Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 11:56 pm BST

PDB ID	:	2HMB
Title	:	THREE-DIMENSIONAL STRUCTURE OF RECOMBINANT HUMAN MUSCLE FATTY ACID-BINDING PROTEIN
Authors	:	Zanotti, G.; Scapin, G.; Spadon, P.; Veerkamp, J.H.; Sacchettini, J.C.
Deposited on	:	1992-09-11
Resolution	:	2.10 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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 ⓘ symbol.

The following versions of software and data (see references ①) were used in the production of this report:

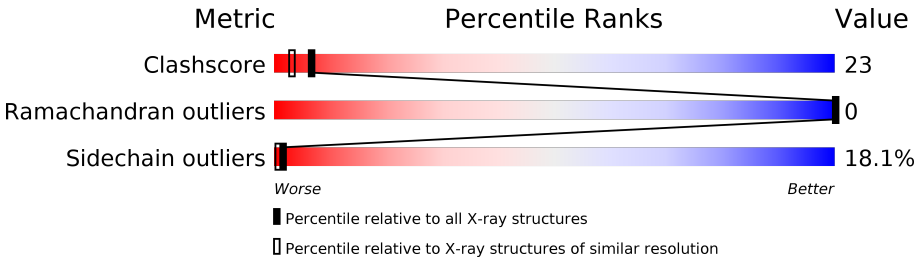
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 2.10 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	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)

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 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	A	132	52% 31% 15%

2 Entry composition

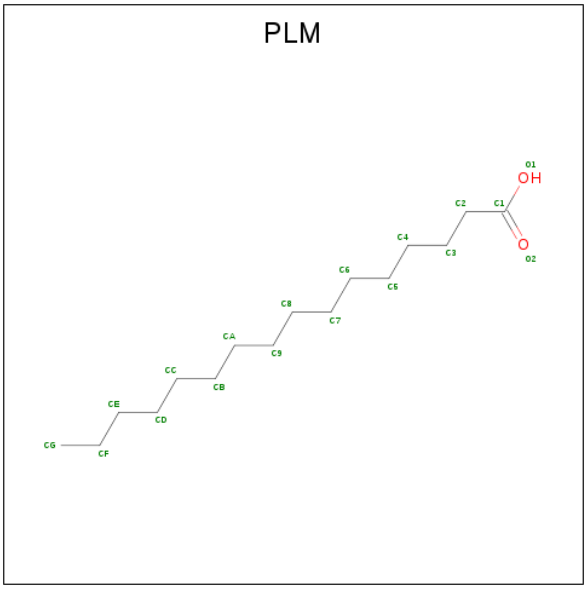
There are 3 unique types of molecules in this entry. The entry contains 1102 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 MUSCLE FATTY ACID BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	131	Total	C	N	O	S	0	0	0
			1030	654	171	202	3			

- Molecule 2 is PALMITIC ACID (three-letter code: PLM) (formula: $C_{16}H_{32}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			18	16	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	54	Total	O	0	0
			54	54		

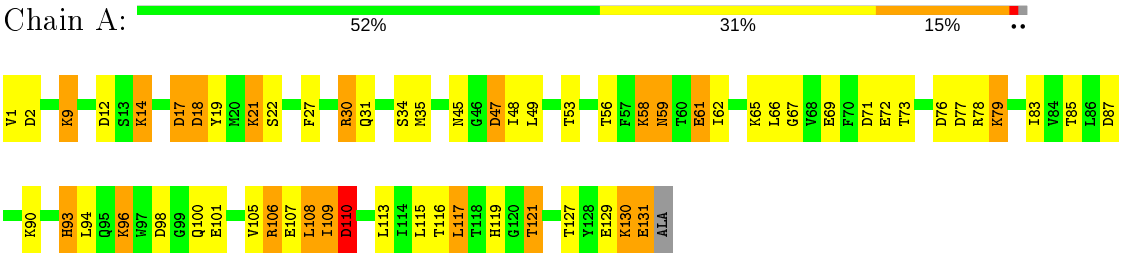
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: MUSCLE FATTY ACID BINDING PROTEIN

Chain A:



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	35.40Å 56.70Å 72.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	9.00 – 2.10	Depositor
% Data completeness (in resolution range)	(Not available) (9.00-2.10)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT, X-PLOR	Depositor
R, R_{free}	0.195 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1102	wwPDB-VP
Average B, all atoms (Å^2)	24.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.93	7/1045 (0.7%)	1.65	29/1410 (2.1%)

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	A	1	0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	61	GLU	CD-OE2	5.88	1.32	1.25
1	A	101	GLU	CD-OE2	5.77	1.31	1.25
1	A	72	GLU	CD-OE2	5.69	1.31	1.25
1	A	129	GLU	CD-OE1	5.57	1.31	1.25
1	A	107	GLU	CD-OE2	5.57	1.31	1.25
1	A	131	GLU	CD-OE2	5.53	1.31	1.25
1	A	69	GLU	CD-OE1	5.06	1.31	1.25

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	53	THR	CA-CB-CG2	15.57	134.19	112.40
1	A	106	ARG	NE-CZ-NH1	15.46	128.03	120.30
1	A	53	THR	OG1-CB-CG2	-13.82	78.21	110.00
1	A	106	ARG	NE-CZ-NH2	-11.44	114.58	120.30
1	A	78	ARG	NE-CZ-NH1	9.65	125.12	120.30
1	A	106	ARG	CD-NE-CZ	8.97	136.16	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	71	ASP	CB-CG-OD1	8.11	125.60	118.30
1	A	121	THR	CA-CB-CG2	7.48	122.87	112.40
1	A	71	ASP	CB-CG-OD2	-7.38	111.66	118.30
1	A	18	ASP	CB-CG-OD2	-6.96	112.04	118.30
1	A	78	ARG	NE-CZ-NH2	-6.51	117.05	120.30
1	A	76	ASP	CB-CG-OD2	6.37	124.04	118.30
1	A	76	ASP	CB-CG-OD1	-6.36	112.58	118.30
1	A	30	ARG	NE-CZ-NH2	-6.33	117.14	120.30
1	A	2	ASP	CB-CG-OD2	-6.08	112.83	118.30
1	A	17	ASP	CB-CG-OD2	-6.06	112.85	118.30
1	A	18	ASP	CB-CG-OD1	5.84	123.56	118.30
1	A	17	ASP	CB-CG-OD1	5.79	123.51	118.30
1	A	47	ASP	CB-CG-OD2	-5.66	113.21	118.30
1	A	107	GLU	N-CA-CB	5.55	120.59	110.60
1	A	110	ASP	CB-CG-OD2	5.32	123.09	118.30
1	A	2	ASP	CB-CG-OD1	5.31	123.08	118.30
1	A	98	ASP	CB-CG-OD2	-5.31	113.52	118.30
1	A	12	ASP	CB-CG-OD1	5.29	123.06	118.30
1	A	77	ASP	CB-CG-OD2	5.26	123.03	118.30
1	A	87	ASP	CB-CG-OD2	5.22	123.00	118.30
1	A	110	ASP	N-CA-CB	5.17	119.90	110.60
1	A	56	THR	CA-CB-OG1	5.06	119.62	109.00
1	A	56	THR	CA-CB-CG2	5.03	119.44	112.40

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	56	THR	CB

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	1030	0	1052	44	3
2	A	18	0	31	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	54	0	0	14	1
All	All	1102	0	1083	50	4

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 23.

All (50) 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:61:GLU:OE2	3:A:228:HOH:O	1.54	1.19
1:A:47:ASP:OD1	3:A:207:HOH:O	1.76	1.03
1:A:1:VAL:N	3:A:203:HOH:O	2.00	0.94
1:A:130:LYS:HE2	3:A:204:HOH:O	1.69	0.92
2:A:133:PLM:HG1	2:A:133:PLM:H22	1.50	0.91
1:A:9:LYS:NZ	1:A:9:LYS:HB3	1.98	0.76
1:A:131:GLU:O	3:A:202:HOH:O	2.03	0.76
2:A:133:PLM:HG1	2:A:133:PLM:C2	2.16	0.75
1:A:100:GLN:HE21	1:A:119:HIS:HE1	1.35	0.75
1:A:90:LYS:HD2	3:A:212:HOH:O	1.89	0.72
1:A:62:ILE:HA	3:A:233:HOH:O	1.89	0.71
1:A:18:ASP:HA	1:A:21:LYS:HD2	1.74	0.69
1:A:100:GLN:HE21	1:A:119:HIS:CE1	2.13	0.67
1:A:9:LYS:HZ2	1:A:9:LYS:HB3	1.59	0.65
1:A:130:LYS:CE	3:A:204:HOH:O	2.35	0.64
1:A:18:ASP:O	1:A:21:LYS:HD3	1.98	0.64
1:A:130:LYS:CD	3:A:204:HOH:O	2.48	0.61
1:A:67:GLY:HA2	1:A:83:ILE:HD11	1.79	0.61
1:A:61:GLU:CG	3:A:228:HOH:O	2.47	0.61
2:A:133:PLM:H22	2:A:133:PLM:CG	2.29	0.60
1:A:17:ASP:OD1	1:A:30:ARG:HD2	2.02	0.59
2:A:133:PLM:HG1	2:A:133:PLM:C3	2.32	0.59
1:A:1:VAL:HG21	1:A:49:LEU:HD11	1.86	0.58
1:A:108:LEU:CD1	1:A:113:LEU:HD12	2.36	0.55
1:A:106:ARG:NH2	3:A:140:HOH:O	2.40	0.54
1:A:14:LYS:HD2	1:A:14:LYS:O	2.10	0.51
1:A:110:ASP:OD1	1:A:110:ASP:O	2.27	0.51
1:A:48:ILE:HD11	1:A:65:LYS:HE2	1.93	0.51
1:A:108:LEU:HD13	1:A:113:LEU:HD12	1.92	0.51
1:A:22:SER:OG	1:A:119:HIS:HD2	1.94	0.50
1:A:58:LYS:C	1:A:59:ASN:HD22	2.15	0.50
1:A:48:ILE:CD1	1:A:65:LYS:HG2	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:HIS:HD2	3:A:140:HOH:O	1.94	0.49
1:A:31:GLN:O	1:A:34:SER:OG	2.33	0.47
1:A:130:LYS:HD2	3:A:204:HOH:O	2.12	0.47
1:A:94:LEU:HG	1:A:96:LYS:HE3	1.96	0.46
1:A:90:LYS:CE	3:A:212:HOH:O	2.64	0.46
1:A:109:ILE:HG13	1:A:109:ILE:O	2.16	0.46
2:A:133:PLM:HG1	2:A:133:PLM:H31	1.95	0.45
2:A:133:PLM:HG2	2:A:133:PLM:H42	1.99	0.44
1:A:108:LEU:HD12	1:A:108:LEU:HA	1.63	0.44
1:A:100:GLN:NE2	1:A:119:HIS:CE1	2.84	0.43
1:A:108:LEU:CD1	1:A:113:LEU:CD1	2.97	0.43
1:A:73:THR:OG1	1:A:73:THR:O	2.31	0.43
1:A:59:ASN:HD22	1:A:59:ASN:N	2.17	0.42
1:A:9:LYS:CB	1:A:9:LYS:NZ	2.79	0.40
1:A:19:TYR:CE1	1:A:117:LEU:HB3	2.56	0.40
1:A:79:LYS:HG2	1:A:79:LYS:O	2.20	0.40
1:A:106:ARG:HG3	1:A:115:LEU:HD13	2.02	0.40
1:A:79:LYS:HE3	1:A:79:LYS:H	1.87	0.40

All (4) 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:27:PHE:CE2	1:A:90:LYS:NZ[2_665]	1.56	0.64
1:A:27:PHE:CZ	1:A:90:LYS:NZ[2_665]	2.10	0.10
3:A:240:HOH:O	3:A:242:HOH:O[3_756]	2.10	0.10
1:A:27:PHE:CD2	1:A:90:LYS:NZ[2_665]	2.14	0.06

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	129/132 (98%)	127 (98%)	2 (2%)	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	116/116 (100%)	95 (82%)	21 (18%)	1 1

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	LYS
1	A	14	LYS
1	A	21	LYS
1	A	35	MET
1	A	45	ASN
1	A	58	LYS
1	A	59	ASN
1	A	66	LEU
1	A	79	LYS
1	A	85	THR
1	A	93	HIS
1	A	96	LYS
1	A	105	VAL
1	A	108	LEU
1	A	109	ILE
1	A	110	ASP
1	A	116	THR
1	A	117	LEU
1	A	121	THR
1	A	127	THR
1	A	130	LYS

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

Mol	Chain	Res	Type
1	A	45	ASN
1	A	59	ASN
1	A	93	HIS
1	A	100	GLN
1	A	119	HIS

5.3.3 RNA

There are no RNA molecules 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

1 ligand is 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	PLM	A	133	-	14,17,17	0.13	0	13,17,17	0.41	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
2	PLM	A	133	-	-	8/13/15/15	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

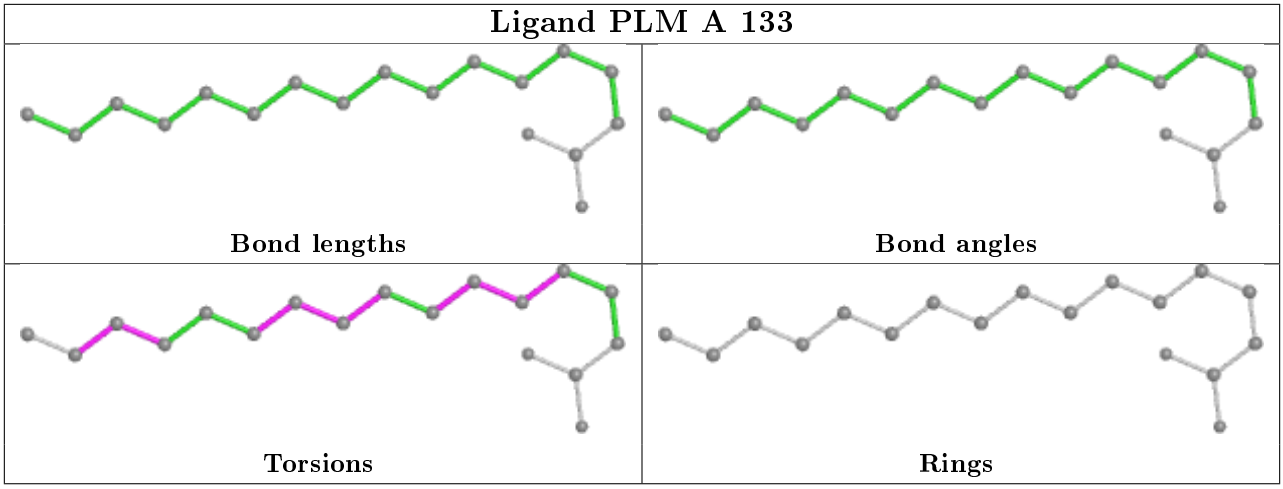
Mol	Chain	Res	Type	Atoms
2	A	133	PLM	C4-C5-C6-C7
2	A	133	PLM	C5-C6-C7-C8
2	A	133	PLM	C8-C9-CA-CB
2	A	133	PLM	C9-CA-CB-CC
2	A	133	PLM	CC-CD-CE-CF
2	A	133	PLM	CD-CE-CF-CG
2	A	133	PLM	C3-C4-C5-C6
2	A	133	PLM	C7-C8-C9-CA

There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	133	PLM	6	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.