



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:10 AM UTC

PDB ID : 3DRJ / pdb_00003drj
Title : Crystal structure of Lactococcal OppA co-crystallized with pTH-related peptide in an open conformation
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Deposited on : 2008-07-11
Resolution : 1.50 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

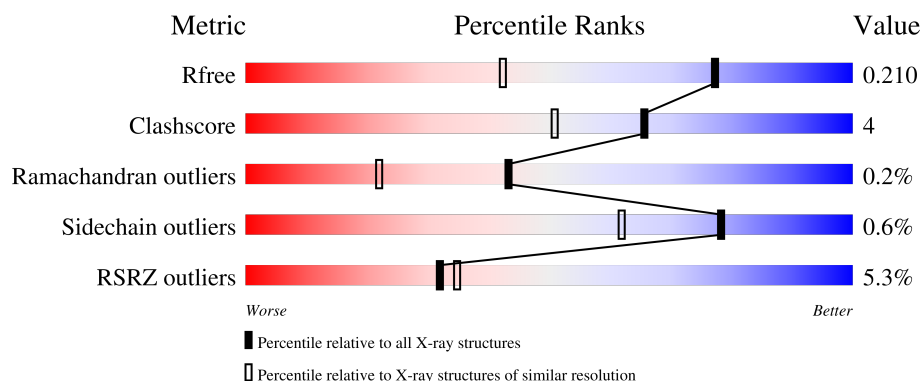
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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	590	 4% 89% 5% • 5%
2	B	5	 80% 60% 40%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5338 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 Oligopeptide-binding protein oppA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	560	Total	C	N	O	S	0	6	0
			4403	2791	727	871	14			

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP A2RJ53
A	579	GLY	-	expression tag	UNP A2RJ53
A	580	SER	-	expression tag	UNP A2RJ53
A	581	ILE	-	expression tag	UNP A2RJ53
A	582	GLU	-	expression tag	UNP A2RJ53
A	583	GLY	-	expression tag	UNP A2RJ53
A	584	ARG	-	expression tag	UNP A2RJ53
A	585	HIS	-	expression tag	UNP A2RJ53
A	586	HIS	-	expression tag	UNP A2RJ53
A	587	HIS	-	expression tag	UNP A2RJ53
A	588	HIS	-	expression tag	UNP A2RJ53
A	589	HIS	-	expression tag	UNP A2RJ53
A	590	HIS	-	expression tag	UNP A2RJ53

- Molecule 2 is a protein called pTH-related peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	5	Total	C	N	O	0	0	0
			34	21	8	5			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	897	Total	O	0	0
			897	897		

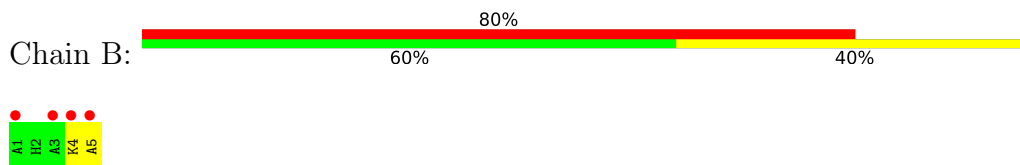
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	4	Total	O	0	0
			4	4		

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- Molecule 1: Oligopeptide-binding protein oppA



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	40.14Å 123.36Å 59.70Å 90.00° 103.91° 90.00°	Depositor
Resolution (Å)	37.16 – 1.50 37.16 – 1.50	Depositor EDS
% Data completeness (in resolution range)	95.8 (37.16-1.50) 95.8 (37.16-1.50)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.07 (at 1.50Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.162 , 0.208 0.171 , 0.210	Depositor DCC
R_{free} test set	4288 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	12.8	Xtriage
Anisotropy	0.105	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5338	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.42% 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	0.84	3/4504 (0.1%)	0.89	1/6103 (0.0%)
2	B	0.87	0/34	0.85	0/44
All	All	0.84	3/4538 (0.1%)	0.89	1/6147 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	361	ILE	CA-CB	-8.63	1.47	1.54
1	A	364	ILE	CA-CB	5.55	1.62	1.54
1	A	369	THR	CA-CB	5.12	1.62	1.53

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	283	TYR	N-CA-C	5.76	117.56	111.28

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	4403	0	4294	31	0
2	B	34	0	37	3	0
3	A	897	0	0	12	0
3	B	4	0	0	0	0
All	All	5338	0	4331	33	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 (33) 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:341:ASN:HD21	1:A:428[B]:ASN:HD21	1.09	0.92
1:A:396:ASN:HD22	1:A:397:LYS:N	1.67	0.92
1:A:503:LYS:HE2	3:A:1457:HOH:O	1.70	0.91
1:A:503:LYS:CE	3:A:1457:HOH:O	2.24	0.84
1:A:341:ASN:HD21	1:A:428[B]:ASN:ND2	1.78	0.81
1:A:341:ASN:ND2	1:A:428[B]:ASN:HD21	1.82	0.78
2:B:4:LYS:O	2:B:5:ALA:HB2	1.90	0.69
1:A:396:ASN:HD22	1:A:396:ASN:C	1.99	0.67
1:A:281:SER:C	3:A:1477:HOH:O	2.44	0.59
1:A:278:MET:HB2	3:A:660:HOH:O	2.04	0.57
1:A:341:ASN:OD1	1:A:428[B]:ASN:OD1	2.24	0.55
1:A:174:GLN:CG	3:A:883:HOH:O	2.55	0.54
1:A:423:GLU:HB2	3:A:838:HOH:O	2.07	0.54
1:A:280:SER:HA	1:A:353:LEU:HD13	1.90	0.54
1:A:260:THR:HG22	3:A:788:HOH:O	2.07	0.54
1:A:299:ALA:HB3	1:A:542[B]:MET:HG2	1.90	0.53
1:A:260:THR:CG2	3:A:788:HOH:O	2.57	0.52
1:A:95:SER:O	1:A:174:GLN:NE2	2.40	0.51
1:A:300:MET:SD	1:A:364:ILE:HD13	2.51	0.50
1:A:363:PRO:O	1:A:366:LYS:HD3	2.12	0.50
1:A:355[A]:THR:HG23	3:A:817:HOH:O	2.12	0.48
1:A:474:SER:OG	2:B:5:ALA:HB3	2.13	0.48
1:A:396:ASN:C	1:A:396:ASN:ND2	2.66	0.47
1:A:234:SER:HB3	1:A:256:GLU:HG2	1.96	0.47
2:B:4:LYS:O	2:B:5:ALA:CB	2.55	0.47
1:A:282:GLN:HA	3:A:1477:HOH:O	2.13	0.46
1:A:299:ALA:HB3	1:A:542[B]:MET:CG	2.45	0.46
1:A:282:GLN:N	3:A:1477:HOH:O	2.48	0.45
1:A:174:GLN:CD	3:A:883:HOH:O	2.63	0.41
1:A:396:ASN:HD22	1:A:397:LYS:H	1.57	0.41
1:A:45:MET:O	1:A:232:GLY:HA2	2.21	0.40
1:A:304:LEU:HD12	1:A:536:VAL:HB	2.03	0.40
1:A:25:ASN:OD1	1:A:26:PRO:HD2	2.21	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	564/590 (96%)	551 (98%)	12 (2%)	1 (0%)	43	22
2	B	3/5 (60%)	3 (100%)	0	0	100	100
All	All	567/595 (95%)	554 (98%)	12 (2%)	1 (0%)	43	22

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	282	GLN

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	476/496 (96%)	473 (99%)	3 (1%)	78	62
2	B	2/2 (100%)	2 (100%)	0	100	100
All	All	478/498 (96%)	475 (99%)	3 (1%)	78	62

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	SER
1	A	280	SER
1	A	396	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	GLN
1	A	69	GLN
1	A	167	ASN
1	A	285	GLN
1	A	396	ASN
1	A	421	ASN
1	A	431	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	560/590 (94%)	0.18	26 (4%) 37 41	4, 11, 25, 34	6 (1%)
2	B	5/5 (100%)	2.34	4 (80%) 0 0	28, 28, 30, 33	0
All	All	565/595 (94%)	0.20	30 (5%) 32 35	4, 11, 26, 34	6 (1%)

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	398	SER	5.2
1	A	420	ALA	4.5
2	B	5	ALA	3.5
1	A	279	VAL	3.5
1	A	288	ASN	3.5
1	A	277	GLY	3.2
1	A	280	SER	2.9
1	A	14	LEU	2.8
1	A	283	TYR	2.8
1	A	26	PRO	2.7
1	A	395	LEU	2.5
1	A	286	VAL	2.5
1	A	418	GLY	2.4
1	A	105	ASP	2.4
1	A	364	ILE	2.3
1	A	419	ASP	2.3
1	A	281	SER	2.3
1	A	167	ASN	2.3
1	A	285	GLN	2.3
1	A	290	LYS	2.3
1	A	34	ASN	2.3
1	A	399	THR	2.2
1	A	397	LYS	2.2
2	B	4	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	284	LYS	2.2
1	A	260	THR	2.1
2	B	3	ALA	2.1
1	A	24	GLN	2.1
2	B	1	ALA	2.1
1	A	390	GLU	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 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.