wwPDB X-ray Structure Validation Summary Report (i)

Jan 23, 2021 – 02:01 PM EST

PDB ID	:	1AM9
Title	:	HUMAN SREBP-1A BOUND TO LDL RECEPTOR PROMOTER
Authors	:	Parraga, A.; Burley, S.K.
Deposited on	:	1997-06-25
Resolution	:	2.30 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (i)) were used in the production of this report:

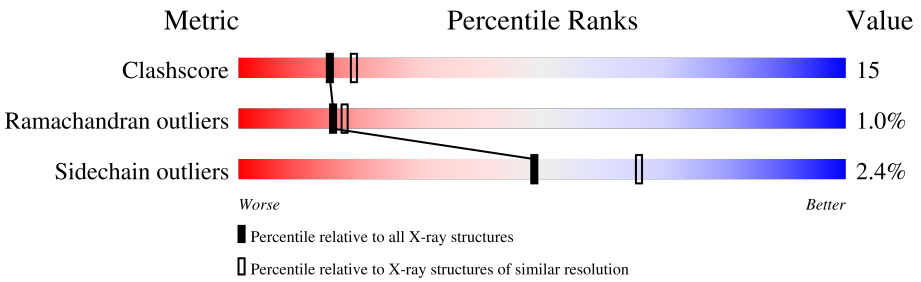
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.16

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 2.30 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))
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)

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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	E	17	29% 35% 35%
1	G	17	18% 41% 35% 6%
2	F	21	24% 43% 29% 5%
2	H	21	19% 52% 19% 10%
3	A	82	68% 29% •
3	B	82	56% 34% • 9%
3	C	82	72% 27% •
3	D	82	72% 20% • 7%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 4288 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 DNA chain called DNA (5'-D(*TP*TP*GP*CP*AP*GP*TP*GP*GP*GP*GP*TP*GP*AP*TP*CP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	E	17	351	168	63	104	16	0	0	0
1	G	17	351	168	63	104	16	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	F	21	422	202	80	120	20	0	0	0
2	H	21	422	202	80	120	20	0	0	0

- Molecule 3 is a protein called PROTEIN (STEROL REGULATORY ELEMENT BINDING PROTEIN 1A).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	A	80	621	386	122	113	0	0	0
3	B	75	587	365	113	109	0	0	0
3	C	82	640	398	125	117	0	0	0
3	D	76	593	369	114	110	0	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Mg 1 1	0	0
4	C	1	Total Mg 1 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	E	23	Total O 23 23	0	0
5	F	33	Total O 33 33	0	0
5	G	27	Total O 27 27	0	0
5	H	40	Total O 40 40	0	0
5	A	48	Total O 48 48	0	0
5	B	38	Total O 38 38	0	0
5	C	48	Total O 48 48	0	0
5	D	42	Total O 42 42	0	0

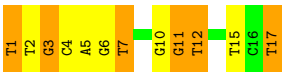
3 Residue-property plots (i)

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. 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. 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.

Note EDS was not executed.

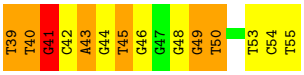
• Molecule 1: DNA (5'-D(*TP*TP*GP*CP*AP*GP*TP*GP*GP*GP*GP*TP*GP*AP*TP*CP *T)-3')

Chain E: 29% 35% 35%



• Molecule 1: DNA (5'-D(*TP*TP*GP*CP*AP*GP*TP*GP*GP*GP*GP*TP*GP*AP*TP*CP *T)-3')

Chain G: 18% 41% 35% 6%

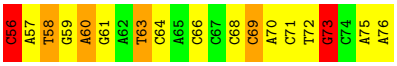


• Molecule 2: DNA (5'-D(*CP*AP*TP*GP*AP*GP*AP*TP*CP*AP*CP*CP*CP*CP*AP*CP *T P*GP*CP*AP*A)-3')

Chain F: 24% 43% 29% 5%

• Molecule 2: DNA (5'-D(*CP*AP*TP*GP*AP*GP*AP*TP*CP*AP*CP*CP*CP*CP*AP*CP *T P*GP*CP*AP*A)-3')

Chain H: 19% 52% 19% 10%



• Molecule 3: PROTEIN (STEROL REGULATORY ELEMENT BINDING PROTEIN 1A)

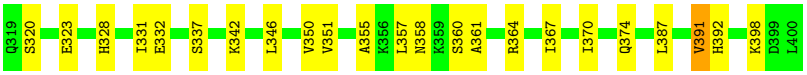
Chain A: 68% 29%

• Molecule 3: PROTEIN (STEROL REGULATORY ELEMENT BINDING PROTEIN 1A)

Chain B:

• Molecule 3: PROTEIN (STEROL REGULATORY ELEMENT BINDING PROTEIN 1A)

Chain C:



• Molecule 3: PROTEIN (STEROL REGULATORY ELEMENT BINDING PROTEIN 1A)

Chain D:

4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α, β, γ	94.63Å 94.63Å 459.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	6.00 – 2.30	Depositor
% Data completeness (in resolution range)	85.0 (6.00-2.30)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.219 , 0.278	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4288	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

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: MG

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	E	1.30	2/393 (0.5%)	1.93	16/607 (2.6%)
1	G	1.35	5/393 (1.3%)	1.83	13/607 (2.1%)
2	F	1.17	1/473 (0.2%)	1.99	19/726 (2.6%)
2	H	1.22	0/473	1.94	18/726 (2.5%)
3	A	0.62	0/626	0.72	0/838
3	B	0.66	0/591	0.79	1/791 (0.1%)
3	C	0.63	0/645	0.69	0/862
3	D	0.63	0/597	0.75	0/800
All	All	0.94	8/4191 (0.2%)	1.40	67/5957 (1.1%)

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	E	0	3
1	G	0	4
2	F	0	4
2	H	0	4
All	All	0	15

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	45	DT	C5-C7	6.41	1.53	1.50
1	E	1	DT	C5-C7	5.87	1.53	1.50
1	E	12	DT	C5-C7	5.36	1.53	1.50
1	G	39	DT	C5-C7	5.35	1.53	1.50
2	F	20	DT	C5-C7	5.24	1.53	1.50

The worst 5 of 67 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	F	22	DA	O4'-C1'-N9	9.98	114.99	108.00
2	F	38	DA	O4'-C1'-N9	9.79	114.85	108.00
2	F	18	DC	O4'-C1'-N1	8.62	114.04	108.00
2	H	56	DC	O4'-C1'-N1	7.82	113.48	108.00
1	E	15	DT	C6-C5-C7	-7.78	118.23	122.90

There are no chirality outliers.

5 of 15 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	1	DT	Sidechain
1	E	11	DG	Sidechain
1	E	3	DG	Sidechain
2	F	18	DC	Sidechain
2	F	19	DA	Sidechain

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	E	351	0	195	9	0
1	G	351	0	195	11	0
2	F	422	0	236	15	0
2	H	422	0	236	15	0
3	A	621	0	628	27	0
3	B	587	0	600	27	0
3	C	640	0	658	21	0
3	D	593	0	604	16	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
5	A	48	0	0	1	0
5	B	38	0	0	2	0
5	C	48	0	0	2	0
5	D	42	0	0	1	0
5	E	23	0	0	0	0
5	F	33	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	G	27	0	0	1	0
5	H	40	0	0	1	0
All	All	4288	0	3352	107	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 15.

The worst 5 of 107 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)
2:H:58:DT:H4'	2:H:58:DT:OP1	1.82	0.78
2:H:58:DT:H6	2:H:58:DT:H5"	1.47	0.78
1:E:7:DT:O4	3:D:328:HIS:HE1	1.70	0.74
3:B:339:ILE:HG21	5:B:2011:HOH:O	1.90	0.70
2:H:57:DA:H2"	2:H:58:DT:H5"	1.76	0.67

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
3	A	78/82 (95%)	75 (96%)	1 (1%)	2 (3%)	5 4
3	B	73/82 (89%)	69 (94%)	4 (6%)	0	100 100
3	C	80/82 (98%)	76 (95%)	3 (4%)	1 (1%)	12 12
3	D	74/82 (90%)	73 (99%)	1 (1%)	0	100 100
All	All	305/328 (93%)	293 (96%)	9 (3%)	3 (1%)	15 17

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	396	SER
3	A	394	SER
3	C	398	LYS

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	
3	A	63/74 (85%)	63 (100%)	0	100	100
3	B	61/74 (82%)	60 (98%)	1 (2%)	62	78
3	C	67/74 (90%)	65 (97%)	2 (3%)	41	57
3	D	61/74 (82%)	58 (95%)	3 (5%)	25	35
All	All	252/296 (85%)	246 (98%)	6 (2%)	49	66

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	391	VAL
3	D	382	GLN
3	D	325	ARG
3	C	337	SER
3	D	379	LYS

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

Mol	Chain	Res	Type
3	A	319	GLN
3	A	384	ASN
3	B	384	ASN
3	C	358	ASN
3	D	382	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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 (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)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.