



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:46 PM GMT

PDB ID : 1QMF
Title : PENICILLIN-BINDING PROTEIN 2X (PBP-2X) ACYL-ENZYME COMPLEX
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Deposited on : 1999-09-28
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
<http://wwpdb.org/validation/2016/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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

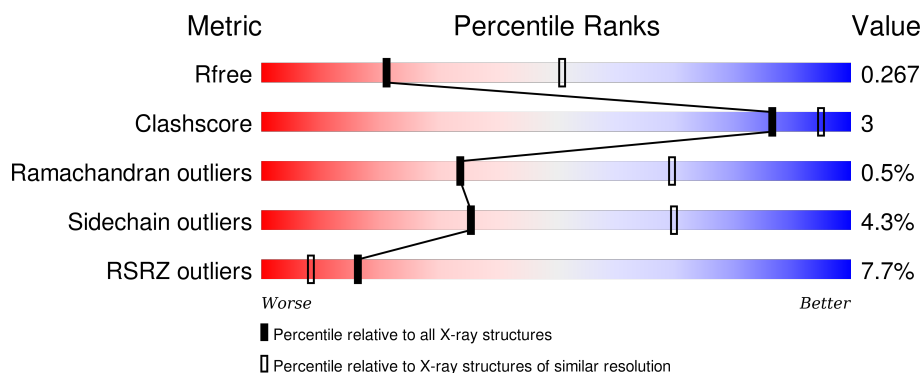
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

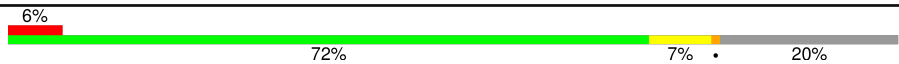
The reported resolution of this entry is 2.80 Å.

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}	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (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 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	702	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	KEF	A	900	X	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4373 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.

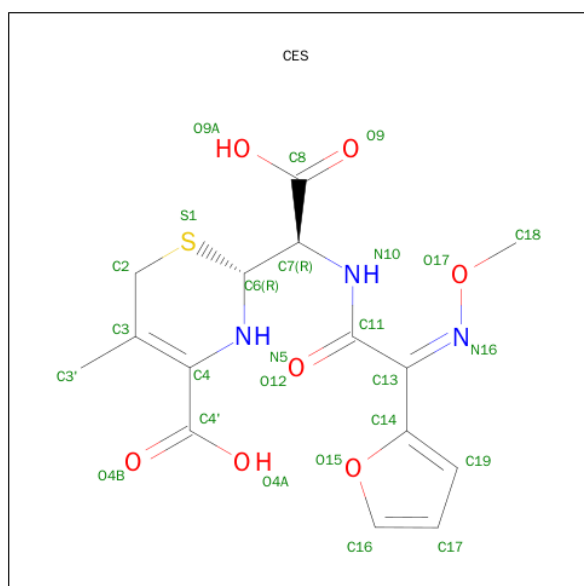
- Molecule 1 is a protein called PENICILLIN-BINDING PROTEIN 2X.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	559	Total	C	N	O	S	0	0	0
			4262	2669	712	861	20			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	385	MET	THR	CONFLICT	UNP P14677

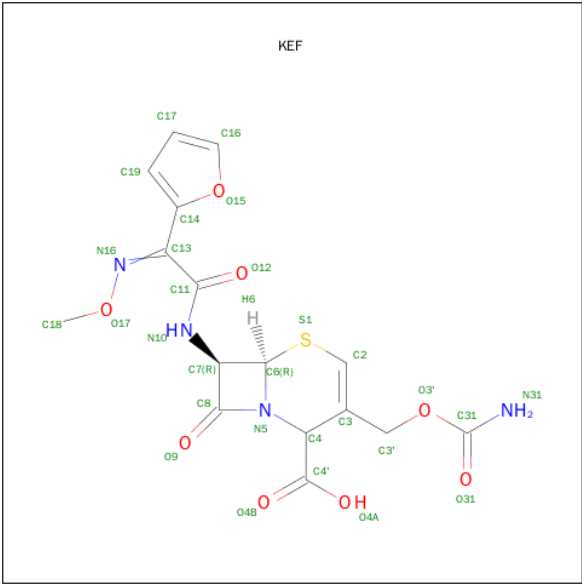
- Molecule 2 is 2-[CARBOXY-(2-FURAN-2-YL-2-METHOXYIMINO-ACETYLAMINO)-METHYL]-5-METHYL-3,6-DIHYDRO-2H-[1,3]THIAZINE-4-CARBOXYLIC ACID (three-letter code: CES) (formula: C₁₅H₁₇N₃O₇S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			25	15	3	6	1		

- Molecule 3 is CEFUROXIME (OCT-3-ENE FORM) (three-letter code: KEF) (formula:

C₁₆H₁₆N₄O₈S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			29	16	4	8	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	57	Total	O	0	0
			57	57		

4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	129.91Å 129.91Å 139.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.80 39.42 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.1 (50.00-2.80) 97.1 (39.42-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 2.81Å)	Xtriage
Refinement program	CNS 0.5	Depositor
R, R_{free}	0.239 , 0.271 0.236 , 0.267	Depositor DCC
R_{free} test set	2910 reflections (9.95%)	DCC
Wilson B-factor (Å ²)	53.9	Xtriage
Anisotropy	0.444	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 29276 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4373	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.92% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: KEF, CES

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	0.62	0/4336	0.77	0/5877

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

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	4262	0	4190	23	0
2	A	25	0	15	1	0
3	A	29	0	14	0	0
4	A	57	0	0	2	0
All	All	4373	0	4219	23	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 3.

All (23) 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:281:GLN:HE22	1:A:286:GLY:N	1.90	0.69
1:A:281:GLN:HE22	1:A:286:GLY:H	1.42	0.66
1:A:281:GLN:NE2	1:A:286:GLY:H	1.97	0.63
1:A:76:ARG:HH21	1:A:220:GLU:HG3	1.62	0.62
1:A:492:ARG:HG3	1:A:634:SER:HB2	1.81	0.62
1:A:348:ILE:HD13	1:A:510:LEU:HB3	1.81	0.61
1:A:545:ALA:HB3	1:A:575:MET:HB2	1.82	0.60
1:A:374:TRP:HB3	2:A:800:CES:H22	1.84	0.60
1:A:306:GLN:O	1:A:309:THR:HG23	2.05	0.56
1:A:536:THR:HG22	1:A:537:VAL:HG23	1.89	0.53
1:A:570:PHE:HB2	1:A:589:VAL:HG22	1.90	0.53
1:A:288:TYR:HE1	1:A:323:VAL:HG21	1.76	0.50
1:A:183:SER:N	4:A:2004:HOH:O	2.45	0.50
1:A:329:TYR:HB2	1:A:432:GLU:HG3	1.93	0.50
1:A:467:ALA:HB2	1:A:474:MET:HB3	1.93	0.49
1:A:498:ILE:HG21	1:A:700:TYR:HD2	1.78	0.48
1:A:468:ILE:HG21	1:A:515:MET:HE1	1.98	0.45
1:A:512:ARG:HD2	1:A:578:ALA:O	2.17	0.45
1:A:739:LYS:HD2	1:A:739:LYS:H	1.82	0.45
1:A:289:MET:HB2	1:A:307:ARG:HB2	1.98	0.44
1:A:340:LYS:HG2	1:A:400:MET:HG2	2.00	0.43
1:A:376:VAL:HG11	4:A:2016:HOH:O	2.20	0.41
1:A:79:ILE:HD11	1:A:227:LEU:HD21	2.03	0.41

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	551/702 (78%)	523 (95%)	25 (4%)	3 (0%)	34 69

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	556	GLU
1	A	554	ALA
1	A	700	TYR

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	468/590 (79%)	448 (96%)	20 (4%)	35 70

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	213	LEU
1	A	227	LEU
1	A	282	GLU
1	A	293	LEU
1	A	309	THR
1	A	323	VAL
1	A	339	MET
1	A	364	LEU
1	A	368	ASP
1	A	427	PHE
1	A	510	LEU
1	A	532	THR
1	A	536	THR
1	A	552	GLN
1	A	553	ILE
1	A	589	VAL
1	A	599	GLN
1	A	615	LYS
1	A	671	ASN
1	A	689	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	281	GLN
1	A	458	GLN
1	A	558	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 ⓘ

2 ligands are 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	CES	A	800	1	17,26,27	4.67	8 (47%)	16,35,37	2.60	8 (50%)
3	KEF	A	900	-	21,31,31	2.73	6 (28%)	24,44,44	2.51	9 (37%)

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	CES	A	800	1	-	1/10/38/40	0/0/2/2
3	KEF	A	900	-	1/1/10/12	0/15/53/53	0/1/3/3

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	800	CES	C3'-C3	-9.32	1.35	1.51
2	A	800	CES	C2-C3	-7.69	1.41	1.51
3	A	900	KEF	C4-N5	-3.08	1.42	1.47
3	A	900	KEF	O31-C31	-2.74	1.17	1.21
2	A	800	CES	C13-C11	-2.23	1.46	1.50
2	A	800	CES	C17-C19	-2.06	1.33	1.39
3	A	900	KEF	O17-N16	3.55	1.48	1.40
3	A	900	KEF	C6-C7	3.66	1.63	1.56
3	A	900	KEF	O3'-C31	3.80	1.41	1.35
2	A	800	CES	C13-N16	3.88	1.35	1.29
2	A	800	CES	C4'-C4	4.39	1.59	1.52
2	A	800	CES	C7-N10	4.56	1.51	1.46
3	A	900	KEF	C31-N31	9.07	1.53	1.33
2	A	800	CES	C3-C4	12.29	1.50	1.34

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	800	CES	C19-C17-C16	-4.29	96.52	112.78
3	A	900	KEF	C19-C17-C16	-4.13	97.11	112.78
3	A	900	KEF	O3'-C31-O31	-3.35	119.94	123.28
2	A	800	CES	C2-C3-C4	-2.68	118.99	123.50
2	A	800	CES	C3'-C3-C4	-2.63	120.11	123.48
3	A	900	KEF	O3'-C3'-C3	-2.56	103.26	109.34
2	A	800	CES	C14-C13-N16	-2.53	109.55	122.17
3	A	900	KEF	C7-C6-N5	-2.50	84.76	87.98
2	A	800	CES	C18-O17-N16	-2.22	105.94	108.47
3	A	900	KEF	O31-C31-N31	2.13	129.43	125.34
3	A	900	KEF	C6-N5-C8	2.68	97.39	93.89
3	A	900	KEF	C7-C6-S1	2.74	121.06	116.45
2	A	800	CES	O12-C11-C13	2.84	123.69	120.53
3	A	900	KEF	O17-N16-C13	4.25	116.18	111.11
2	A	800	CES	O17-N16-C13	4.48	116.45	111.11
2	A	800	CES	C3'-C3-C2	5.01	120.89	113.22
3	A	900	KEF	C3-C4-N5	7.84	119.01	109.08

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	900	KEF	C4

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	800	CES	O12-C11-C13-C14

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	800	CES	1	0

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	559/702 (79%)	0.29	43 (7%) 16 8	28, 49, 97, 100	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	558	ASN	12.6
1	A	559	GLY	10.5
1	A	380	LEU	7.9
1	A	381	THR	6.4
1	A	560	GLY	6.2
1	A	377	ASN	6.0
1	A	557	LYS	5.8
1	A	234	ILE	5.7
1	A	232	GLY	4.9
1	A	620	LEU	4.8
1	A	376	VAL	4.6
1	A	561	TYR	4.5
1	A	233	ILE	4.5
1	A	378	GLU	4.2
1	A	563	VAL	4.2
1	A	71	THR	4.1
1	A	382	GLY	3.8
1	A	556	GLU	3.8
1	A	552	GLN	3.7
1	A	254	ARG	3.6
1	A	554	ALA	3.5
1	A	379	GLY	3.5
1	A	255	THR	3.3
1	A	383	GLY	3.3
1	A	368	ASP	3.2
1	A	208	ASP	3.2
1	A	72	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	374	TRP	3.2
1	A	210	SER	2.8
1	A	439	ALA	2.8
1	A	211	LYS	2.8
1	A	183	SER	2.6
1	A	555	ASP	2.6
1	A	740	ASP	2.5
1	A	92	THR	2.4
1	A	442	ILE	2.4
1	A	565	LEU	2.3
1	A	375	ASP	2.3
1	A	739	LYS	2.3
1	A	562	LEU	2.2
1	A	720	PHE	2.2
1	A	459	THR	2.2
1	A	73	PRO	2.1

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
3	KEF	A	900	29/29	0.85	0.48	7.54	28,43,58,63	29
2	CES	A	800	25/26	0.88	0.24	0.06	42,67,87,92	0

6.5 Other polymers [i](#)

There are no such residues in this entry.