



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 5, 2026 – 01:23 PM EDT

PDB ID : 1N5X / pdb_00001n5x
Title : Xanthine Dehydrogenase from Bovine Milk with Inhibitor TEI-6720 Bound
Authors : Okamoto, K.; Eger, B.T.; Nishino, T.; Kondo, S.; Pai, E.F.; Nishino, T.
Deposited on : 2002-11-07
Resolution : 2.80 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 ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

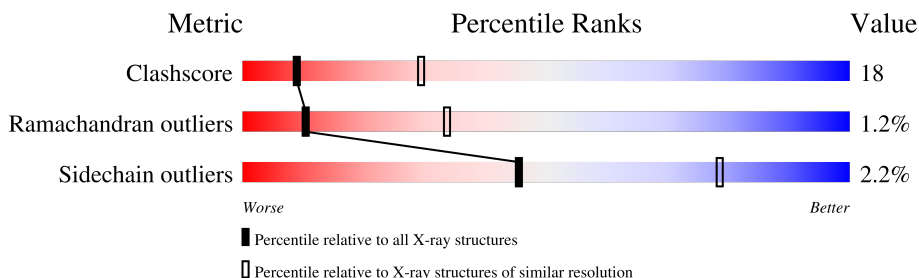
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	1331	
1	B	1331	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	ZW9	A	3007	-	-	X	-
5	ZW9	B	4007	-	-	X	-

2 Entry composition [i](#)

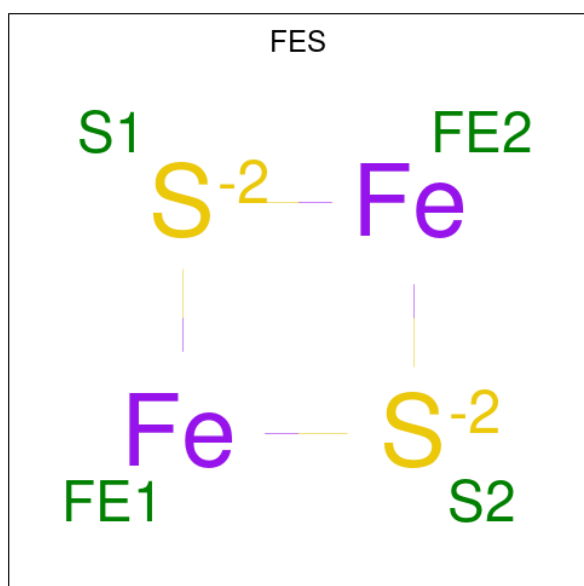
There are 5 unique types of molecules in this entry. The entry contains 20268 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Xanthine Dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1290	Total	C	N	O	S	0	0	0
			10023	6373	1718	1873	59			
1	B	1290	Total	C	N	O	S	0	0	0
			10023	6373	1718	1873	59			

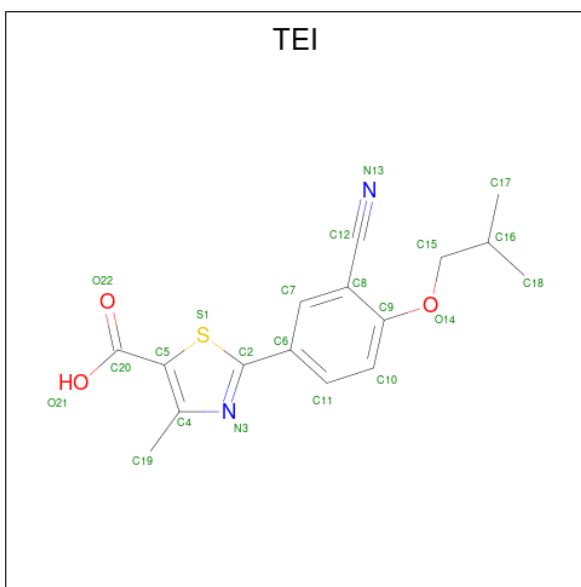
- Molecule 2 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			4	2	2		
2	A	1	Total	Fe	S	0	0
			4	2	2		
2	B	1	Total	Fe	S	0	0
			4	2	2		
2	B	1	Total	Fe	S	0	0
			4	2	2		

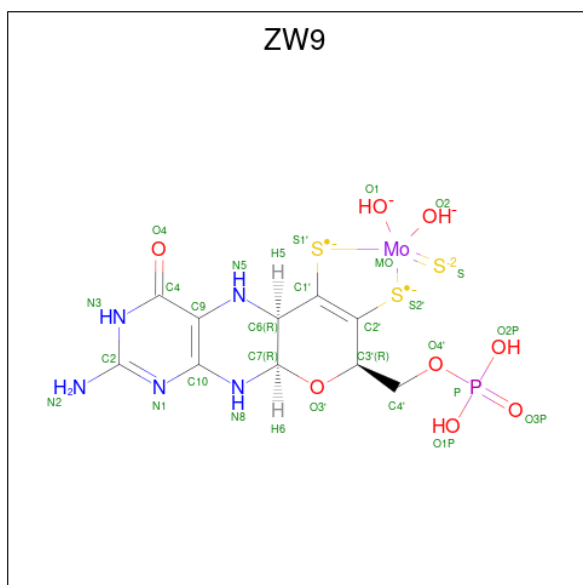
- # FAD

- Molecule 4 is 2-(3-CYANO-4-ISOBUTOXY-PHENYL)-4-METHYL-5-THIAZOLE-CARB OXYLIC ACID (CCD ID: TEI) (formula: $C_{16}H_{16}N_2O_3S$).

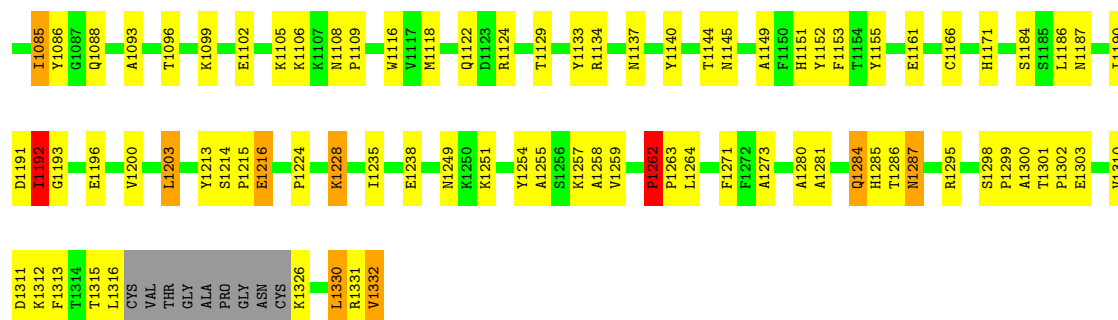


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			22	16	2	3	1		
4	B	1	Total	C	N	O	S	0	0
			22	16	2	3	1		

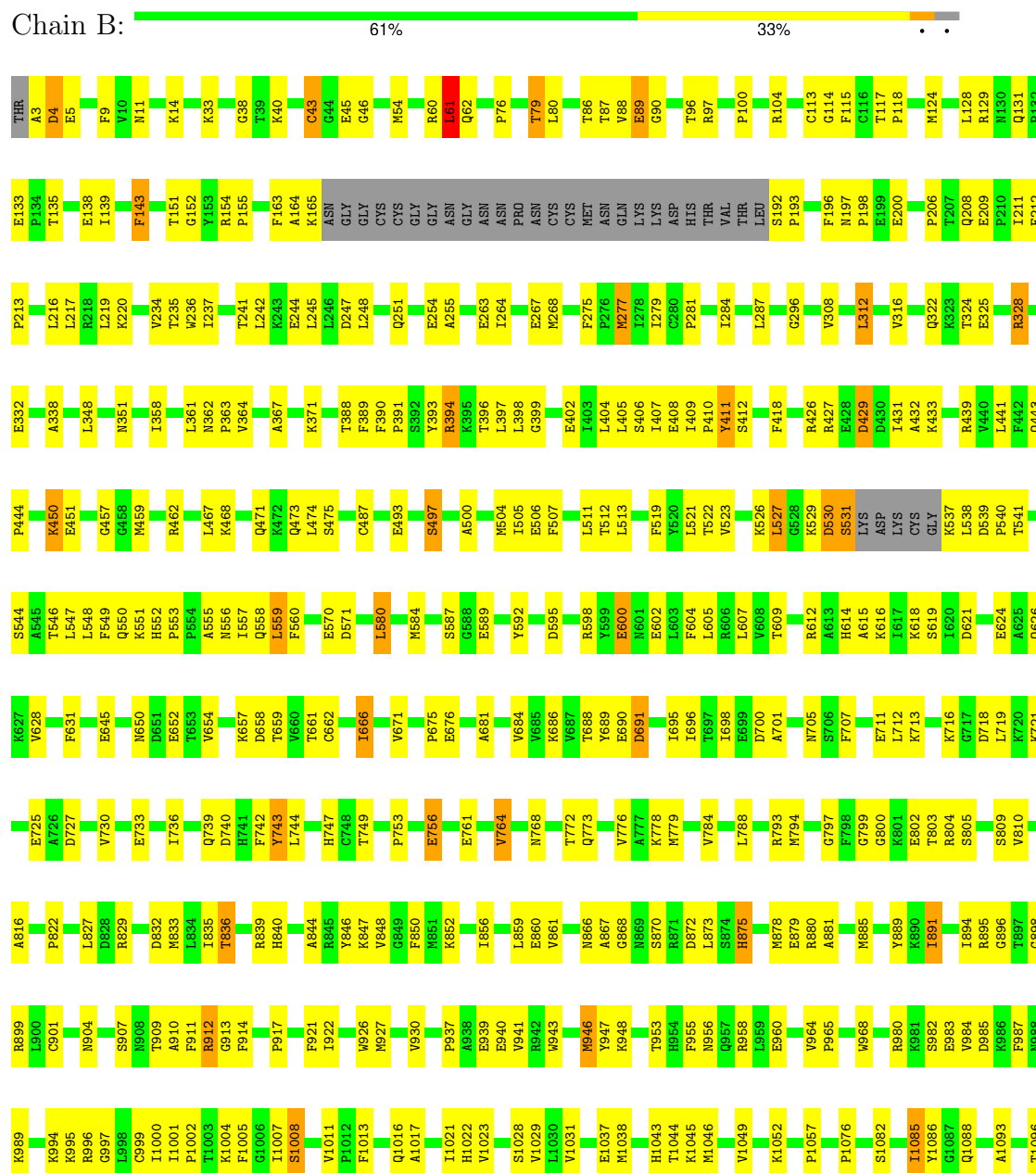
- Molecule 5 is Molybdopterin dihydroxy-thioxo-molybdenum (CCD ID: ZW9) (formula: $C_{10}H_{14}MoN_5O_8PS_3$).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
5	A	1	Total	C	Mo	N	O	P	S	0	0
			28	10	1	5	8	1	3		
5	B	1	Total	C	Mo	N	O	P	S	0	0
			28	10	1	5	8	1	3		



● Molecule 1: Xanthine Dehydrogenase





4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	168.27Å 124.66Å 147.32Å 90.00° 90.99° 90.00°	Depositor
Resolution (Å)	20.00 – 2.80	Depositor
% Data completeness (in resolution range)	93.7 (20.00-2.80)	Depositor
R_{merge}	0.11	Depositor
R_{sym}	0.11	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.244 , 0.275	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	20268	wwPDB-VP
Average B, all atoms (Å ²)	20.0	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: TEI, ZW9, FAD, FES

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	0.50	0/10242	1.00	48/13860 (0.3%)
1	B	0.50	0/10242	1.00	47/13860 (0.3%)
All	All	0.50	0/20484	1.00	95/27720 (0.3%)

There are no bond length outliers.

The worst 5 of 95 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	835	ILE	N-CA-C	10.90	122.03	111.67
1	A	835	ILE	N-CA-C	10.89	122.02	111.67
1	B	151	THR	N-CA-C	9.70	124.38	112.54
1	A	151	THR	N-CA-C	9.68	124.36	112.54
1	B	409	ILE	CA-C-N	8.34	128.35	119.76

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	10023	0	10026	368	5
1	B	10023	0	10026	370	4
2	A	8	0	0	2	0
2	B	8	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	53	0	29	2	0
3	B	53	0	29	2	0
4	A	22	0	15	0	0
4	B	22	0	15	0	0
5	A	28	0	0	10	0
5	B	28	0	0	10	0
All	All	20268	0	20140	747	5

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 747 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:1330:LEU:HD22	1:A:1331:ARG:N	1.62	1.14
1:B:1330:LEU:HD22	1:B:1331:ARG:N	1.62	1.13
1:A:537:LYS:HG3	1:A:538:LEU:H	1.18	1.07
1:A:1286:THR:HG22	1:A:1287:ASN:H	1.19	1.07
1:B:1286:THR:HG22	1:B:1287:ASN:H	1.19	1.04

All (5) 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:499:ASP:OD2	1:B:1326:LYS:O[1_545]	1.88	0.32
1:A:501:PRO:CA	1:B:1328:TRP:CB[1_545]	1.96	0.24
1:A:1213:TYR:CB	1:B:1332:VAL:CG1[1_545]	2.09	0.11
1:A:565:ASN:CB	1:A:565:ASN:CB[2_655]	2.17	0.03
1:A:501:PRO:CB	1:B:1328:TRP:CG[1_545]	2.19	0.01

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	1282/1331 (96%)	1173 (92%)	94 (7%)	15 (1%)	10	34
1	B	1282/1331 (96%)	1172 (91%)	95 (7%)	15 (1%)	10	34
All	All	2564/2662 (96%)	2345 (92%)	189 (7%)	30 (1%)	10	34

5 of 30 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4	ASP
1	A	530	ASP
1	A	1008	SER
1	A	1192	ILE
1	A	1287	ASN

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	1095/1127 (97%)	1070 (98%)	25 (2%)	44	78
1	B	1095/1127 (97%)	1071 (98%)	24 (2%)	45	78
All	All	2190/2254 (97%)	2141 (98%)	49 (2%)	45	78

5 of 49 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	312	LEU
1	B	666	ILE
1	B	328	ARG
1	B	546	THR
1	B	756	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 50 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	226	GLN

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Mol	Chain	Res	Type
1	B	626	GLN
1	B	1284	GLN
1	B	251	GLN
1	B	473	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

10 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
5	ZW9	B	4007	-	21,31,31	6.35	10 (47%)	22,52,52	4.29	13 (59%)
4	TEI	B	4006	-	23,23,23	4.16	8 (34%)	32,32,32	2.39	9 (28%)
2	FES	A	3002	1	0,4,4	-	-	-	-	-
5	ZW9	A	3007	-	21,31,31	6.35	10 (47%)	22,52,52	4.30	13 (59%)
3	FAD	A	3005	-	58,58,58	5.37	40 (68%)	85,89,89	2.38	29 (34%)
2	FES	A	3001	1	0,4,4	-	-	-	-	-
2	FES	B	4002	1	0,4,4	-	-	-	-	-
4	TEI	A	3006	-	23,23,23	4.16	8 (34%)	32,32,32	2.40	9 (28%)
2	FES	B	4001	1	0,4,4	-	-	-	-	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	FAD	B	4005	-	58,58,58	5.37	40 (68%)	85,89,89	2.39	30 (35%)

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
5	ZW9	B	4007	-	-	0/6/46/46	0/4/4/4
4	TEI	B	4006	-	-	0/15/15/15	0/2/2/2
5	ZW9	A	3007	-	-	0/6/46/46	0/4/4/4
2	FES	A	3002	1	-	-	0/1/1/1
3	FAD	A	3005	-	-	6/34/50/50	0/6/6/6
2	FES	A	3001	1	-	-	0/1/1/1
2	FES	B	4002	1	-	-	0/1/1/1
4	TEI	A	3006	-	-	0/15/15/15	0/2/2/2
2	FES	B	4001	1	-	-	0/1/1/1
3	FAD	B	4005	-	-	6/34/50/50	0/6/6/6

The worst 5 of 116 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	4005	FAD	P-O3P	-21.71	1.36	1.59
3	A	3005	FAD	P-O3P	-21.68	1.36	1.59
5	A	3007	ZW9	C7-C6	17.21	1.67	1.53
5	B	4007	ZW9	C7-C6	17.21	1.67	1.53
4	B	4006	TEI	C11-C6	15.43	1.62	1.39

The worst 5 of 103 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	3007	ZW9	C2'-C1'-S1'	12.32	127.14	120.15
5	B	4007	ZW9	C2'-C1'-S1'	12.26	127.11	120.15
4	A	3006	TEI	C2-N3-C4	8.72	115.41	111.20
4	B	4006	TEI	C2-N3-C4	8.65	115.37	111.20
5	A	3007	ZW9	C1'-C2'-S2'	-8.45	115.35	120.15

There are no chirality outliers.

5 of 12 torsion outliers are listed below:

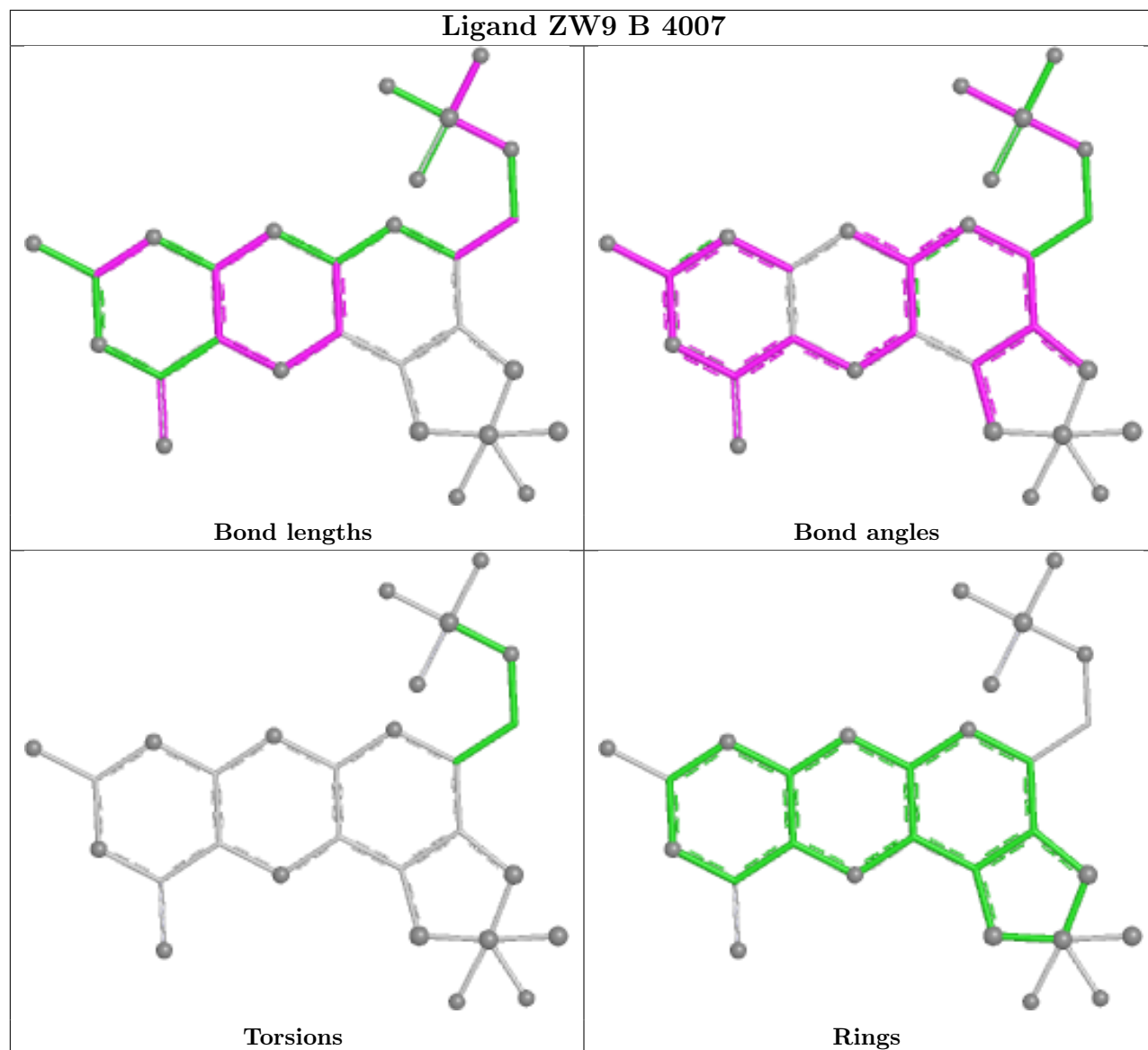
Mol	Chain	Res	Type	Atoms
3	A	3005	FAD	N10-C1'-C2'-O2'
3	A	3005	FAD	C2'-C3'-C4'-O4'
3	A	3005	FAD	O3'-C3'-C4'-O4'
3	A	3005	FAD	O3'-C3'-C4'-C5'
3	B	4005	FAD	N10-C1'-C2'-O2'

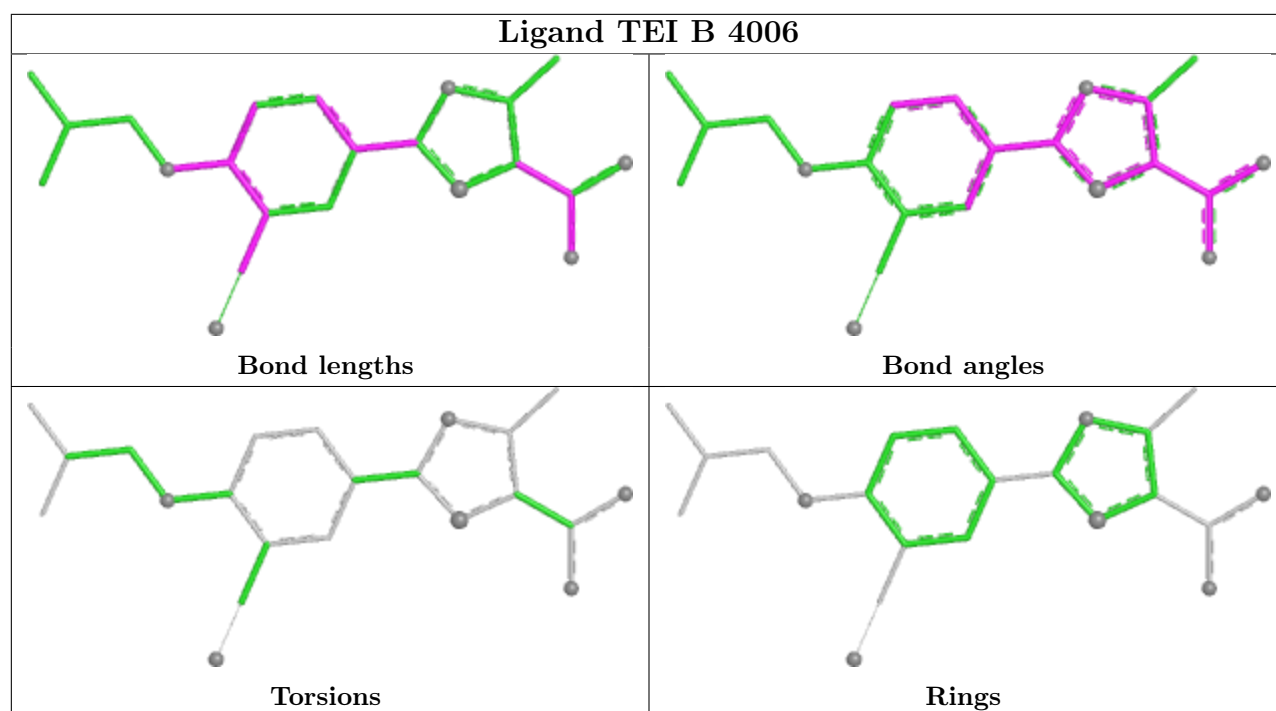
There are no ring outliers.

8 monomers are involved in 28 short contacts:

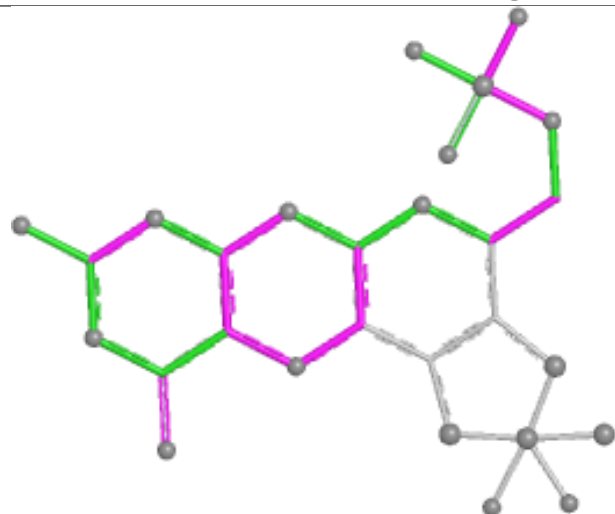
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	4007	ZW9	10	0
2	A	3002	FES	1	0
5	A	3007	ZW9	10	0
3	A	3005	FAD	2	0
2	A	3001	FES	1	0
2	B	4002	FES	1	0
2	B	4001	FES	1	0
3	B	4005	FAD	2	0

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 than 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.

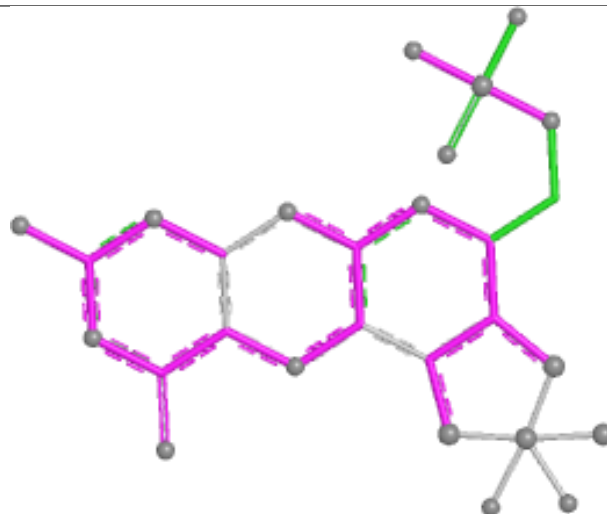




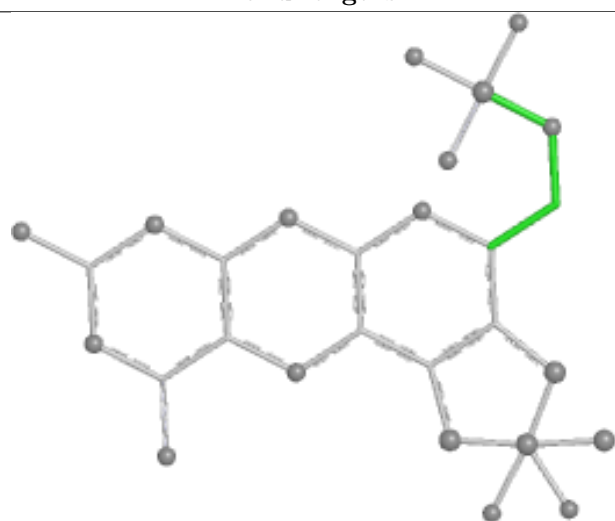
Ligand ZW9 A 3007



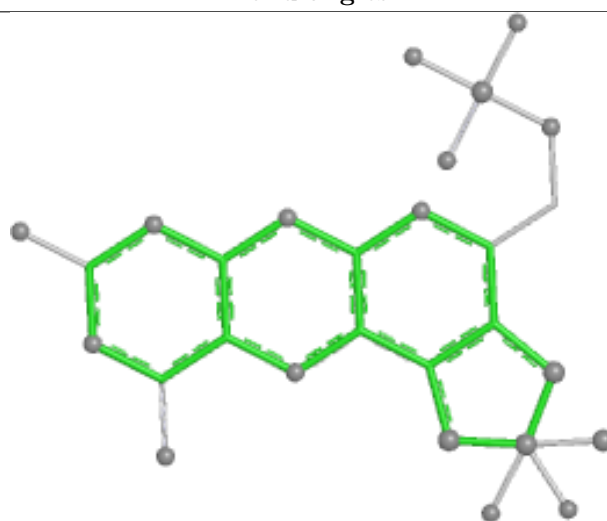
Bond lengths



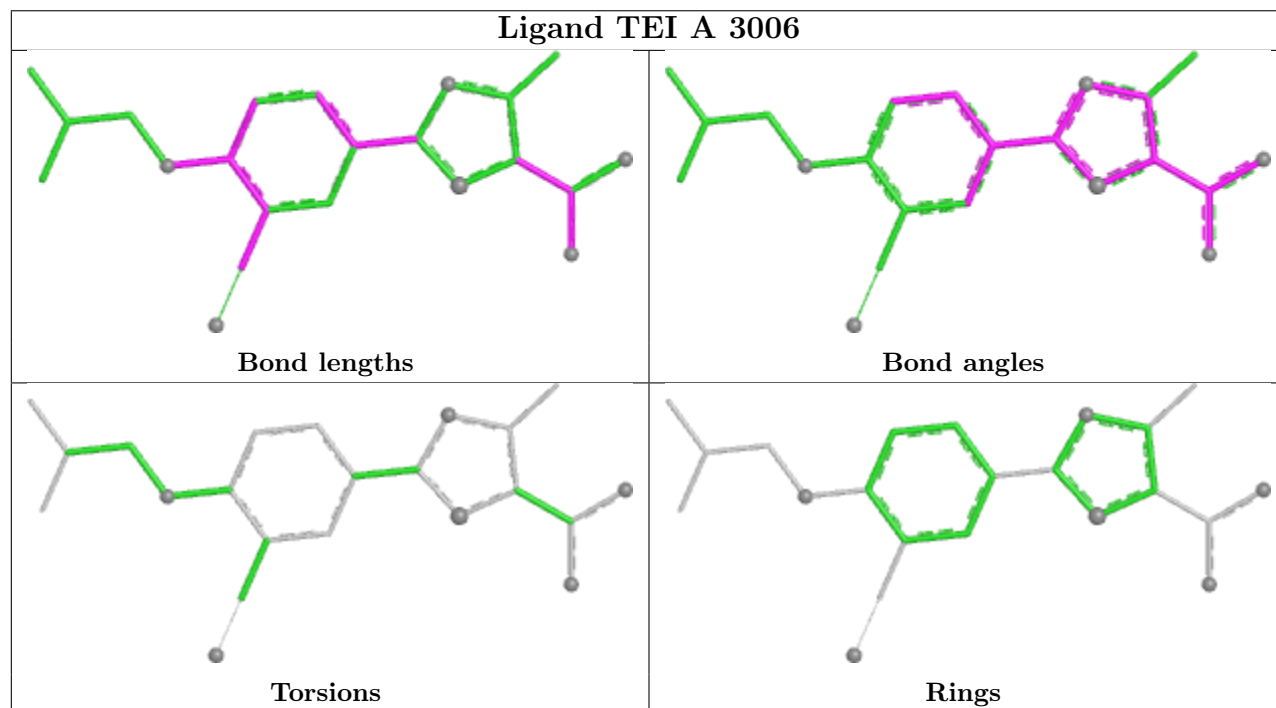
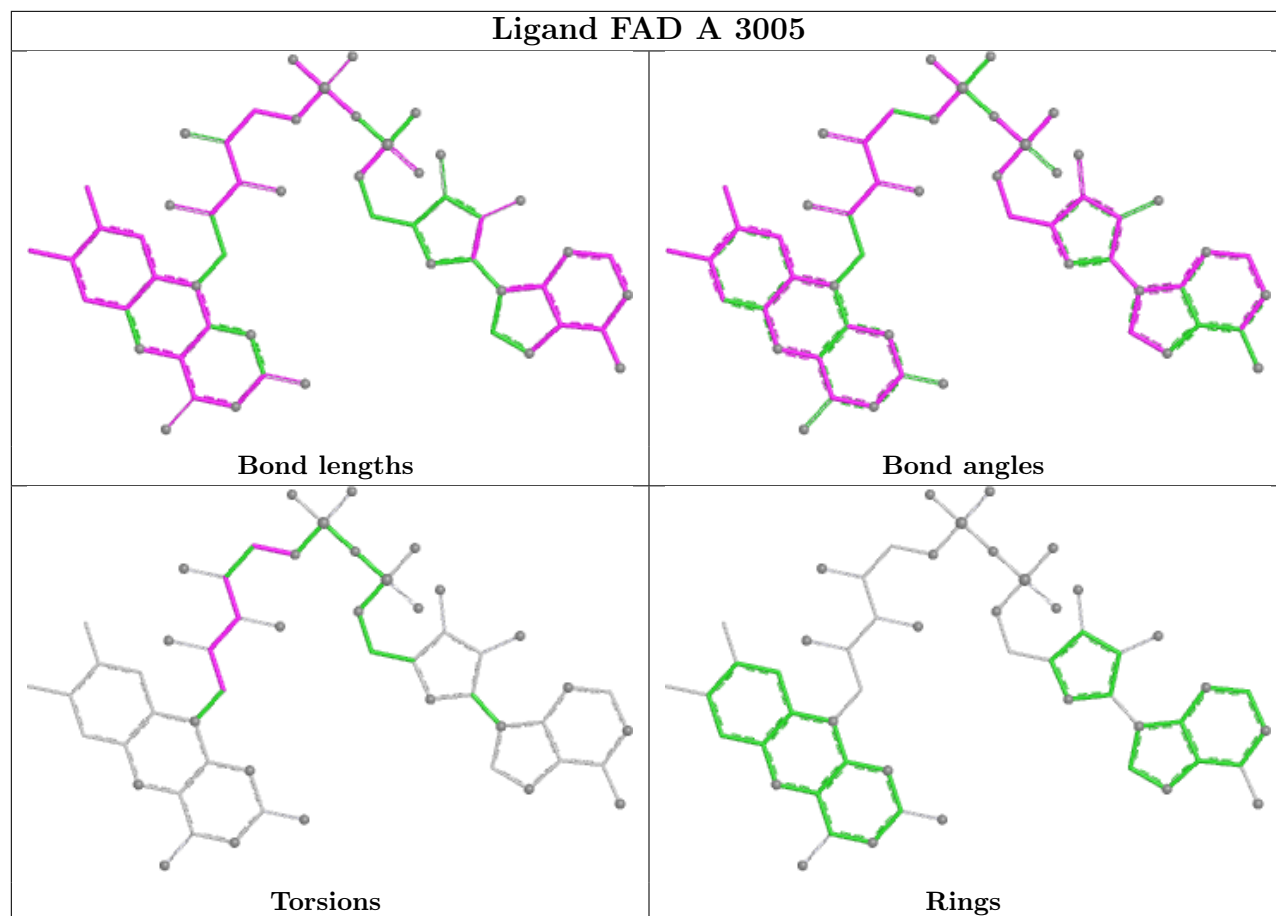
Bond angles

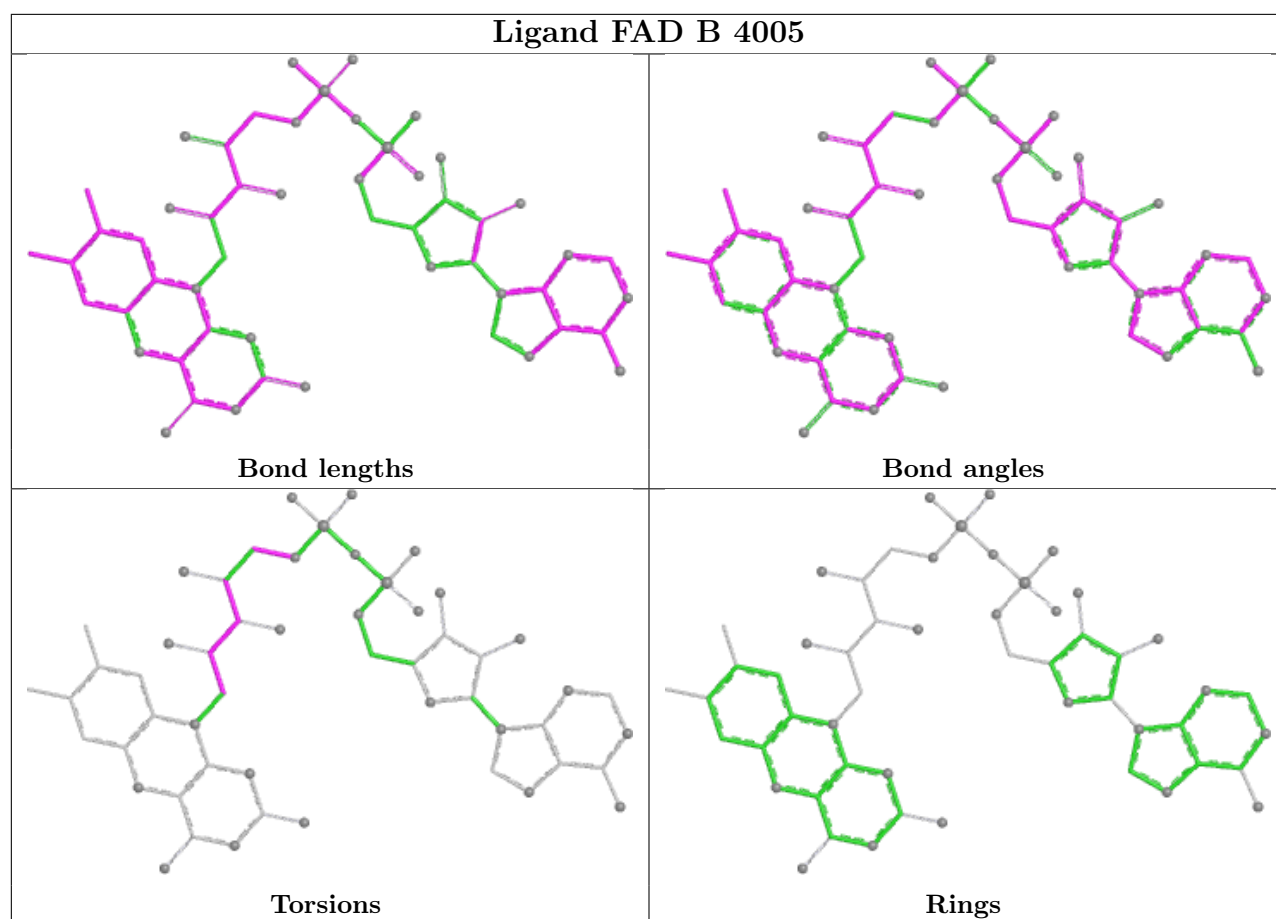


Torsions



Rings





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.