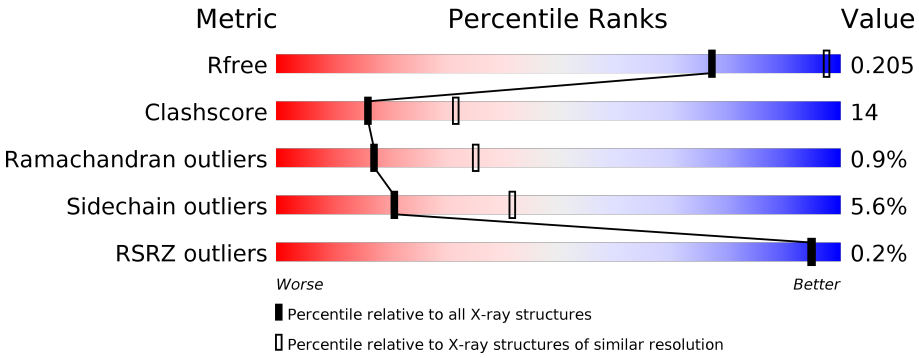
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 2.60 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}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 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	479	69% 24%
1	B	479	68% 26%
1	C	479	64% 29%
1	D	479	67% 25%
1	E	479	61% 32%
1	F	479	% 61% 31%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 22669 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 BETA-GLUCOSIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	461	Total	C	N	O	S	0	0	0
			3767	2435	636	686	10			
1	B	461	Total	C	N	O	S	0	0	0
			3767	2435	636	686	10			
1	C	461	Total	C	N	O	S	0	0	0
			3767	2435	636	686	10			
1	D	461	Total	C	N	O	S	0	0	0
			3767	2435	636	686	10			
1	E	461	Total	C	N	O	S	0	0	0
			3767	2435	636	686	10			
1	F	461	Total	C	N	O	S	0	0	0
			3767	2435	636	686	10			

- Molecule 2 is water.

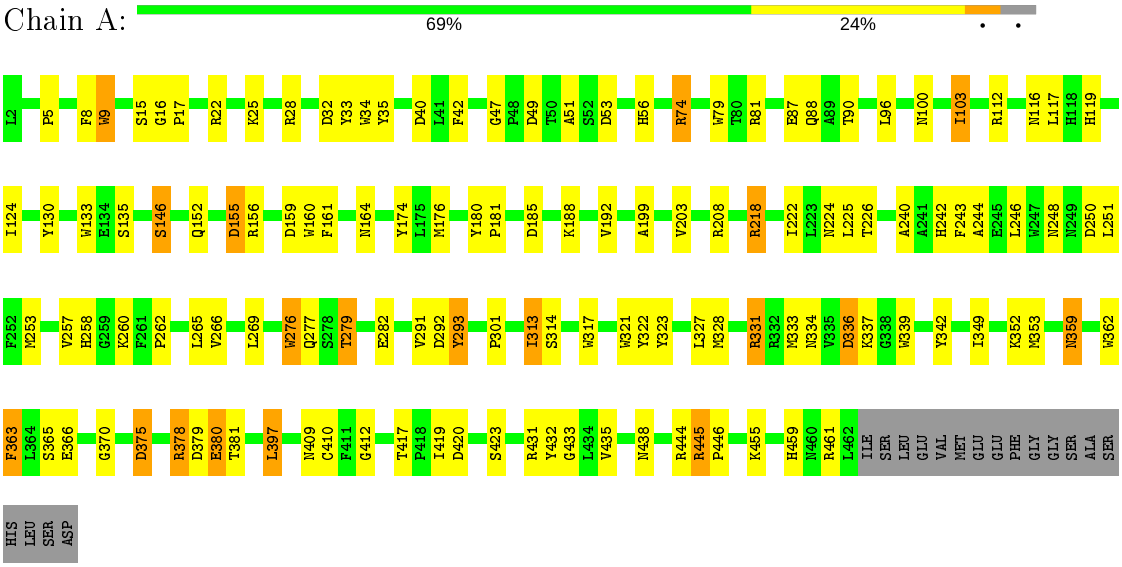
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	13	Total	O	0	0
			13	13		
2	B	17	Total	O	0	0
			17	17		
2	C	10	Total	O	0	0
			10	10		
2	D	14	Total	O	0	0
			14	14		
2	E	9	Total	O	0	0
			9	9		
2	F	4	Total	O	0	0
			4	4		

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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.

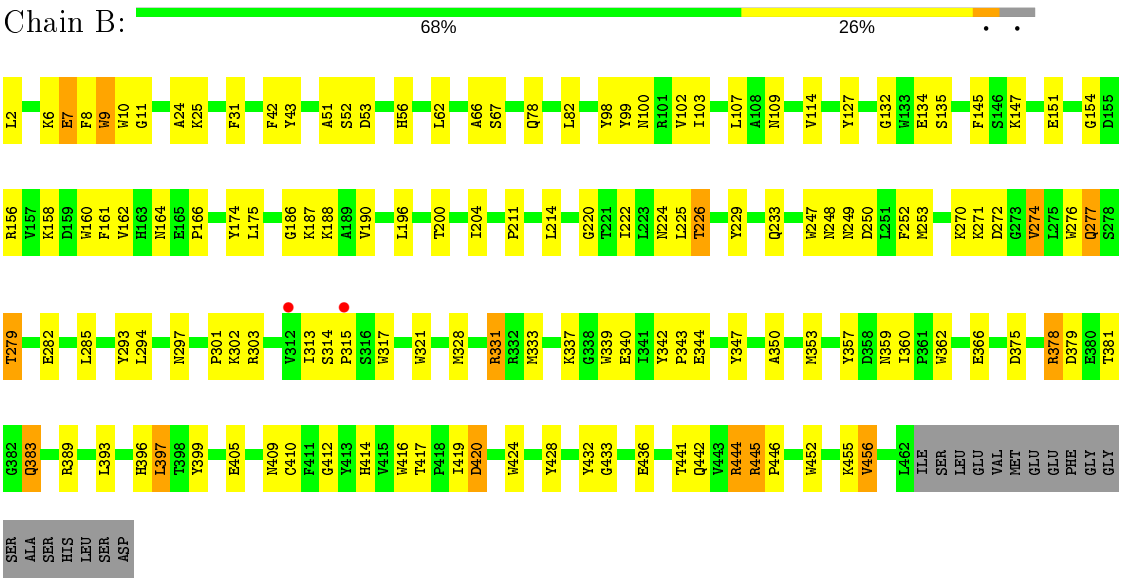
- Molecule 1: BETA-GLUCOSIDASE

Chain A:



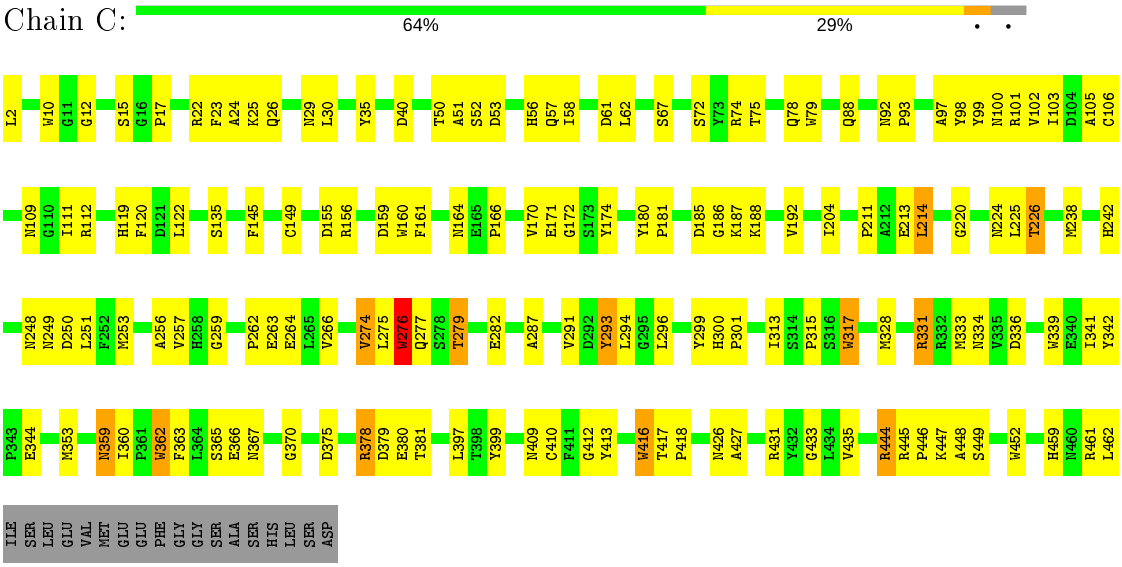
- Molecule 1: BETA-GLUCOSIDASE

Chain B:



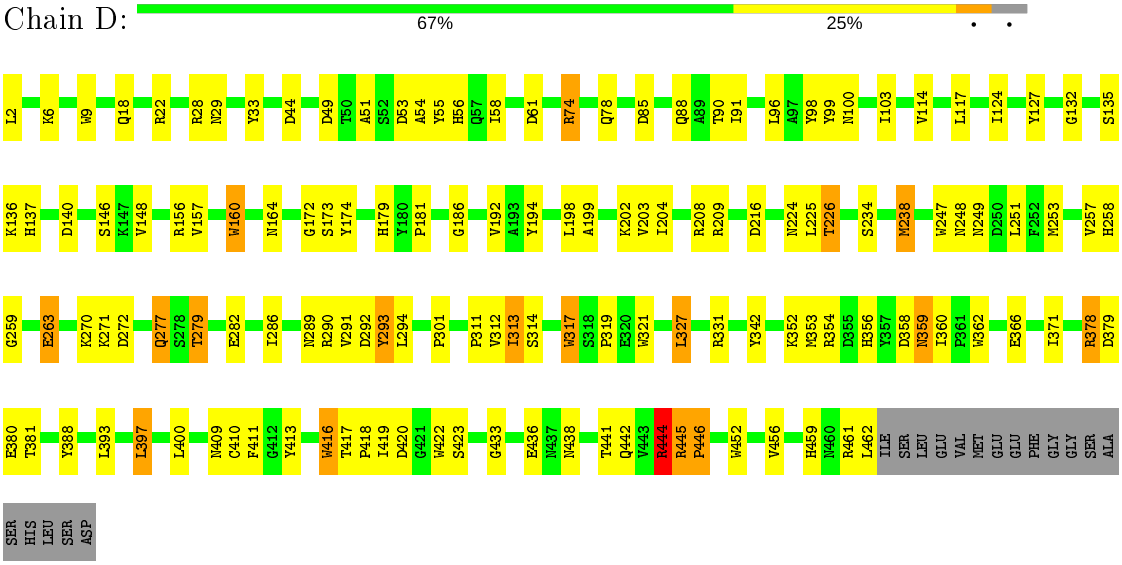
- Molecule 1: BETA-GLUCOSIDASE

Chain C:



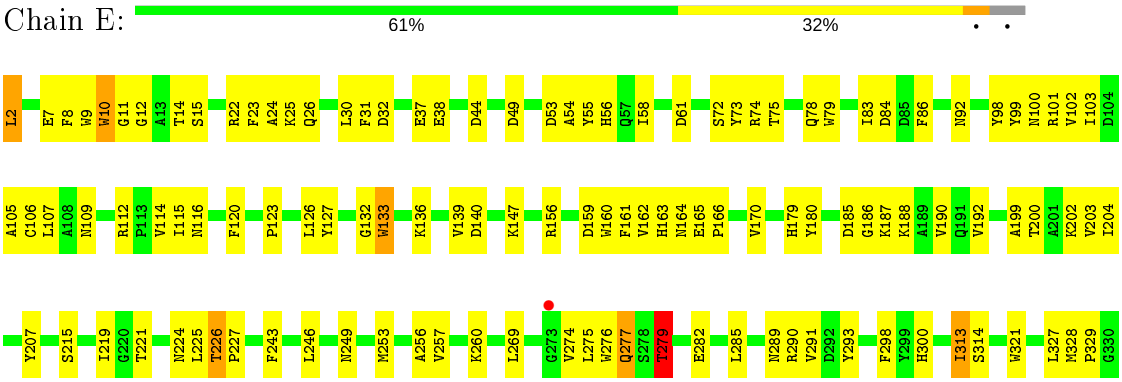
• Molecule 1: BETA-GLUCOSIDASE

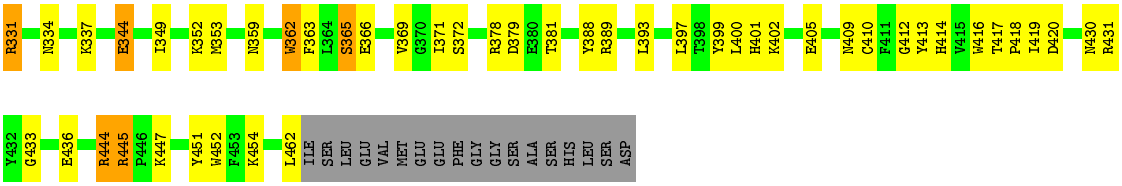
Chain D:



• Molecule 1: BETA-GLUCOSIDASE

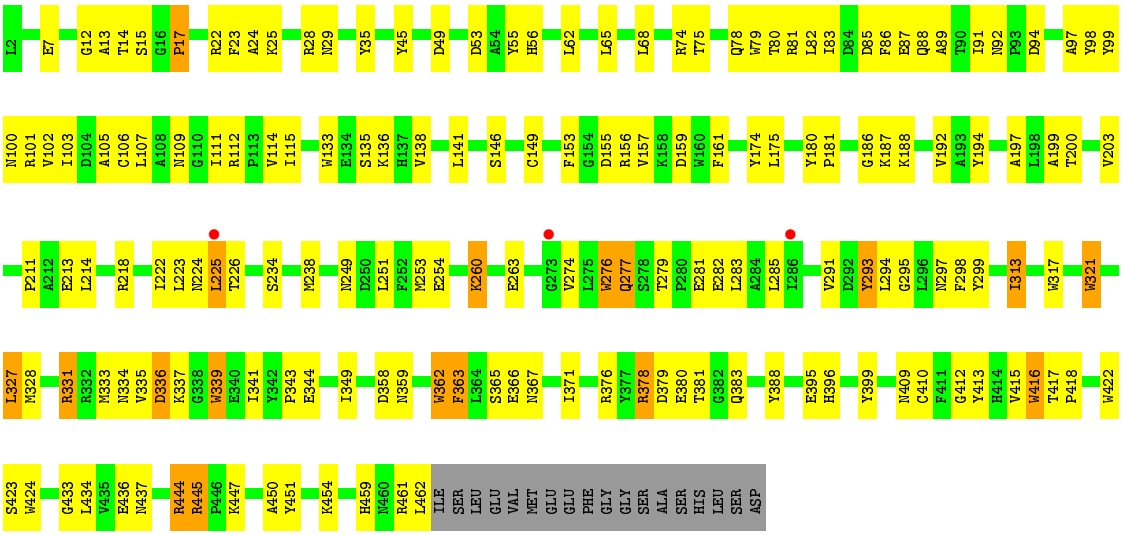
Chain E:





- Molecule 1: BETA-GLUCOSIDASE

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	106.42Å 190.20Å 106.17Å 90.00° 118.88° 90.00°	Depositor
Resolution (Å)	66.56 – 2.60 66.56 – 2.60	Depositor EDS
% Data completeness (in resolution range)	93.7 (66.56-2.60) 93.0 (66.56-2.60)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.94 (at 2.62Å)	Xtriage
Refinement program	REFMAC 5.7.0025	Depositor
R, R_{free}	0.223 , 0.286 0.182 , 0.205	Depositor DCC
R_{free} test set	5282 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	41.5	Xtriage
Anisotropy	0.066	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 5.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.40$, $< L^2 > = 0.23$	Xtriage
Estimated twinning fraction	0.069 for l,k,-h-l 0.069 for -h-l,k,h 0.074 for -h-l,-k,l 0.074 for h,-k,-h-l 0.348 for l,-k,h	Xtriage
Reported twinning fraction	0.677 for H, K, L 0.323 for L, -K, H	Depositor
Outliers	0 of 105767 reflections	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	22669	wwPDB-VP
Average B, all atoms (Å^2)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.41% 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

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	0.81	9/3891 (0.2%)	0.83	2/5304 (0.0%)
1	B	0.81	7/3891 (0.2%)	0.85	1/5304 (0.0%)
1	C	0.79	9/3891 (0.2%)	0.81	2/5304 (0.0%)
1	D	0.78	6/3891 (0.2%)	0.81	5/5304 (0.1%)
1	E	0.77	8/3891 (0.2%)	0.79	1/5304 (0.0%)
1	F	0.74	8/3891 (0.2%)	0.76	0/5304
All	All	0.78	47/23346 (0.2%)	0.81	11/31824 (0.0%)

The worst 5 of 47 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	133	TRP	CD2-CE2	6.40	1.49	1.41
1	D	416	TRP	CD2-CE2	6.38	1.49	1.41
1	A	34	TRP	CD2-CE2	6.27	1.48	1.41
1	B	10	TRP	CD2-CE2	6.10	1.48	1.41
1	C	79	TRP	CD2-CE2	6.07	1.48	1.41

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	444	ARG	NE-CZ-NH2	-6.47	117.07	120.30
1	B	250	ASP	CB-CG-OD1	5.99	123.69	118.30
1	D	444	ARG	NE-CZ-NH1	5.97	123.29	120.30
1	A	28	ARG	NE-CZ-NH2	-5.85	117.37	120.30
1	C	296	LEU	CA-CB-CG	5.81	128.65	115.30

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	3767	0	3574	91	0
1	B	3767	0	3574	94	0
1	C	3767	0	3574	101	0
1	D	3767	0	3574	105	0
1	E	3767	0	3574	126	0
1	F	3767	0	3574	122	0
2	A	13	0	0	2	0
2	B	17	0	0	0	0
2	C	10	0	0	0	0
2	D	14	0	0	2	0
2	E	9	0	0	0	0
2	F	4	0	0	0	0
All	All	22669	0	21444	633	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 14.

The worst 5 of 633 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:224:ASN:OD1	1:D:226:THR:HG23	1.40	1.20
1:A:224:ASN:OD1	1:A:226:THR:HG23	1.44	1.15
1:E:331:ARG:HG2	1:E:331:ARG:HH11	1.05	1.13
1:C:331:ARG:HH11	1:C:331:ARG:HG2	1.13	1.11
1:A:225:LEU:HD22	1:A:353:MET:CE	1.86	1.04

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	459/479 (96%)	423 (92%)	32 (7%)	4 (1%)	17	35
1	B	459/479 (96%)	420 (92%)	34 (7%)	5 (1%)	14	30
1	C	459/479 (96%)	420 (92%)	37 (8%)	2 (0%)	34	57
1	D	459/479 (96%)	416 (91%)	38 (8%)	5 (1%)	14	30
1	E	459/479 (96%)	404 (88%)	51 (11%)	4 (1%)	17	35
1	F	459/479 (96%)	408 (89%)	47 (10%)	4 (1%)	17	35
All	All	2754/2874 (96%)	2491 (90%)	239 (9%)	24 (1%)	17	35

5 of 24 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	24	ALA
1	B	271	LYS
1	D	209	ARG
1	E	344	GLU
1	F	24	ALA

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	389/404 (96%)	365 (94%)	24 (6%)	18	37
1	B	389/404 (96%)	371 (95%)	18 (5%)	27	51
1	C	389/404 (96%)	364 (94%)	25 (6%)	17	35
1	D	389/404 (96%)	367 (94%)	22 (6%)	20	41
1	E	389/404 (96%)	372 (96%)	17 (4%)	28	53
1	F	389/404 (96%)	365 (94%)	24 (6%)	18	37
All	All	2334/2424 (96%)	2204 (94%)	130 (6%)	21	42

5 of 130 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	331	ARG
1	D	238	MET
1	F	336	ASP
1	C	359	ASN
1	C	445	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 78 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	289	ASN
1	D	100	ASN
1	F	359	ASN
1	C	359	ASN
1	C	430	ASN

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

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95^{th} 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	461/479 (96%)	-0.43	0 100 100	17, 32, 49, 65	1 (0%)
1	B	461/479 (96%)	-0.39	2 (0%) 92 91	16, 31, 50, 68	1 (0%)
1	C	461/479 (96%)	-0.39	0 100 100	21, 36, 56, 81	1 (0%)
1	D	461/479 (96%)	-0.42	0 100 100	19, 34, 56, 70	1 (0%)
1	E	461/479 (96%)	-0.19	1 (0%) 95 95	23, 46, 65, 89	1 (0%)
1	F	461/479 (96%)	-0.11	3 (0%) 87 86	28, 49, 74, 91	1 (0%)
All	All	2766/2874 (96%)	-0.32	6 (0%) 95 95	16, 38, 62, 91	6 (0%)

The worst 5 of 6 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	225	LEU	3.3
1	F	273	GLY	3.0
1	B	315	PRO	2.3
1	F	286	ILE	2.1
1	B	312	VAL	2.1

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

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.