



wwPDB X-ray Structure Validation Summary Report

Mar 1, 2014 – 12:10 AM GMT

PDB ID : 3CD7
Title : Thermodynamic and structure guided design of statin hmg-coa reductase inhibitors
Authors : Pavlovsky, A.; Sarver, R.W.; Harris, M.S.; Finzel, B.C.
Deposited on : 2008-02-26
Resolution : 2.05 Å(reported)

This is a wwPDB 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/ValidationPDFNotes.html>

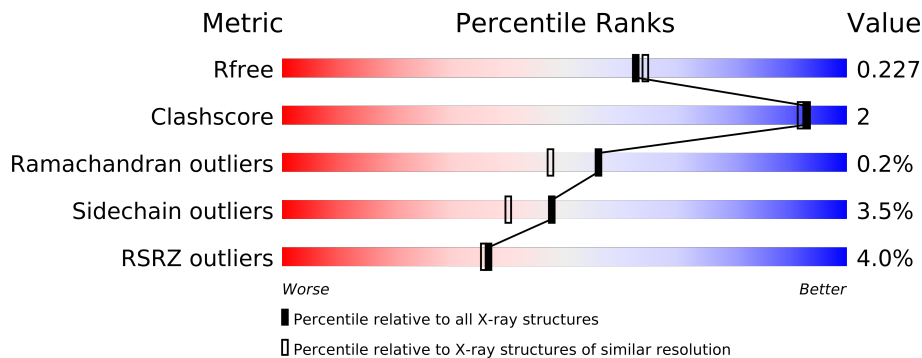
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.05 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}	66092	1380 (2.06-2.02)
Clashscore	79885	1577 (2.06-2.02)
Ramachandran outliers	78287	1565 (2.06-2.02)
Sidechain outliers	78261	1565 (2.06-2.02)
RSRZ outliers	66119	1381 (2.06-2.02)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	441	
1	B	441	
1	C	441	
1	D	441	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12542 atoms, of which 0 are hydrogen and 0 are deuterium.

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 3-hydroxy-3-methylglutaryl-coenzymeA reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	392	Total	C	N	O	S	0	0	0
			2903	1806	511	557	29			
1	B	401	Total	C	N	O	S	0	0	0
			2979	1859	524	567	29			
1	C	416	Total	C	N	O	S	0	0	0
			3087	1923	540	594	30			
1	D	394	Total	C	N	O	S	0	0	0
			2920	1818	514	559	29			

There are 28 discrepancies between the modelled and reference sequences:

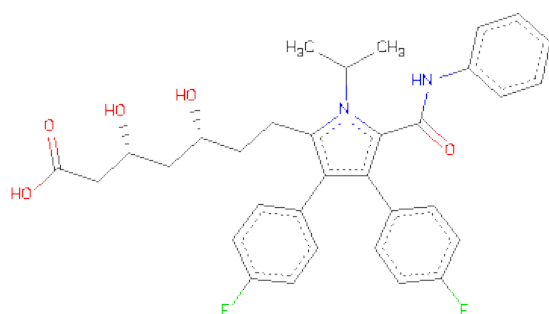
Chain	Residue	Modelled	Actual	Comment	Reference
A	435	HIS	-	expression tag	UNP P04035
A	436	HIS	-	expression tag	UNP P04035
A	437	HIS	-	expression tag	UNP P04035
A	438	HIS	-	expression tag	UNP P04035
A	439	HIS	-	expression tag	UNP P04035
A	440	HIS	-	expression tag	UNP P04035
A	485	ILE	MET	engineered	UNP P04035
B	435	HIS	-	expression tag	UNP P04035
B	436	HIS	-	expression tag	UNP P04035
B	437	HIS	-	expression tag	UNP P04035
B	438	HIS	-	expression tag	UNP P04035
B	439	HIS	-	expression tag	UNP P04035
B	440	HIS	-	expression tag	UNP P04035
B	485	ILE	MET	engineered	UNP P04035
C	435	HIS	-	expression tag	UNP P04035
C	436	HIS	-	expression tag	UNP P04035
C	437	HIS	-	expression tag	UNP P04035
C	438	HIS	-	expression tag	UNP P04035
C	439	HIS	-	expression tag	UNP P04035
C	440	HIS	-	expression tag	UNP P04035
C	485	ILE	MET	engineered	UNP P04035

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Chain	Residue	Modelled	Actual	Comment	Reference
D	435	HIS	-	expression tag	UNP P04035
D	436	HIS	-	expression tag	UNP P04035
D	437	HIS	-	expression tag	UNP P04035
D	438	HIS	-	expression tag	UNP P04035
D	439	HIS	-	expression tag	UNP P04035
D	440	HIS	-	expression tag	UNP P04035
D	485	ILE	MET	engineered	UNP P04035

- Molecule 2 is (3R,5R)-7-[5-(ANILINOCARBONYL)-3,4-BIS(4-FLUOROPHENYL)-1-ISOPROPYL-1H-PYRROL-2-YL]-3,5-DIHYDROXYHEPTANOICACID (three-letter code: 882) (formula: C₃₃H₃₄F₂N₂O₅).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			42	33	2	2	5		
2	A	1	Total	C	F	N	O	0	0
			42	33	2	2	5		
2	C	1	Total	C	F	N	O	0	0
			42	33	2	2	5		
2	D	1	Total	C	F	N	O	0	0
			42	33	2	2	5		

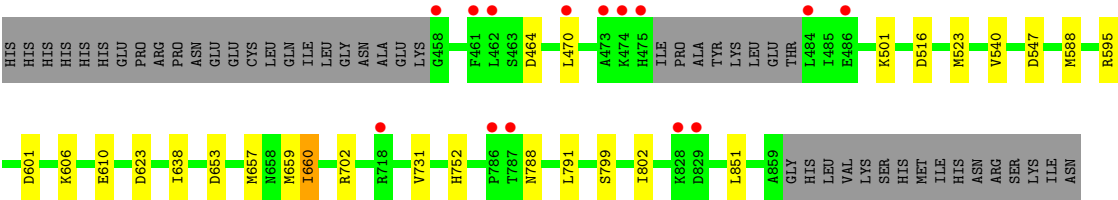
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	132	Total	O	0	0
			132	132		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	113	Total 113	O 113	0	0
3	C	109	Total 109	O 109	0	0
3	D	131	Total 131	O 131	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.73Å 176.96Å 76.73Å 90.00° 118.83° 90.00°	Depositor
Resolution (Å)	30.00 – 2.05 19.99 – 2.22	Depositor EDS
% Data completeness (in resolution range)	79.7 (30.00-2.05) 74.9 (19.99-2.22)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 2.21Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.223 , 0.263 0.227 , 0.227	Depositor DCC
R_{free} test set	3233 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	26.4	Xtriage
Anisotropy	0.250	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 30.4	EDS
Estimated twinning fraction	0.012 for -h-l,k,h 0.012 for l,k,-h-l 0.029 for h,-k,-h-l 0.034 for -h-l,-k,l 0.029 for l,-k,h	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 64413 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	12542	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section:
882

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	0.36	0/2943	0.63	4/3977 (0.1%)
1	B	0.34	0/3022	0.61	6/4084 (0.1%)
1	C	0.36	0/3130	0.62	4/4230 (0.1%)
1	D	0.36	0/2960	0.64	6/3999 (0.2%)
All	All	0.36	0/12055	0.62	20/16290 (0.1%)

There are no bond length outliers.

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	623	ASP	CB-CG-OD2	5.83	123.54	118.30
1	B	623	ASP	CB-CG-OD2	5.81	123.53	118.30
1	C	790	ASP	CB-CG-OD2	5.68	123.41	118.30
1	C	623	ASP	CB-CG-OD2	5.66	123.39	118.30
1	A	623	ASP	CB-CG-OD2	5.62	123.36	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2903	0	2933	12	0
1	B	2979	0	3020	10	0
1	C	3087	0	3121	15	0
1	D	2920	0	2957	6	0
2	A	84	0	66	1	0
2	C	42	0	33	1	0
2	D	42	0	33	1	0
3	A	132	0	0	0	0
3	B	113	0	0	0	0
3	C	109	0	0	0	0
3	D	131	0	0	0	0
All	All	12542	0	12163	41	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 2.

The worst 5 of 41 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:655:MET:SD	1:A:657:MET:HG2	2.27	0.74
1:A:771:ASN:OD1	1:A:775:SER:OG	2.10	0.69
1:C:681:LEU:HD22	1:D:731:VAL:HG22	1.73	0.69
1:C:655:MET:SD	1:C:657:MET:HG2	2.35	0.67
1:C:445:ASN:HD22	1:C:469:GLN:HE21	1.44	0.65

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	388/441 (88%)	374 (96%)	13 (3%)	1 (0%)	50	39
1	B	397/441 (90%)	381 (96%)	15 (4%)	1 (0%)	50	39
1	C	412/441 (93%)	391 (95%)	20 (5%)	1 (0%)	56	47
1	D	390/441 (88%)	375 (96%)	15 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1587/1764 (90%)	1521 (96%)	63 (4%)	3 (0%)	56 47

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	445	ASN
1	B	525	ALA
1	A	860	GLY

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	310/355 (87%)	298 (96%)	12 (4%)	43 35
1	B	318/355 (90%)	306 (96%)	12 (4%)	44 36
1	C	330/355 (93%)	321 (97%)	9 (3%)	57 51
1	D	312/355 (88%)	300 (96%)	12 (4%)	44 36
All	All	1270/1420 (89%)	1225 (96%)	45 (4%)	48 40

5 of 45 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	669	SER
1	C	452	LEU
1	D	788	ASN
1	B	771	ASN
1	C	460	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	488	HIS
1	C	679	GLN
1	D	632	GLN
1	C	469	GLN

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Mol	Chain	Res	Type
1	C	472	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

4 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	882	A	1	-	45,45,45	1.03	3 (6%)	63,63,63	2.14	11 (17%)
2	882	A	2	-	45,45,45	0.96	2 (4%)	63,63,63	1.98	12 (19%)
2	882	C	4	-	45,45,45	0.99	2 (4%)	63,63,63	1.91	12 (19%)
2	882	D	3	-	45,45,45	1.05	4 (8%)	63,63,63	2.16	10 (15%)

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	882	A	1	-	-	1/30/33/33	0/4/4/4
2	882	A	2	-	-	1/30/33/33	0/4/4/4
2	882	C	4	-	-	1/30/33/33	0/4/4/4
2	882	D	3	-	-	1/30/33/33	0/4/4/4

The worst 5 of 11 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	3	882	C2-N1	-3.45	1.36	1.40
2	A	1	882	C2-N1	-3.44	1.36	1.40
2	A	2	882	C2-N1	-3.35	1.36	1.40
2	C	4	882	C2-N1	-3.32	1.36	1.40
2	D	3	882	C1-C2	-2.57	1.38	1.42

The worst 5 of 45 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	882	C3-C2-N1	11.46	129.91	123.32
2	D	3	882	C3-C2-N1	10.95	129.62	123.32
2	A	2	882	C3-C2-N1	9.98	129.06	123.32
2	C	4	882	C3-C2-N1	9.60	128.84	123.32
2	D	3	882	C7-C4-C5	-4.43	124.98	130.95

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	4	882	O1-C3-C2-C1
2	A	2	882	O1-C3-C2-C1
2	D	3	882	O1-C3-C2-C1
2	A	1	882	O1-C3-C2-C1

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	392/441 (88%)	0.09	11 (2%) 50 51	22, 32, 55, 70	0
1	B	401/441 (90%)	0.33	27 (6%) 17 15	22, 37, 71, 89	0
1	C	416/441 (94%)	0.00	12 (2%) 49 49	19, 31, 60, 85	0
1	D	394/441 (89%)	-0.03	14 (3%) 41 40	19, 31, 53, 90	0
All	All	1603/1764 (90%)	0.10	64 (3%) 36 36	19, 33, 58, 90	0

The worst 5 of 64 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	479	TYR	11.6
1	B	861	HIS	7.8
1	B	524	GLY	6.8
1	D	461	PHE	6.4
1	A	861	HIS	6.2

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	882	C	4	42/42	0.14	1.55	27,35,37,38	0
2	882	A	2	42/42	0.12	0.01	25,31,35,39	0
2	882	D	3	42/42	0.10	-0.09	21,27,32,32	0
2	882	A	1	42/42	0.12	-0.18	24,35,41,44	0

6.5 Other polymers ⓘ

There are no such residues in this entry.