

Full wwPDB X-ray Structure Validation Report (i)

Oct 31, 2021 – 11:40 PM EDT

PDB ID	:	1V3S
Title	:	Crystal structure of TT1020 from Thermus thermophilus HB8
Authors	:	Wang, H.; Sakai, H.; Takemoto-Hori, C.; Kaminishi, T.; Terada, T.;
		Kuramitsu, S.; Shirouzu, M.; Yokoyama, S.; RIKEN Structural Ge-
		nomics/Proteomics Initiative (RSGI)
Deposited on	:	2003-11-05
Resolution	:	1.85 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 (i) symbol.

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

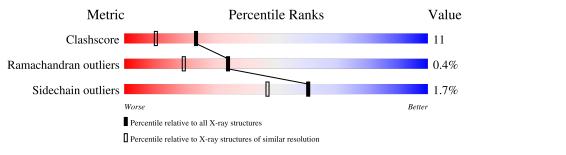
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 1.85 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$		
Clashscore	141614	2625 (1.86-1.86)		
Ramachandran outliers	138981	2592 (1.86-1.86)		
Sidechain outliers	138945	2592 (1.86-1.86)		

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 cha	in	
1	А	116	62%	22%	16%
1	В	116	65%	14% •	20%
1	С	116	59%	17% •	22%



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 2442 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.

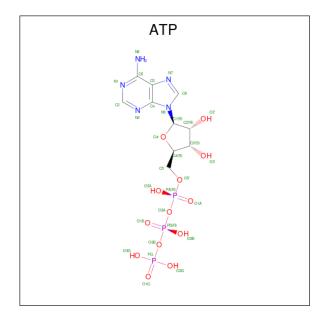
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	97	Total	С	Ν	0	S	0	0	0
1	Л	91	755	481	133	140	1	0	0	0
1	В	93	Total	С	Ν	0	S	0	0	0
	D	90	729	467	129	131	2	0	0	0
1	C	01	Total	С	Ν	0	S	0	0	0
		91	712	457	127	127	1	0		0

• Molecule 1 is a protein called Nitrogen regulatory protein P-II.

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	18	LYS	GLU	engineered mutation	UNP Q5SM86
В	18	LYS	GLU	engineered mutation	UNP Q5SM86
С	18	LYS	GLU	engineered mutation	UNP Q5SM86

• Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	۸	1	Total	С	Ν	Ο	Р	0	0
	A	1	31	10	5	13	3	0	0
0	р	1	Total	С	Ν	Ο	Р	0	0
	D	1	31	10	5	13	3	0	0
9	С	1	Total	С	Ν	Ο	Р	0	0
	U	1	31	10	5	13	3	0	0

• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	36	$\begin{array}{cc} \text{Total} & \text{O} \\ 36 & 36 \end{array}$	0	0
3	В	62	$\begin{array}{cc} \text{Total} & \text{O} \\ 62 & 62 \end{array}$	0	0
3	С	55	$\begin{array}{cc} \text{Total} & \text{O} \\ 55 & 55 \end{array}$	0	0



Residue-property plots (i) 3

• Molecule 1: Nitrogen regulatory protein P-II

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.

Chain A: 62% 22% 16% 3LU ARG 7AL 7AL 7HR 7HR 7HR 7HR 7HR 7HR 7HR 7HR 7HR PRO VAL GLN • Molecule 1: Nitrogen regulatory protein P-II Chain B: 65% 14% 20% ALA ALA ALA ALA /AL THR /AL • Molecule 1: Nitrogen regulatory protein P-II Chain C: 59% 17% 22%

Note EDS was not executed.



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	53.58Å 75.51Å 76.91Å	Depositor	
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor	
Resolution (Å)	37.99 - 1.85	Depositor	
% Data completeness	97.7 (37.99-1.85)	Depositor	
(in resolution range)	51.1 (51.55-1.65)	Depositor	
R_{merge}	0.05	Depositor	
R _{sym}	(Not available)	Depositor	
Refinement program	CNS 1.1	Depositor	
R, R_{free}	0.201 , 0.243	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	2442	wwPDB-VP	
Average B, all atoms $(Å^2)$	28.0	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: ATP

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		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.27	0/763	0.60	0/1025	
1	В	0.29	0/737	0.62	0/989	
1	С	0.28	0/720	0.60	0/967	
All	All	0.28	0/2220	0.61	0/2981	

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	А	755	0	797	22	0
1	В	729	0	779	14	0
1	С	712	0	764	24	0
2	А	31	0	12	1	0
2	В	31	0	12	0	0
2	С	31	0	12	3	0
3	А	36	0	0	1	0
3	В	62	0	0	0	0
3	С	55	0	0	2	0
All	All	2442	0	2376	52	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 11.

All (52) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:C:107:GLU:HG2	1:C:108:ASP:H	1.34	0.90
1:C:1:MET:HG2	1:C:99:VAL:HG21	1.65	0.78
1:C:107:GLU:CG	1:C:108:ASP:H	2.03	0.70
1:C:71:LYS:HZ2	1:C:71:LYS:H	1.40	0.69
1:B:82:ARG:HG3	1:C:102:ILE:O	1.97	0.65
1:A:75:GLU:HG3	1:A:79:LYS:HE3	1.79	0.65
1:C:107:GLU:HG2	1:C:108:ASP:N	2.09	0.65
1:A:71:LYS:HB2	1:A:72:PRO:HD3	1.78	0.65
1:A:26:ARG:HG2	1:A:26:ARG:HH11	1.63	0.62
1:A:15:GLU:HG3	3:A:223:HOH:O	2.01	0.58
1:A:12:LYS:HE3	1:A:83:THR:HG22	1.86	0.58
1:A:38:GLY:N	2:A:200:ATP:O2B	2.37	0.58
1:B:104:THR:OG1	1:B:106:GLU:HG2	2.05	0.57
1:C:71:LYS:H	1:C:71:LYS:NZ	2.03	0.56
1:A:67:GLU:HB2	1:A:68:PRO:HD3	1.86	0.56
1:C:108:ASP:HB3	3:C:452:HOH:O	2.07	0.55
1:A:98:LYS:NZ	1:A:107:GLU:HB2	2.22	0.55
1:B:54:GLU:CD	1:B:54:GLU:H	2.10	0.55
1:C:12:LYS:HD2	1:C:83:THR:HG22	1.89	0.54
1:C:86:VAL:HG13	2:C:400:ATP:O2B	2.07	0.54
1:A:82:ARG:HH11	1:A:82:ARG:HG2	1.75	0.51
1:C:1:MET:HG2	1:C:99:VAL:CG2	2.37	0.51
1:B:82:ARG:HD3	1:C:103:ARG:O	2.11	0.50
1:C:71:LYS:N	1:C:71:LYS:HD3	2.27	0.49
1:C:107:GLU:CG	1:C:108:ASP:N	2.74	0.49
1:A:96:VAL:HG22	1:C:94:LEU:CD2	2.42	0.49
1:C:61:LEU:HD12	1:C:61:LEU:N	2.28	0.49
1:C:67:GLU:HB2	1:C:68:PRO:HD3	1.95	0.48
1:A:14:ASN:O	1:A:18:LYS:HG3	2.14	0.47
1:C:71:LYS:HB2	1:C:72:PRO:HD3	1.96	0.47
1:A:12:LYS:O	1:A:16:VAL:HG23	2.15	0.47
1:A:54:GLU:OE1	1:A:54:GLU:N	2.49	0.45
1:B:36:HIS:HD2	1:C:26:ARG:O	2.00	0.45
1:B:67:GLU:HB2	1:B:68:PRO:HD3	1.99	0.44
1:C:87:GLY:N	2:C:400:ATP:O2B	2.47	0.44
1:B:69:PHE:C	1:B:72:PRO:HD2	2.37	0.44
1:A:26:ARG:HG2	1:A:26:ARG:NH1	2.31	0.44

Continued on next page...



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:LYS:HZ1	1:A:107:GLU:HB2	1.83	0.44
1:B:24:GLU:HA	1:B:26:ARG:HH12	1.83	0.43
1:A:82:ARG:HG2	1:A:82:ARG:NH1	2.33	0.43
1:A:101:ARG:HD3	2:C:400:ATP:O3G	2.18	0.42
1:B:12:LYS:HE3	1:B:80:ALA:O	2.20	0.42
1:A:30:LEU:C	1:A:30:LEU:HD12	2.39	0.42
1:A:55:LEU:HD12	1:A:55:LEU:N	2.35	0.41
1:B:95:PRO:HD3	1:C:97:GLU:HG3	2.02	0.41
1:A:38:GLY:HA2	1:B:26:ARG:HB2	2.00	0.41
1:A:86:VAL:HG22	1:B:103:ARG:NH1	2.36	0.41
1:B:53:MET:N	1:C:21:PHE:CE1	2.89	0.41
1:C:85:GLU:HG3	3:C:447:HOH:O	2.20	0.41
1:A:98:LYS:NZ	1:A:109:GLU:HG3	2.36	0.40
1:C:23:ALA:O	1:C:24:GLU:HB2	2.22	0.40
1:B:96:VAL:HG11	1:B:99:VAL:CG2	2.51	0.40

Continued from previous page...

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	А	93/116~(80%)	91~(98%)	2(2%)	0	100 100
1	В	89/116~(77%)	88~(99%)	1 (1%)	0	100 100
1	С	87/116 (75%)	86 (99%)	0	1 (1%)	14 4
All	All	269/348~(77%)	265~(98%)	3(1%)	1 (0%)	34 19

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	С	107	GLU



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	А	81/98~(83%)	80~(99%)	1 (1%)	71 62
1	В	79/98~(81%)	77~(98%)	2(2%)	47 31
1	С	77/98~(79%)	76~(99%)	1 (1%)	69 58
All	All	237/294~(81%)	233~(98%)	4 (2%)	60 47

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	1	MET
1	В	54	GLU
1	В	82	ARG
1	С	71	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	14	ASN
1	А	36	HIS
1	В	14	ASN
1	В	36	HIS
1	С	14	ASN
1	С	22	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)

3 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	Turne	Chain	Res	Link	Bo	ond leng	\mathbf{ths}	В	ond ang	les
IVIOI	Type	Chain	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
2	ATP	В	300	-	26,33,33	0.97	1 (3%)	$31,\!52,\!52$	1.75	7 (22%)
2	ATP	С	400	-	26,33,33	0.98	1 (3%)	31,52,52	1.75	7 (22%)
2	ATP	А	200	-	26,33,33	0.98	1 (3%)	31,52,52	1.74	7 (22%)

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	ATP	В	300	-	-	3/18/38/38	0/3/3/3
2	ATP	С	400	-	-	4/18/38/38	0/3/3/3
2	ATP	А	200	-	-	1/18/38/38	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	Observed(Å)	Ideal(Å)
2	С	400	ATP	O4'-C1'	2.46	1.44	1.41
2	А	200	ATP	O4'-C1'	2.44	1.44	1.41
2	В	300	ATP	O4'-C1'	2.41	1.44	1.41

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	В	300	ATP	N3-C2-N1	-4.92	121.00	128.68

Continued on next page...



Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	А	200	ATP	N3-C2-N1	-4.88	121.05	128.68
2	С	400	ATP	N3-C2-N1	-4.85	121.11	128.68
2	А	200	ATP	PA-O3A-PB	-3.85	119.63	132.83
2	С	400	ATP	PB-O3B-PG	-3.70	120.14	132.83
2	В	300	ATP	PA-O3A-PB	-3.69	120.18	132.83
2	С	400	ATP	PA-O3A-PB	-3.68	120.21	132.83
2	А	200	ATP	PB-O3B-PG	-3.59	120.50	132.83
2	В	300	ATP	PB-O3B-PG	-3.50	120.81	132.83
2	С	400	ATP	O5'-PA-O1A	-3.09	97.00	109.07
2	В	300	ATP	O5'-PA-O1A	-2.88	97.83	109.07
2	А	200	ATP	O5'-PA-O1A	-2.78	98.21	109.07
2	В	300	ATP	O2A-PA-O1A	2.70	125.60	112.24
2	А	200	ATP	O2A-PA-O1A	2.65	125.34	112.24
2	А	200	ATP	O2B-PB-O1B	2.62	125.21	112.24
2	С	400	ATP	O2B-PB-O1B	2.62	125.19	112.24
2	С	400	ATP	O2A-PA-O1A	2.60	125.10	112.24
2	В	300	ATP	O2B-PB-O1B	2.57	124.96	112.24
2	В	300	ATP	O2A-PA-O5'	-2.28	97.17	107.75
2	А	200	ATP	O2A-PA-O5'	-2.18	97.61	107.75
2	С	400	ATP	C2'-C3'-C4'	-2.03	98.71	102.64

Continued from previous page...

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	В	300	ATP	PB-O3B-PG-O2G
2	В	300	ATP	PB-O3B-PG-O3G
2	С	400	ATP	C5'-O5'-PA-O1A
2	С	400	ATP	C5'-O5'-PA-O2A
2	С	400	ATP	PG-O3B-PB-O1B
2	А	200	ATP	PB-O3B-PG-O3G
2	С	400	ATP	C5'-O5'-PA-O3A
2	В	300	ATP	PB-O3A-PA-O1A

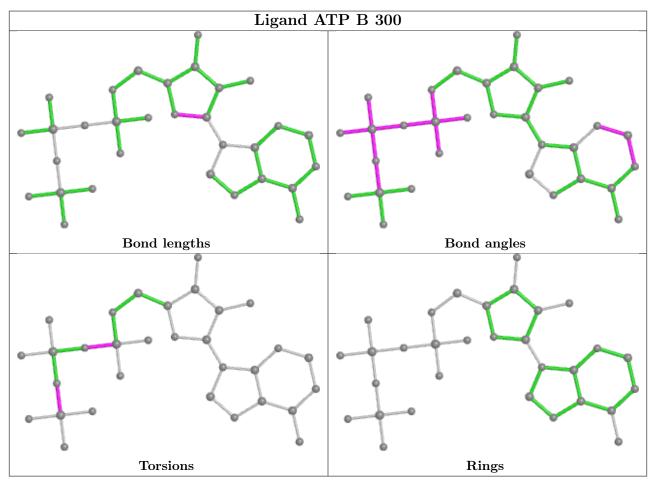
There are no ring outliers.

2 monomers are involved in 4 short contacts:

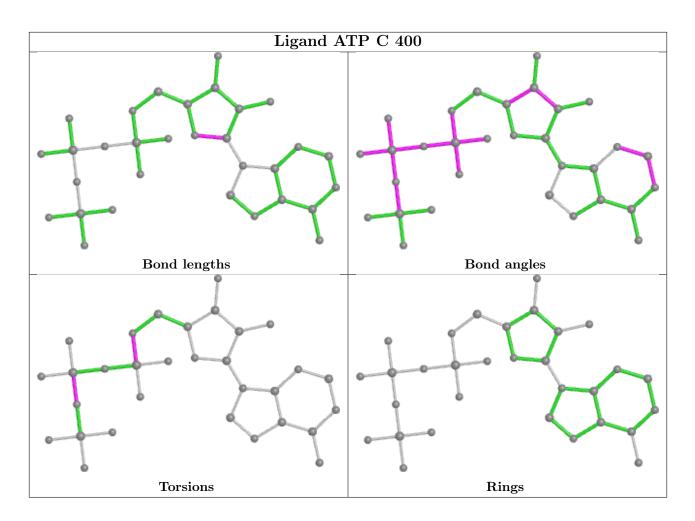
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	С	400	ATP	3	0
2	А	200	ATP	1	0



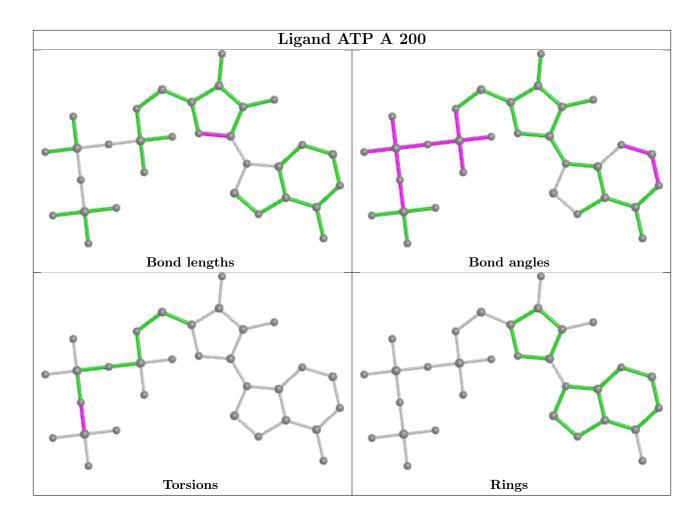
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 (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.

