



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2018 – 09:41 pm GMT

PDB ID : 2EX2
Title : Crystal structure of penicillin binding protein 4 (dacB) from Escherichia coli
Authors : Kishida, H.; Unzai, S.; Roper, D.I.; Lloyd, A.; Park, S.-Y.; Tame, J.R.H.
Deposited on : 2005-11-07
Resolution : 1.55 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : trunk30967
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk30967

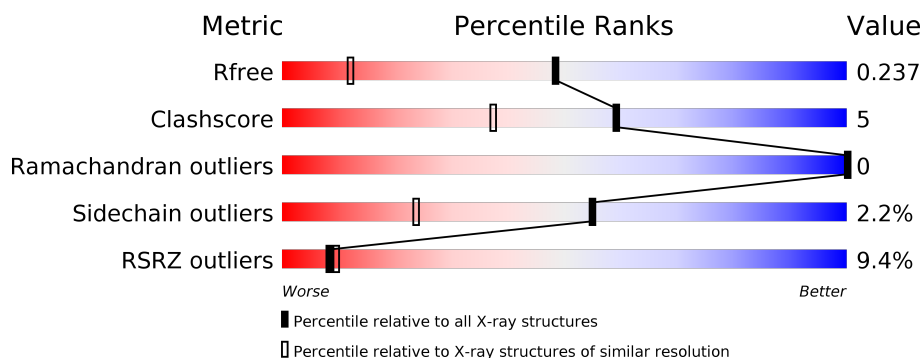
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.55 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}	111664	1224 (1.56-1.56)
Clashscore	122126	1265 (1.56-1.56)
Ramachandran outliers	120053	1240 (1.56-1.56)
Sidechain outliers	120020	1238 (1.56-1.56)
RSRZ outliers	108989	1207 (1.56-1.56)

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	458	<div> <div>9%</div> <div>88%</div> <div>10%</div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3834 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 Penicillin-binding protein 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	456	3489	2202	615	655	17	0	5	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	20	MET	-	INITIATING METHIONINE	UNP P24228
A	261	TYR	ASP	ENGINEERED MUTATION	UNP P24228

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	6	3	3	0	0

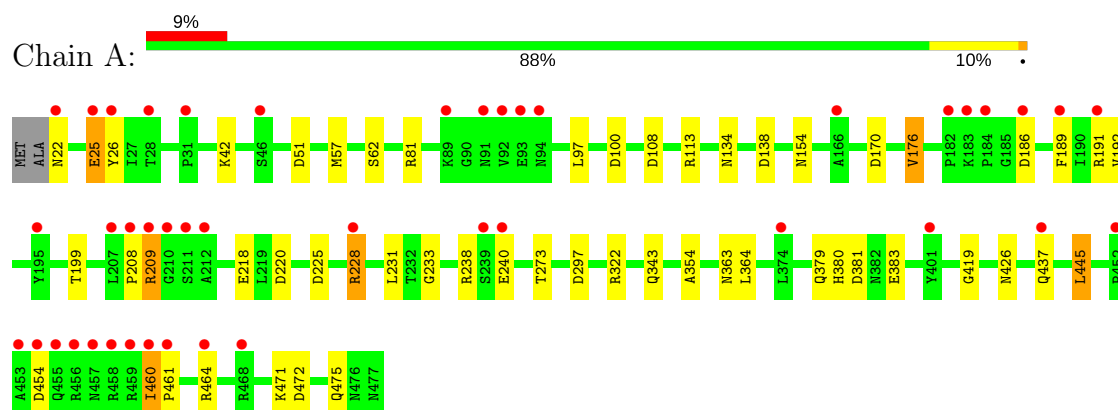
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	339	Total 339	O 339	0	0

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: Penicillin-binding protein 4



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	95.85Å 95.85Å 115.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.55 30.79 – 1.55	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-1.55) 96.0 (30.79-1.55)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 1.55Å)	Xtriage
Refinement program	REFMAC 5.1.19	Depositor
R, R_{free}	0.206 , 0.237 0.206 , 0.237	Depositor DCC
R_{free} test set	3778 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	23.7	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 43.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3834	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.80% 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

Bond lengths and bond angles in the following residue types are not validated in this section: GOL

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	0.57	0/3583	0.81	13/4866 (0.3%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	108	ASP	CB-CG-OD2	7.07	124.66	118.30
1	A	186	ASP	CB-CG-OD2	6.17	123.85	118.30
1	A	472	ASP	CB-CG-OD2	5.91	123.62	118.30
1	A	138	ASP	CB-CG-OD2	5.90	123.61	118.30
1	A	51	ASP	CB-CG-OD2	5.73	123.45	118.30
1	A	100	ASP	CB-CG-OD2	5.72	123.45	118.30
1	A	297	ASP	CB-CG-OD2	5.64	123.38	118.30
1	A	220	ASP	CB-CG-OD2	5.56	123.30	118.30
1	A	170	ASP	CB-CG-OD2	5.47	123.22	118.30
1	A	113	ARG	NE-CZ-NH2	-5.31	117.65	120.30
1	A	381	ASP	CB-CG-OD2	5.25	123.02	118.30
1	A	81	ARG	NE-CZ-NH2	-5.20	117.70	120.30
1	A	454	ASP	CB-CG-OD2	5.03	122.82	118.30

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	3489	0	3483	36	0
2	A	6	0	8	0	0
3	A	339	0	0	3	0
All	All	3834	0	3491	36	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 5.

All (36) 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:189:PHE:HE1	1:A:191:ARG:HG2	1.33	0.93
1:A:189:PHE:CE1	1:A:191:ARG:HG2	2.11	0.86
1:A:189:PHE:HE1	1:A:191:ARG:CG	2.00	0.74
1:A:379:GLN:NE2	3:A:1265:HOH:O	2.21	0.73
1:A:460:ILE:HG13	1:A:461:PRO:HD3	1.70	0.72
1:A:354:ALA:H	1:A:363:ASN:ND2	1.89	0.70
1:A:57:MET:HB3	1:A:364:LEU:HD22	1.79	0.64
1:A:460:ILE:O	1:A:464:ARG:HG2	1.98	0.64
1:A:209:ARG:HE	1:A:209:ARG:H	1.49	0.61
1:A:225:ASP:O	1:A:228:ARG:HD3	2.01	0.61
1:A:134:ASN:HD22	1:A:273:THR:H	1.48	0.60
1:A:22:ASN:HB3	1:A:25:GLU:HG2	1.85	0.59
1:A:189:PHE:CE1	1:A:191:ARG:CG	2.81	0.57
1:A:199:THR:HB	1:A:228:ARG:HG3	1.86	0.57
1:A:42:LYS:HB2	1:A:437:GLN:HE21	1.70	0.56
1:A:42:LYS:HB2	1:A:437:GLN:NE2	2.22	0.55
1:A:471:LYS:NZ	1:A:475:GLN:OE1	2.40	0.53
1:A:209:ARG:NE	1:A:209:ARG:H	2.08	0.52
1:A:154:ASN:H	1:A:154:ASN:ND2	2.09	0.51
1:A:22:ASN:HB2	1:A:26:TYR:CZ	2.45	0.50
1:A:354:ALA:H	1:A:363:ASN:HD21	1.60	0.48
1:A:42:LYS:HE2	3:A:1328:HOH:O	2.13	0.48
1:A:322:ARG:HH12	1:A:343:GLN:HE22	1.62	0.48
1:A:218:GLU:O	1:A:233:GLY:HA3	2.16	0.46
1:A:380:HIS:HE1	3:A:1138:HOH:O	1.99	0.46
1:A:238:ARG:HD3	1:A:240:GLU:O	2.16	0.45
1:A:62:SER:HB2	1:A:419:GLY:HA2	1.99	0.44
1:A:460:ILE:HG13	1:A:461:PRO:CD	2.45	0.44
1:A:176:VAL:HG22	1:A:192:VAL:HB	2.00	0.43
1:A:154:ASN:H	1:A:154:ASN:HD22	1.66	0.43
1:A:25:GLU:HG3	1:A:26:TYR:N	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:209:ARG:HE	1:A:209:ARG:N	2.15	0.43
1:A:380:HIS:HD2	1:A:383:GLU:OE2	2.01	0.43
1:A:426:ASN:HD22	1:A:445:LEU:H	1.67	0.42
1:A:208:PRO:HA	1:A:209:ARG:HH21	1.85	0.41
1:A:134:ASN:ND2	1:A:273:THR:H	2.18	0.41

There are no symmetry-related clashes.

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	458/458 (100%)	446 (97%)	12 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	373/370 (101%)	365 (98%)	8 (2%)	56	25

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	GLU

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Mol	Chain	Res	Type
1	A	97	LEU
1	A	176	VAL
1	A	209	ARG
1	A	228	ARG
1	A	231	LEU
1	A	445	LEU
1	A	460	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	GLN
1	A	114	GLN
1	A	130	GLN
1	A	134	ASN
1	A	154	ASN
1	A	249	GLN
1	A	277	GLN
1	A	343	GLN
1	A	363	ASN
1	A	380	HIS
1	A	382	ASN
1	A	385	ASN
1	A	400	GLN
1	A	422	GLN
1	A	426	ASN
1	A	437	GLN
1	A	457	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

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	GOL	A	501	-	5,5,5	0.40	0	5,5,5	0.81	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	501	-	-	0/4/4/4	0/0/0/0

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

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, 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	456/458 (99%)	0.51	43 (9%) 8 9	18, 27, 50, 69	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	453	ALA	7.8
1	A	209	ARG	7.6
1	A	456	ARG	7.4
1	A	455	GLN	6.4
1	A	210	GLY	6.3
1	A	212	ALA	5.7
1	A	184	PRO	5.5
1	A	46	SER	5.4
1	A	26	TYR	5.0
1	A	182	PRO	5.0
1	A	208	PRO	4.9
1	A	401[A]	TYR	4.8
1	A	460	ILE	4.8
1	A	452	PRO	4.7
1	A	189	PHE	4.7
1	A	457	ASN	4.7
1	A	183	LYS	4.6
1	A	239	SER	4.0
1	A	22	ASN	3.8
1	A	454	ASP	3.6
1	A	25	GLU	3.6
1	A	89	LYS	3.2
1	A	240	GLU	3.2
1	A	437	GLN	3.1
1	A	464	ARG	3.1
1	A	92	VAL	2.9
1	A	207	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	28	THR	2.8
1	A	211	SER	2.7
1	A	94	ASN	2.5
1	A	191	ARG	2.5
1	A	93	GLU	2.5
1	A	458	ARG	2.5
1	A	186	ASP	2.5
1	A	468	ARG	2.5
1	A	461	PRO	2.4
1	A	31	PRO	2.4
1	A	166	ALA	2.3
1	A	374	LEU	2.1
1	A	195	TYR	2.1
1	A	91	ASN	2.1
1	A	459	ARG	2.1
1	A	228	ARG	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 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. 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	B-factors(\AA^2)	Q<0.9
2	GOL	A	501	6/6	0.77	0.16	33,40,44,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.