



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 10:40 AM GMT

PDB ID : 3MP5
Title : Crystal Structure of Human Lyase R41M in complex with HMG-CoA
Authors : Fu, Z.; Runquist, J.A.; Montgomery, C.; Miziorko, H.M.; Kim, J.-J.P.
Deposited on : 2010-04-24
Resolution : 2.25 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

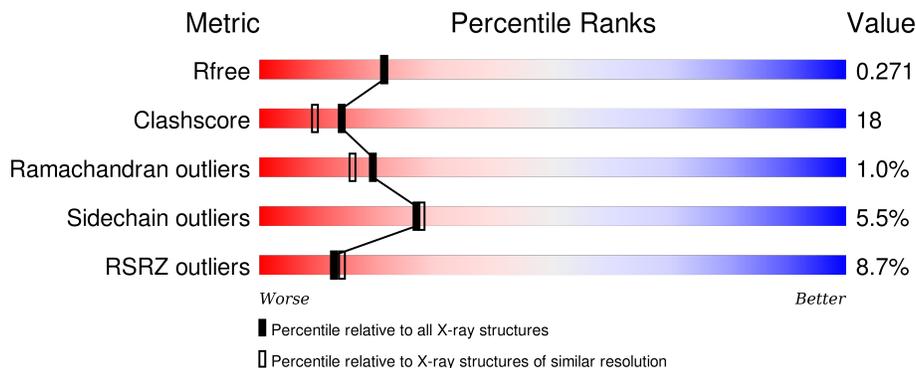
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1640 (2.28-2.24)
Clashscore	102246	1095 (2.26-2.26)
Ramachandran outliers	100387	1063 (2.26-2.26)
Sidechain outliers	100360	1063 (2.26-2.26)
RSRZ outliers	91569	1647 (2.28-2.24)

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	298	 4% 75% 21% ..
1	B	298	 7% 65% 31% ..
1	C	298	 12% 59% 35% ..
1	D	298	 5% 68% 26% ..
1	E	298	 4% 67% 28% ..

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Mol	Chain	Length	Quality of chain
1	F	298	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into three segments: a red segment on the left labeled '19%', a green segment in the middle labeled '59%', and a yellow segment on the right labeled '36%'. There are two small black dots at the far right end of the bar.</p>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13438 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.

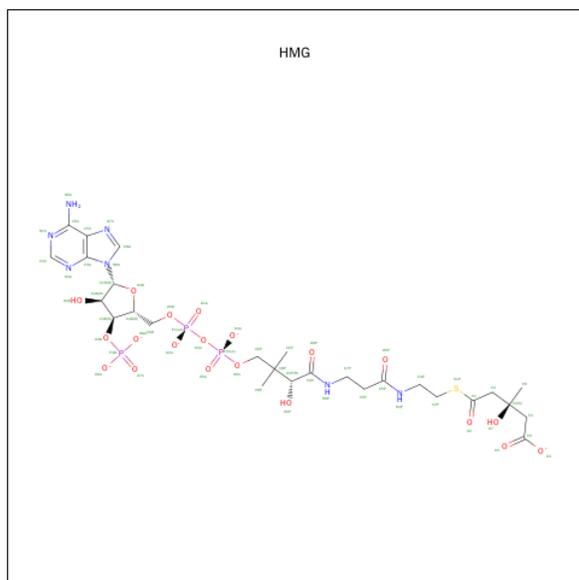
- Molecule 1 is a protein called Hydroxymethylglutaryl-CoA lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	296	2186	1389	362	417	18	0	0	0
1	B	296	2186	1389	362	417	18	0	0	0
1	C	288	2134	1357	353	407	17	0	0	0
1	D	294	2173	1382	359	414	18	0	0	0
1	E	296	2186	1389	362	417	18	0	0	0
1	F	294	2169	1377	360	414	18	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	41	MET	ARG	ENGINEERED	UNP P35914
B	41	MET	ARG	ENGINEERED	UNP P35914
C	41	MET	ARG	ENGINEERED	UNP P35914
D	41	MET	ARG	ENGINEERED	UNP P35914
E	41	MET	ARG	ENGINEERED	UNP P35914
F	41	MET	ARG	ENGINEERED	UNP P35914

- Molecule 2 is 3-HYDROXY-3-METHYLGLUTARYL-COENZYME A (three-letter code: HMG) (formula: C₂₇H₃₉N₇O₂₀P₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	N	O	P			S
2	B	1	58	27	7	20	3	1	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

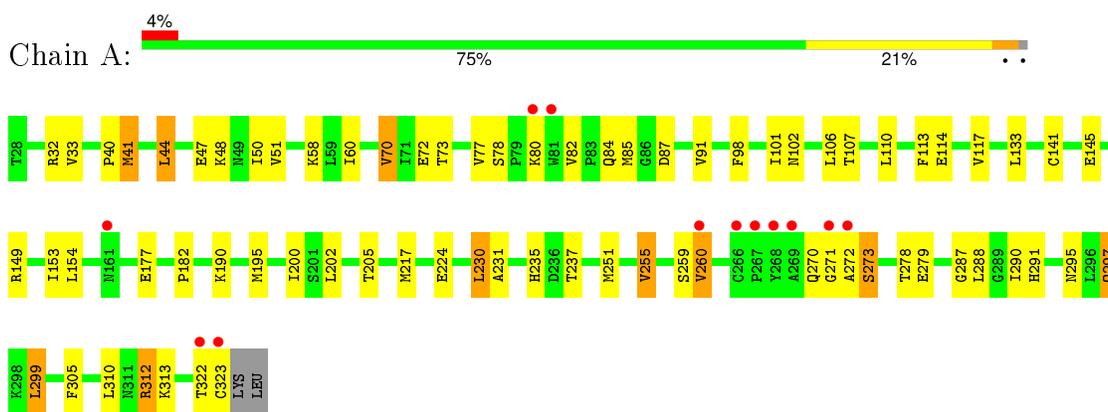
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	105	105	105	0	0
4	B	62	62	62	0	0
4	C	30	30	30	0	0
4	D	66	66	66	0	0
4	E	59	59	59	0	0
4	F	22	22	22	0	0

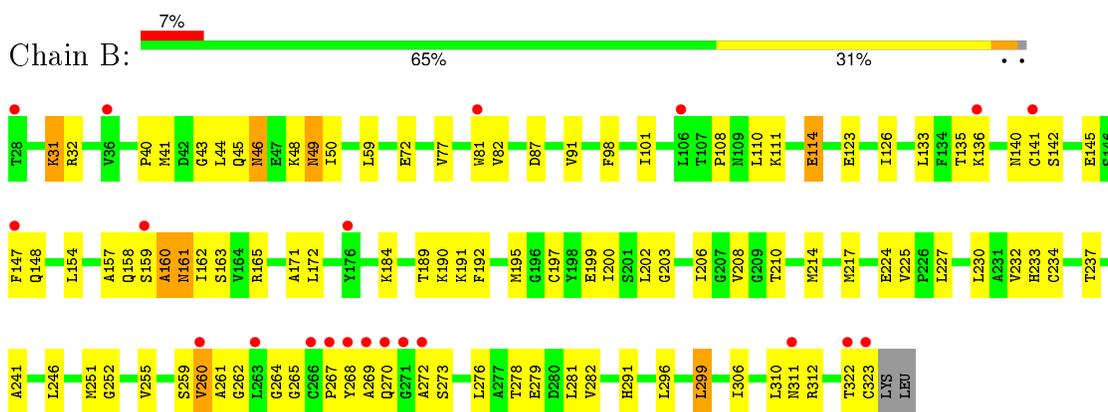
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

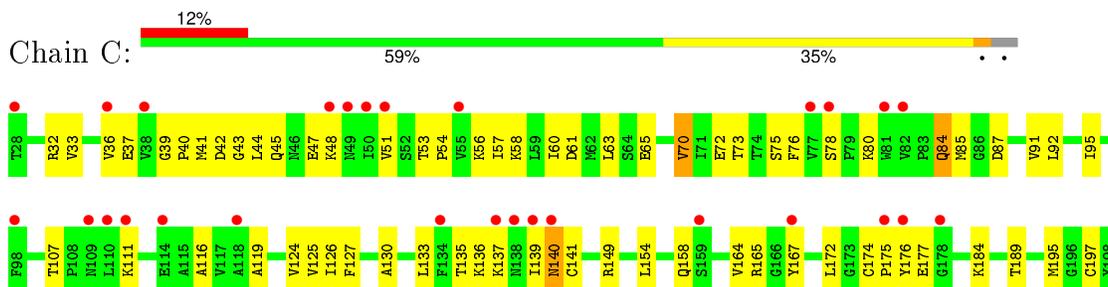
- Molecule 1: Hydroxymethylglutaryl-CoA lyase

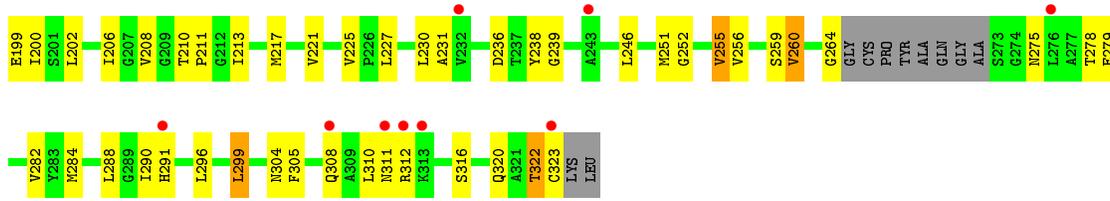


- Molecule 1: Hydroxymethylglutaryl-CoA lyase

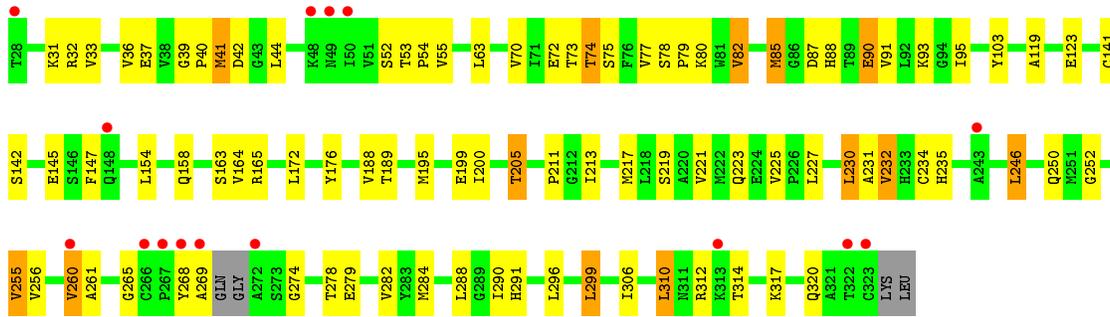


- Molecule 1: Hydroxymethylglutaryl-CoA lyase

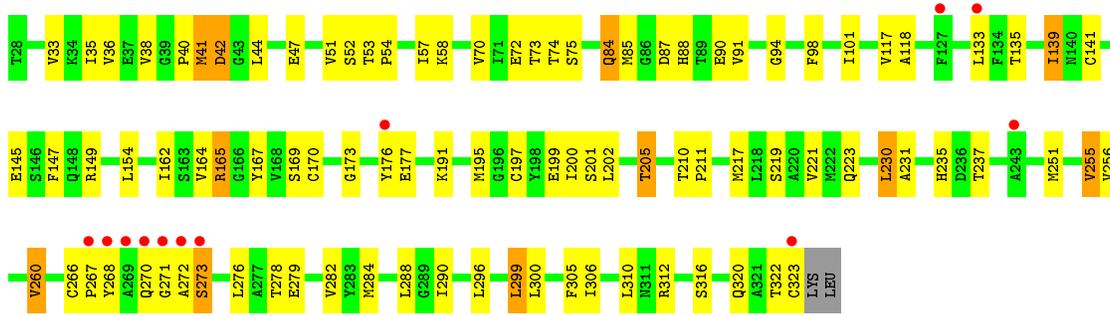




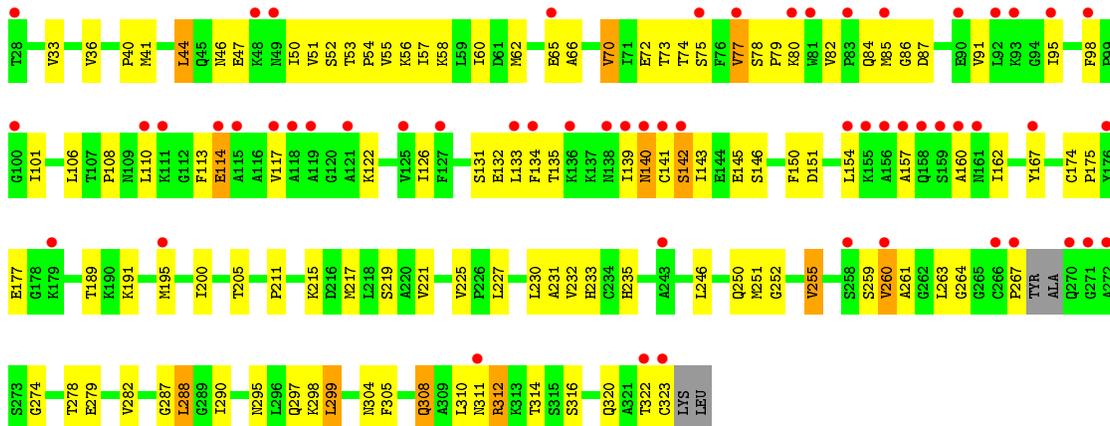
• Molecule 1: Hydroxymethylglutaryl-CoA lyase



• Molecule 1: Hydroxymethylglutaryl-CoA lyase



• Molecule 1: Hydroxymethylglutaryl-CoA lyase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	197.98Å 116.66Å 86.52Å 90.00° 112.16° 90.00°	Depositor
Resolution (Å)	28.35 – 2.25 28.35 – 2.25	Depositor EDS
% Data completeness (in resolution range)	94.8 (28.35-2.25) 94.9 (28.35-2.25)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 2.24Å)	Xtrriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.221 , 0.271 0.221 , 0.271	Depositor DCC
R_{free} test set	8219 reflections (10.04%)	DCC
Wilson B-factor (Å ²)	29.3	Xtrriage
Anisotropy	0.129	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.3	EDS
Estimated twinning fraction	No twinning to report.	Xtrriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Outliers	1 of 81864 reflections (0.001%)	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13438	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.35% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.40	0/2221	0.63	0/3007
1	B	0.38	0/2221	0.62	0/3007
1	C	0.35	0/2166	0.56	0/2930
1	D	0.37	0/2207	0.59	0/2987
1	E	0.37	0/2221	0.59	0/3007
1	F	0.33	0/2202	0.56	0/2979
All	All	0.37	0/13238	0.59	0/17917

There are no bond length outliers.

There are no bond angle outliers.

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	2186	0	2240	59	0
1	B	2186	0	2240	81	0
1	C	2134	0	2194	92	0
1	D	2173	0	2228	76	0
1	E	2186	0	2240	80	0
1	F	2169	0	2225	107	0
2	B	58	0	38	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	105	0	0	4	0
4	B	62	0	0	2	0
4	C	30	0	0	1	0
4	D	66	0	0	0	0
4	E	59	0	0	1	0
4	F	22	0	0	1	0
All	All	13438	0	13405	484	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 18.

The worst 5 of 484 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:D:73:THR:HG22	1:D:74:THR:HG22	1.32	1.11
1:E:139:ILE:H	1:E:139:ILE:HD13	1.30	0.95
1:D:36:VAL:HA	1:D:70:VAL:HG13	1.49	0.95
1:C:288:LEU:HD11	1:D:211:PRO:HD2	1.47	0.95
1:C:211:PRO:HG2	1:D:288:LEU:HD21	1.49	0.93

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
1	A	294/298 (99%)	281 (96%)	10 (3%)	3 (1%)	19 16
1	B	294/298 (99%)	283 (96%)	8 (3%)	3 (1%)	19 16
1	C	284/298 (95%)	265 (93%)	16 (6%)	3 (1%)	17 13

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	290/298 (97%)	275 (95%)	14 (5%)	1 (0%)	46	52
1	E	294/298 (99%)	282 (96%)	9 (3%)	3 (1%)	19	16
1	F	290/298 (97%)	258 (89%)	28 (10%)	4 (1%)	14	9
All	All	1746/1788 (98%)	1644 (94%)	85 (5%)	17 (1%)	19	16

5 of 17 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	260	VAL
1	A	273	SER
1	B	160	ALA
1	B	260	VAL
1	C	260	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	238/240 (99%)	228 (96%)	10 (4%)	36	42
1	B	238/240 (99%)	223 (94%)	15 (6%)	22	21
1	C	234/240 (98%)	223 (95%)	11 (5%)	32	36
1	D	237/240 (99%)	221 (93%)	16 (7%)	20	18
1	E	238/240 (99%)	225 (94%)	13 (6%)	27	27
1	F	237/240 (99%)	224 (94%)	13 (6%)	27	27
All	All	1422/1440 (99%)	1344 (94%)	78 (6%)	27	27

5 of 78 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	312	ARG
1	D	205	THR
1	F	288	LEU
1	D	41	MET

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Mol	Chain	Res	Type
1	D	85	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 58 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	250	GLN
1	D	161	ASN
1	F	297	GLN
1	C	275	ASN
1	C	297	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 3 ligands modelled in this entry, 2 are 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$
2	HMG	B	401	3	47,60,60	3.80	22 (46%)	58,90,90	3.49	22 (37%)

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	HMG	B	401	3	-	0/54/77/77	0/3/3/3

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	HMG	O7-C3	-5.32	1.35	1.44
2	B	401	HMG	C6-C3	-2.80	1.49	1.52
2	B	401	HMG	CEP-CBP	-2.62	1.48	1.53
2	B	401	HMG	C7P-N8P	-2.58	1.40	1.46
2	B	401	HMG	C2P-C3P	-2.05	1.42	1.51

The worst 5 of 22 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	HMG	O2-C1-S1P	-12.61	112.83	122.83
2	B	401	HMG	N3A-C2A-N1A	-12.48	119.34	128.89
2	B	401	HMG	C1B-N9A-C4A	-4.41	120.29	126.94
2	B	401	HMG	CDP-CBP-CCP	-3.93	103.41	108.50
2	B	401	HMG	OAP-CAP-C9P	-3.88	101.47	110.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	HMG	2	0

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	296/298 (99%)	0.07	12 (4%) 41 44	14, 25, 49, 65	0
1	B	296/298 (99%)	0.45	21 (7%) 19 20	17, 32, 56, 64	0
1	C	288/298 (96%)	0.78	37 (12%) 5 5	22, 43, 65, 73	0
1	D	294/298 (98%)	0.17	15 (5%) 32 35	16, 32, 54, 70	0
1	E	296/298 (99%)	0.18	12 (4%) 41 44	18, 31, 56, 69	0
1	F	294/298 (98%)	0.99	57 (19%) 1 1	24, 50, 68, 78	0
All	All	1764/1788 (98%)	0.44	154 (8%) 13 14	14, 34, 62, 78	0

The worst 5 of 154 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	140	ASN	9.6
1	E	271	GLY	7.3
1	F	118	ALA	6.4
1	D	268	TYR	5.8
1	C	50	ILE	5.7

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(Å ²)	Q<0.9
2	HMG	B	401	58/58	0.64	0.31	1.91	54,78,88,88	0
3	MG	C	403	1/1	0.81	0.32	-	42,42,42,42	0
3	MG	B	402	1/1	0.94	0.19	-	29,29,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.