

wwPDB X-ray Structure Validation Summary Report (i)

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PDB ID 3G9I

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2009-02-13 Deposited on

1.85 Å(reported) Resolution

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 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

> 1.8.5 (274361), CSD as541be (2020) Mogul

Xtriage (Phenix) 1.13 EDS 2.35

20191225.v01 (using entries in the PDB archive December 25th 2019) Percentile statistics

> Refmac 5.8.0158

CCP4 7.0.044 (Gargrove)

Ideal geometry (proteins) Engh & Huber (2001) Ideal geometry (DNA, RNA) Parkinson et al. (1996)

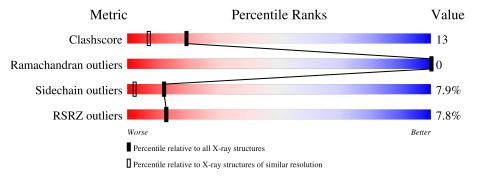
Validation Pipeline (wwPDB-VP) 2.35

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 1.85 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	Similar resolution
	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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	90	11%	66%		17%	6%	12%	
1	В	90	6%	69%		19%		10%	
2	D	16	50%	, D		50%			
3	С	16	25%		56%	_	19	%	



2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 2042 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 Glucocorticoid receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	79	Total 601	_		O 105	S 12	0	0	0
1	В	81	Total 611	_	N 120	_	S 12	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	436	GLY	-	expression tag	UNP P06536
A	437	SER	-	expression tag	UNP P06536
A	438	HIS	-	expression tag	UNP P06536
A	439	MET	-	expression tag	UNP P06536
В	436	GLY	-	expression tag	UNP P06536
В	437	SER	-	expression tag	UNP P06536
В	438	HIS	-	expression tag	UNP P06536
В	439	MET	-	expression tag	UNP P06536

• Molecule 2 is a DNA chain called DNA (5'-D(*TP*AP*GP*AP*AP*AP*AP*AP*AP*AP*TP*GP*TP*TP*CP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	16	Total 326	C 158	N 61	O 92	P 15	0	0	0

• Molecule 3 is a DNA chain called DNA (5'-D(*AP*AP*GP*AP*AP*CP*AP*TP*TP*TP* TP*GP*TP*TP*CP*T)-3').

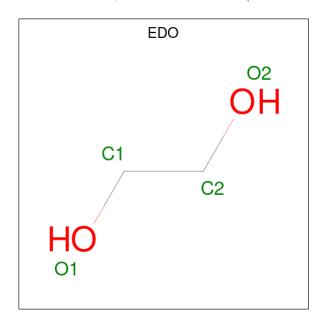
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	C	16	Total	С	N	О	Р	0	0	0
3		16	324	158	55	96	15	0	U	

• Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	2	Total Zn 2 2	0	0
4	В	2	Total Zn 2 2	0	0

 \bullet Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $\mathrm{C_2H_6O_2}).$



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	В	1	Total C O 4 2 2	0	0

• Molecule 6 is water.

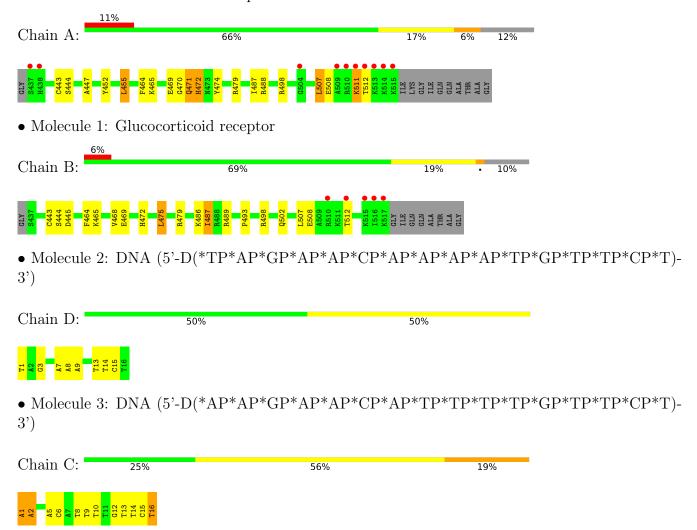
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	50	Total O 50 50	0	0
6	В	69	Total O 69 69	0	0
6	D	22	Total O 22 22	0	0
6	C	23	Total O 23 23	0	0



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.







4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	116.51Å 37.65Å 95.45Å	Donositor
a, b, c, α , β , γ	90.00° 123.08° 90.00°	Depositor
Resolution (Å)	48.81 - 1.85	Depositor
rtesolution (A)	48.81 - 1.85	EDS
% Data completeness	80.7 (48.81-1.85)	Depositor
(in resolution range)	80.8 (48.81-1.85)	EDS
R_{merge}	0.06	Depositor
R_{sym}	0.05	Depositor
$< I/\sigma(I) > 1$	1.16 (at 1.86Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
D D.	0.170 , 0.201	Depositor
R, R_{free}	0.174 , (Not available)	DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	28.6	Xtriage
Anisotropy	0.673	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.34, 62.5	EDS
L-test for twinning ²	$ < L >=0.48, < L^2>=0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	2042	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 15.64% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities estimated from amplitudes.

5 Model quality (i)

5.1 Standard geometry (i)

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

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	Boı	nd lengths	Bond angles		
IVIOI		RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.62	0/609	0.66	0/813	
1	В	0.72	0/619	0.78	1/827 (0.1%)	
2	D	1.20	1/366 (0.3%)	2.27	$14/563 \ (2.5\%)$	
3	С	1.37	2/362~(0.6%)	2.30	19/557 (3.4%)	
All	All	0.95	3/1956~(0.2%)	1.56	$34/2760 \ (1.2\%)$	

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(\mathbf{\mathring{A}})$	$Ideal(\AA)$
3	С	12	DG	C6-N1	5.93	1.43	1.39
2	D	3	DG	C3'-O3'	-5.81	1.36	1.44
3	С	1	DA	N9-C4	-5.43	1.34	1.37

The worst 5 of 34 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
2	D	9	DA	O4'-C1'-N9	-13.30	98.69	108.00
2	D	9	DA	N1-C6-N6	11.36	125.42	118.60
2	D	13	DT	O4'-C1'-N1	-10.87	100.39	108.00
3	С	14	DT	O4'-C1'-N1	-10.49	100.66	108.00
3	С	1	DA	C2-N3-C4	-10.06	105.57	110.60

There are no chirality outliers.

There are no planarity outliers.

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	601	0	580	16	0
1	В	611	0	584	14	0
2	D	326	0	183	1	0
3	С	324	0	185	13	0
4	A	2	0	0	0	0
4	В	2	0	0	0	0
5	A	8	0	12	3	1
5	В	4	0	6	0	1
6	A	50	0	0	2	0
6	В	69	0	0	2	0
6	С	23	0	0	3	0
6	D	22	0	0	0	0
All	All	2042	0	1550	44	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 13.

The worst 5 of 44 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} \text{Interatomic} \\ \text{distance (Å)} \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\r{A}}) \end{array}$
3:C:16:DT:H5"	3:C:16:DT:H6	1.21	1.04
3:C:16:DT:H5"	3:C:16:DT:C6	1.94	1.03
3:C:16:DT:H6	3:C:16:DT:C5'	1.80	0.94
3:C:1:DA:H8	3:C:1:DA:H5"	1.38	0.88
1:B:465:LYS:HE2	1:B:469:GLU:CD	2.01	0.81

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	$egin{aligned} ext{Interatomic} \ ext{distance} & (ext{Å}) \end{aligned}$	$egin{array}{c} \operatorname{Clash} \ \operatorname{overlap}\ (ext{\AA}) \end{array}$
5:A:3:EDO:O1	5:B:1:EDO:O2[4_445]	1.86	0.34



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	Perce	\mathbf{ntiles}
1	A	77/90 (86%)	72 (94%)	5 (6%)	0	100	100
1	В	79/90 (88%)	75 (95%)	4 (5%)	0	100	100
All	All	156/180 (87%)	147 (94%)	9 (6%)	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	63/73~(86%)	57 (90%)	6 (10%)	8 1
1	В	63/73 (86%)	59 (94%)	4 (6%)	18 4
All	All	126/146~(86%)	116 (92%)	10 (8%)	12 2

5 of 10 residues with a non-rotameric side chain are listed below:

Mol	Chain	Res	Type
1	В	487	ILE
1	В	508	GLU
1	В	512	THR
1	A	487	ILE
1	A	507	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:



Mol	Chain	Res	Type
1	A	438	HIS
1	A	471	GLN
1	A	472	HIS
1	В	438	HIS
1	В	480	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)

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 for Mogul analysis.

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	B	ond leng	$_{ m gths}$	Е	ond ang	gles
WIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	A	2	-	3,3,3	0.51	0	2,2,2	0.32	0
5	EDO	A	3	-	3,3,3	0.53	0	2,2,2	0.48	0
5	EDO	В	1	-	3,3,3	0.62	0	2,2,2	0.13	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
5	EDO	A	2	-	-	0/1/1/1	-
5	EDO	A	3	-	-	1/1/1/1	-
5	EDO	В	1	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	3	EDO	O1-C1-C2-O2
5	В	1	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	2	EDO	1	0
5	A	3	EDO	2	1
5	В	1	EDO	0	1

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, 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>	$\#\mathrm{RSRZ}{>}2$	$\mathbf{OWAB}(\mathrm{\AA}^2)$	Q < 0.9
1	A	79/90 (87%)	0.40	10 (12%) 3 3	24, 51, 121, 160	0
1	В	81/90 (90%)	0.28	5 (6%) 20 20	23, 36, 89, 136	0
2	D	16/16 (100%)	-0.42	0 100 100	39, 53, 87, 90	0
3	С	16/16 (100%)	-0.28	0 100 100	33, 61, 82, 103	0
All	All	192/212 (90%)	0.22	15 (7%) 13 13	23, 45, 106, 160	0

The worst 5 of 15 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	515	LYS	4.7
1	A	438	HIS	4.4
1	В	516	ILE	3.9
1	A	437	SER	3.8
1	A	511	LYS	3.3

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)

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, 95^{th} 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
5	EDO	A	3	4/4	0.43	0.39	72,72,78,90	0
5	EDO	В	1	4/4	0.65	0.35	78,81,84,87	0
5	EDO	A	2	4/4	0.86	0.11	57,62,64,77	0
4	ZN	В	527	1/1	1.00	0.14	26,26,26,26	0
4	ZN	A	526	1/1	1.00	0.11	46,46,46,46	0
4	ZN	A	527	1/1	1.00	0.13	28,28,28,28	0
4	ZN	В	526	1/1	1.00	0.12	31,31,31,31	0

6.5 Other polymers (i)

There are no such residues in this entry.

