



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 07:39 PM GMT

PDB ID : 1GM5
Title : STRUCTURE OF RECG BOUND TO THREE-WAY DNA JUNCTION
Authors : Singleton, M.R.; Scaife, S.; Wigley, D.B.
Deposited on : 2001-09-11
Resolution : 3.24 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
<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

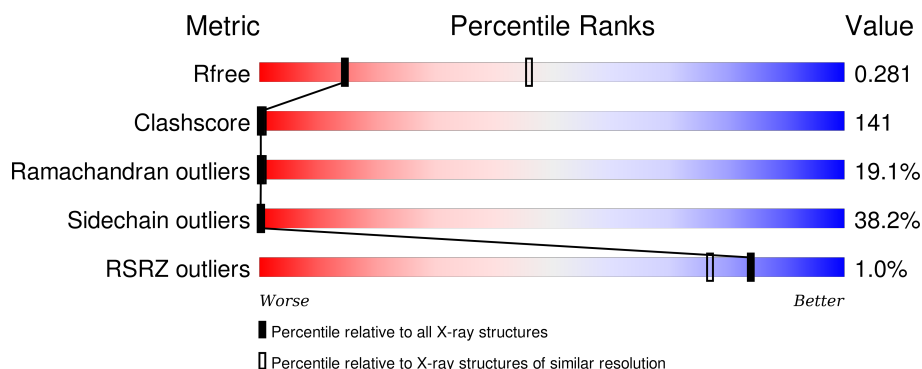
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.24 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	1092 (3.28-3.20)
Clashscore	102246	1227 (3.28-3.20)
Ramachandran outliers	100387	1204 (3.28-3.20)
Sidechain outliers	100360	1203 (3.28-3.20)
RSRZ outliers	91569	1097 (3.28-3.20)

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	780	<div> <div>8%</div> <div>44%</div> <div>32%</div> <div>9%</div> <div>7%</div> </div>
2	X	20	<div> <div>5%</div> <div>5%</div> <div>30%</div> <div>25%</div> <div>40%</div> </div>
3	Y	20	<div> <div>25%</div> <div>75%</div> </div>
4	Z	9	<div> <div>22%</div> <div>78%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6779 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 RECG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	729	Total	C	N	O	S	0	0	0
			5919	3794	1017	1083	25			

- Molecule 2 is a DNA chain called DNA (5'-(*CP*AP*GP*CP*TP*CP*CP*AP*TP*GP*AP*TP* CP*AP*TP*TP*GP*GP*CP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	X	12	Total	C	N	O	P	0	0	0
			239	116	43	69	11			

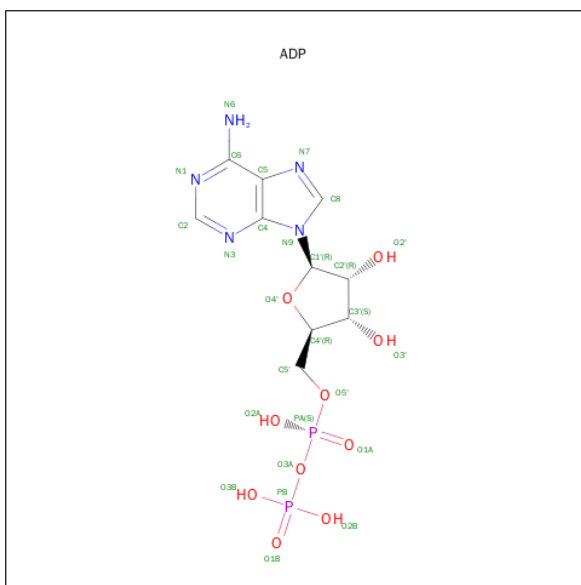
- Molecule 3 is a DNA chain called DNA (5'-(*GP*CP*AP*GP*TP*GP*CP*TP*CP*GP*CP*AP* TP*GP*GP*AP*GP*CP*TP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	Y	20	Total	C	N	O	P	0	0	0
			411	195	78	119	19			

- Molecule 4 is a DNA chain called DNA (5'-(*GP*AP*GP*CP*AP*CP*TP*GP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	Z	9	Total	C	N	O	P	0	0	0
			182	87	36	51	8			

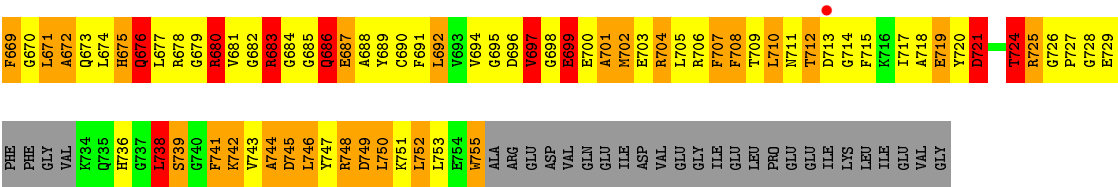
- Molecule 5 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



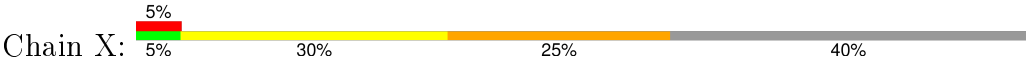
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

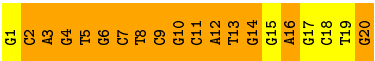
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Mg	0	0
			1	1		



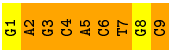
• Molecule 2: DNA (5'-(*CP*AP*GP*CP*TP*CP*CP*AP*TP*GP*AP*TP* CP*AP*TP*TP* GP*GP*CP*A)-3')



• Molecule 3: DNA (5'-(*GP*CP*AP*GP*TP*GP*CP*TP*CP*GP*CP*AP* TP*GP*GP*AP* GP*CP*TP*G)-3')



• Molecule 4: DNA (5'-(*GP*AP*GP*CP*AP*CP*TP*GP*C)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	133.70 Å 144.60 Å 84.02 Å 90.00° 113.82° 90.00°	Depositor
Resolution (Å)	20.00 – 3.24 19.59 – 3.25	Depositor EDS
% Data completeness (in resolution range)	99.3 (20.00-3.24) 98.2 (19.59-3.25)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.53 (at 3.22 Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.272 , 0.328 0.273 , 0.281	Depositor DCC
R_{free} test set	1153 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	96.2	Xtriage
Anisotropy	0.209	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 94.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 44482 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6779	wwPDB-VP
Average B, all atoms (Å ²)	94.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% 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: MG, ADP

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	0.96	8/6025 (0.1%)	1.26	52/8101 (0.6%)
2	X	1.19	0/267	1.99	11/410 (2.7%)
3	Y	1.67	5/461 (1.1%)	3.13	66/711 (9.3%)
4	Z	1.81	4/204 (2.0%)	4.22	58/313 (18.5%)
All	All	1.07	17/6957 (0.2%)	1.68	187/9535 (2.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	10

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	55	GLU	CD-OE1	8.25	1.34	1.25
1	A	553	GLU	CD-OE2	6.75	1.33	1.25
3	Y	10	DG	C2-N3	6.19	1.37	1.32
3	Y	9	DC	N1-C6	-6.14	1.33	1.37
1	A	55	GLU	CG-CD	6.09	1.61	1.51

The worst 5 of 187 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Y	5	DT	N3-C4-O4	16.48	129.79	119.90
4	Z	2	DA	O4'-C1'-N9	-15.33	97.27	108.00
4	Z	7	DT	O4'-C1'-N1	14.47	118.13	108.00
3	Y	8	DT	N3-C4-O4	14.32	128.49	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Y	8	DT	O4'-C1'-N1	14.00	117.80	108.00

There are no chirality outliers.

5 of 10 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	444	PHE	Peptide
1	A	449	ALA	Peptide
1	A	471	GLN	Peptide
1	A	546	GLU	Peptide
1	A	607	SER	Peptide

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	5919	0	6036	1776	23
2	X	239	0	133	30	0
3	Y	411	0	226	59	5
4	Z	182	0	102	18	0
5	A	27	0	12	0	0
6	A	1	0	0	0	0
All	All	6779	0	6509	1868	23

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 141.

The worst 5 of 1868 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:7:PHE:CD2	1:A:11:LEU:HD11	1.14	1.63
1:A:621:GLY:HA3	1:A:647:THR:CG2	1.31	1.61
1:A:615:LYS:HD3	1:A:641:ASP:CB	1.15	1.59
1:A:615:LYS:CD	1:A:641:ASP:HB2	1.29	1.58
1:A:7:PHE:HD2	1:A:11:LEU:CD1	1.14	1.57

The worst 5 of 23 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	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:ARG:CZ	1:A:40:ARG:NH1[2_557]	0.36	1.84
1:A:12:PHE:CE1	1:A:563:ARG:NH2[2_556]	0.56	1.64
1:A:40:ARG:NE	1:A:40:ARG:NH1[2_557]	1.02	1.18
1:A:40:ARG:CZ	1:A:40:ARG:CZ[2_557]	1.13	1.07
1:A:12:PHE:CE1	1:A:563:ARG:CZ[2_556]	1.28	0.92

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	719/780 (92%)	405 (56%)	177 (25%)	137 (19%)	0 0

5 of 137 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	8	THR
1	A	10	SER
1	A	27	ASN
1	A	34	LYS
1	A	37	VAL

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	646/692 (93%)	399 (62%)	247 (38%)	0 0

5 of 247 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	323	ARG
1	A	413	ASN
1	A	706	ARG
1	A	334	LEU
1	A	363	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 26 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	309	HIS
1	A	377	HIS
1	A	686	GLN
1	A	346	HIS
1	A	373	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	ADP	A	1756	6	22,29,29	1.23	1 (4%)	27,45,45	2.47	3 (11%)

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	ADP	A	1756	6	-	0/12/32/32	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1756	ADP	C2-N3	2.94	1.37	1.32

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1756	ADP	N3-C2-N1	-11.03	120.45	128.89
5	A	1756	ADP	O4'-C4'-C3'	-2.16	100.79	105.15
5	A	1756	ADP	O3A-PA-O5'	2.74	110.20	102.94

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 [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, 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	729/780 (93%)	-0.31	7 (0%) 84 77	34, 89, 140, 175	0
2	X	12/20 (60%)	-0.13	1 (8%) 14 9	83, 103, 145, 160	0
3	Y	20/20 (100%)	-0.28	0 100 100	44, 101, 153, 165	0
4	Z	9/9 (100%)	-0.33	0 100 100	53, 64, 90, 97	0
All	All	770/829 (92%)	-0.30	8 (1%) 84 77	34, 90, 142, 175	0

The worst 5 of 8 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	501	ARG	2.6
1	A	588	ILE	2.5
1	A	229	SER	2.3
1	A	108	GLY	2.1
1	A	409	ALA	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
5	ADP	A	1756	27/27	0.91	0.18	-0.54	50,83,100,103	0
6	MG	A	1757	1/1	0.90	0.93	-	93,93,93,93	0

6.5 Other polymers [i](#)

There are no such residues in this entry.