



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:02 AM UTC

PDB ID : 1BVQ / pdb_00001bvq
Title : THREE-DIMENSIONAL STRUCTURE OF 4-HYDROXYBENZOYL COA THIOESTERASE FROM PSEUDOMONAS SP. STRAIN CBS-3.
Authors : Holden, H.M.; Benning, M.M.; Dunaway-Mariano, D.
Deposited on : 1998-09-16
Resolution : 2.00 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 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
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)	:	2.49

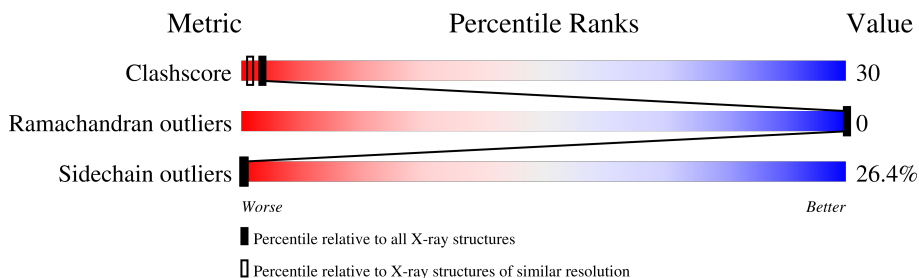
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)

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	141	 39% 38% 14% 7%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 1196 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.

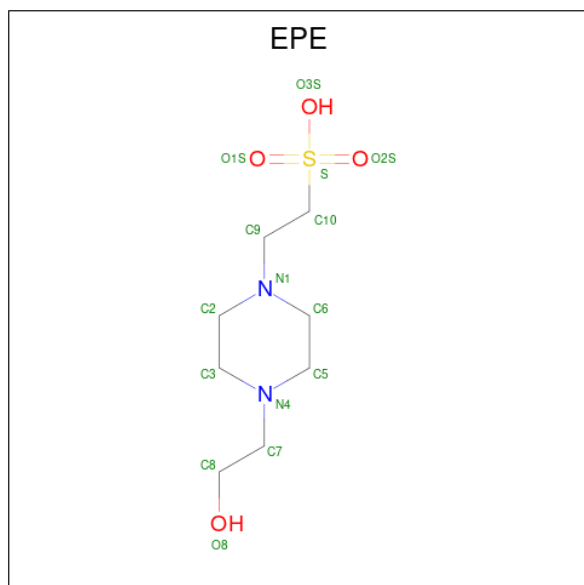
- Molecule 1 is a protein called PROTEIN (4-HYDROXYBENZOYL COA THIOESTERASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	139	Total	C	N	O	S	0	0	0
			1115	702	201	203	9			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	TYR	PHE	conflict	UNP P56653

- Molecule 2 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 3 is water.

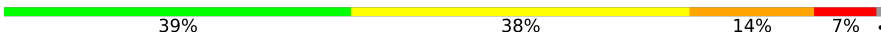
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	66	Total 66	O 66	0	0

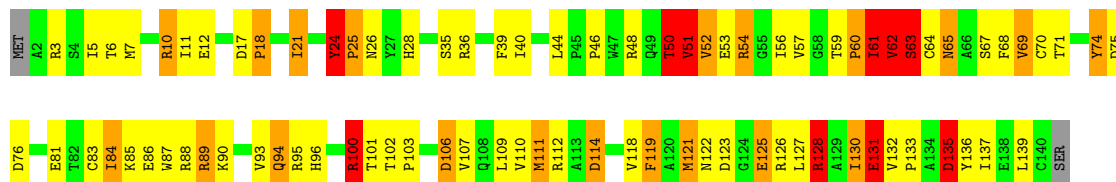
3 Residue-property plots [i](#)

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.

Note EDS was not executed.

• Molecule 1: PROTEIN (4-HYDROXYBENZOYL COA THIOESTERASE)

Chain A: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	49.60Å 56.00Å 93.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.00	Depositor
% Data completeness (in resolution range)	96.0 (30.00-2.00)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT 5E	Depositor
R, R_{free}	(Not available) , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1196	wwPDB-VP
Average B, all atoms (Å ²)	30.0	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: EPE

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	1.36	8/1141 (0.7%)	1.97	39/1550 (2.5%)

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	0	1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	54	ARG	C-N	-6.06	1.25	1.33
1	A	70	CYS	N-CA	5.69	1.53	1.45
1	A	52	VAL	C-N	-5.62	1.26	1.33
1	A	106	ASP	C-N	5.58	1.40	1.33
1	A	69	VAL	N-CA	-5.38	1.39	1.45
1	A	130	ILE	CA-C	-5.24	1.46	1.52
1	A	128	ARG	NE-CZ	5.18	1.38	1.33
1	A	12	GLU	CD-OE1	5.00	1.34	1.25

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	128	ARG	CD-NE-CZ	14.90	145.26	124.40
1	A	100	ARG	CD-NE-CZ	-12.14	107.41	124.40
1	A	128	ARG	NE-CZ-NH1	11.13	132.63	121.50
1	A	76	ASP	CA-CB-CG	9.14	121.74	112.60
1	A	128	ARG	NE-CZ-NH2	-8.65	111.42	119.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	119	PHE	CA-CB-CG	-8.51	105.29	113.80
1	A	61	ILE	CB-CA-C	7.95	122.10	111.21
1	A	50	THR	N-CA-CB	-7.22	99.22	110.06
1	A	135	ASP	N-CA-CB	6.89	121.62	110.40
1	A	114	ASP	CA-CB-CG	6.82	119.42	112.60
1	A	96	HIS	CA-CB-CG	6.79	120.59	113.80
1	A	60	PRO	CB-CA-C	-6.40	103.20	111.39
1	A	103	PRO	N-CA-C	-6.12	101.59	111.14
1	A	69	VAL	CB-CA-C	-6.10	105.69	111.06
1	A	76	ASP	N-CA-CB	6.09	118.94	110.17
1	A	63	SER	N-CA-CB	-6.06	102.17	111.56
1	A	61	ILE	N-CA-CB	-5.90	102.42	110.56
1	A	24	TYR	N-CA-C	5.84	121.48	113.77
1	A	84	ILE	N-CA-CB	5.75	117.56	111.00
1	A	83	CYS	N-CA-CB	5.71	120.58	111.56
1	A	39	PHE	CA-CB-CG	-5.68	108.12	113.80
1	A	25	PRO	N-CA-CB	5.57	109.10	103.25
1	A	18	PRO	N-CA-CB	5.56	109.04	103.48
1	A	10	ARG	N-CA-CB	5.53	119.19	110.23
1	A	119	PHE	N-CA-C	-5.53	101.47	110.20
1	A	35	SER	N-CA-CB	-5.38	102.21	110.01
1	A	51	VAL	CA-C-N	-5.32	117.50	122.66
1	A	51	VAL	C-N-CA	-5.32	117.50	122.66
1	A	26	ASN	N-CA-CB	5.32	118.52	110.28
1	A	46	PRO	N-CA-CB	5.29	108.36	103.39
1	A	131	GLU	N-CA-CB	-5.24	101.91	109.83
1	A	48	ARG	NE-CZ-NH2	5.14	123.83	119.20
1	A	74	TYR	CB-CA-C	5.12	118.70	109.64
1	A	39	PHE	N-CA-CB	-5.11	102.59	110.16
1	A	102	THR	N-CA-CB	5.09	116.16	109.55
1	A	62	VAL	CA-C-O	5.07	123.55	119.29
1	A	28	HIS	CA-CB-CG	-5.06	108.74	113.80
1	A	63	SER	CB-CA-C	-5.05	101.23	110.62
1	A	130	ILE	O-C-N	5.01	128.68	123.02

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	128	ARG	Sidechain

5.2 Too-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 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	1115	0	1092	67	0
2	A	15	0	17	3	0
3	A	66	0	0	7	0
All	All	1196	0	1109	67	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 30.

All (67) 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:89:ARG:HD3	1:A:89:ARG:H	1.18	1.04
1:A:95:ARG:HH12	2:A:208:EPE:H52	1.28	0.97
1:A:68:PHE:CE2	1:A:111:MET:HE3	2.08	0.89
1:A:61:ILE:HD13	1:A:62:VAL:N	1.94	0.81
1:A:95:ARG:NH1	2:A:208:EPE:H52	1.95	0.81
1:A:64:CYS:HB2	3:A:197:HOH:O	1.83	0.78
1:A:50:THR:HG23	1:A:56:ILE:HB	1.68	0.76
1:A:50:THR:HG22	1:A:56:ILE:O	1.87	0.74
1:A:130:ILE:HG13	1:A:131:GLU:O	1.89	0.73
1:A:125:GLU:H	1:A:125:GLU:CD	1.98	0.72
1:A:87:TRP:CZ3	1:A:132:VAL:HG11	2.24	0.71
1:A:61:ILE:HD11	1:A:63:SER:O	1.91	0.69
1:A:123:ASP:HB2	1:A:128:ARG:HG2	1.76	0.68
1:A:44:LEU:HD12	1:A:139:LEU:HD13	1.76	0.67
1:A:89:ARG:HD3	1:A:89:ARG:N	2.01	0.67
1:A:111:MET:HG2	1:A:112:ARG:N	2.09	0.66
1:A:89:ARG:H	1:A:89:ARG:CD	2.04	0.65
1:A:95:ARG:HH12	2:A:208:EPE:C5	2.07	0.65
1:A:44:LEU:HD12	1:A:139:LEU:CD1	2.26	0.65
1:A:87:TRP:HZ3	1:A:132:VAL:HG11	1.61	0.64
1:A:86:GLU:HG2	1:A:93:VAL:CG2	2.29	0.63
1:A:127:LEU:HD12	1:A:127:LEU:N	2.15	0.61
1:A:24:TYR:CG	1:A:25:PRO:HD3	2.36	0.59
1:A:24:TYR:CD2	1:A:25:PRO:HD3	2.38	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:THR:CG2	1:A:56:ILE:HB	2.35	0.57
1:A:123:ASP:HB2	1:A:128:ARG:CG	2.35	0.57
1:A:84:ILE:HD13	1:A:94:GLN:HG3	1.87	0.56
1:A:59:THR:HA	1:A:118:VAL:O	2.06	0.55
1:A:123:ASP:CB	1:A:128:ARG:HG2	2.37	0.54
1:A:61:ILE:HD13	1:A:61:ILE:C	2.32	0.54
1:A:135:ASP:O	1:A:139:LEU:HG	2.10	0.52
1:A:17:ASP:HB2	1:A:18:PRO:HD2	1.94	0.50
1:A:86:GLU:HG2	1:A:93:VAL:HB	1.94	0.49
1:A:95:ARG:NH1	3:A:181:HOH:O	2.45	0.48
1:A:100:ARG:HB3	1:A:110:VAL:CG2	2.43	0.48
1:A:119:PHE:CE2	1:A:132:VAL:HG13	2.48	0.48
1:A:100:ARG:HB3	1:A:110:VAL:HG23	1.95	0.48
1:A:90:LYS:O	1:A:118:VAL:HG12	2.14	0.48
1:A:136:TYR:N	1:A:136:TYR:CD1	2.79	0.48
1:A:135:ASP:CB	3:A:178:HOH:O	2.61	0.47
1:A:60:PRO:HD2	1:A:118:VAL:O	2.14	0.47
1:A:44:LEU:CD1	1:A:139:LEU:HD12	2.45	0.47
1:A:74:TYR:O	1:A:75:ASP:HB2	2.15	0.46
1:A:86:GLU:CD	1:A:88:ARG:HE	2.23	0.46
1:A:133:PRO:HG2	1:A:136:TYR:CE2	2.51	0.45
1:A:61:ILE:HD13	1:A:62:VAL:H	1.75	0.45
1:A:63:SER:O	1:A:64:CYS:HB3	2.17	0.45
1:A:127:LEU:HD12	1:A:127:LEU:H	1.81	0.45
1:A:36:ARG:NE	3:A:148:HOH:O	2.32	0.45
1:A:86:GLU:CG	1:A:93:VAL:CG2	2.94	0.45
1:A:54:ARG:NH2	1:A:135:ASP:OD1	2.50	0.45
1:A:10:ARG:HG2	1:A:11:ILE:N	2.32	0.44
1:A:101:THR:HA	1:A:106:ASP:O	2.17	0.44
1:A:65:ASN:ND2	3:A:204:HOH:O	2.51	0.44
1:A:61:ILE:HD13	1:A:63:SER:H	1.83	0.44
1:A:86:GLU:HG2	1:A:93:VAL:HG21	1.98	0.43
1:A:123:ASP:HB2	1:A:128:ARG:CD	2.47	0.43
1:A:132:VAL:HG21	3:A:175:HOH:O	2.19	0.43
1:A:44:LEU:HD12	1:A:139:LEU:HD12	2.00	0.42
1:A:86:GLU:HG2	1:A:93:VAL:CB	2.49	0.42
1:A:121:MET:HE3	1:A:121:MET:HB3	1.85	0.42
1:A:17:ASP:HB2	1:A:18:PRO:CD	2.50	0.41
1:A:6:THR:HG22	1:A:7:MET:N	2.35	0.41
1:A:51:VAL:HG23	1:A:56:ILE:O	2.21	0.41
1:A:64:CYS:HA	1:A:114:ASP:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:ASP:HB3	3:A:178:HOH:O	2.21	0.40
1:A:21:ILE:HG21	1:A:71:THR:HB	2.03	0.40

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	137/141 (97%)	132 (96%)	5 (4%)	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	121/123 (98%)	89 (74%)	32 (26%)	0	0

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ARG
1	A	5	ILE
1	A	21	ILE
1	A	24	TYR

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Mol	Chain	Res	Type
1	A	40	ILE
1	A	50	THR
1	A	51	VAL
1	A	52	VAL
1	A	53	GLU
1	A	57	VAL
1	A	61	ILE
1	A	62	VAL
1	A	63	SER
1	A	65	ASN
1	A	67	SER
1	A	69	VAL
1	A	81	GLU
1	A	85	LYS
1	A	89	ARG
1	A	94	GLN
1	A	100	ARG
1	A	107	VAL
1	A	109	LEU
1	A	111	MET
1	A	121	MET
1	A	122	ASN
1	A	125	GLU
1	A	126	ARG
1	A	128	ARG
1	A	131	GLU
1	A	135	ASP
1	A	137	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	94	GLN
1	A	122	ASN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	EPE	A	208	-	15,15,15	0.67	0	19,20,20	1.74	3 (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	EPE	A	208	-	-	9/9/19/19	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	208	EPE	O3S-S-O2S	-4.88	99.18	111.40
2	A	208	EPE	O2S-S-O1S	3.65	125.70	113.82
2	A	208	EPE	C3-C2-N1	-2.87	104.87	110.65

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	208	EPE	C10-C9-N1-C6
2	A	208	EPE	C9-C10-S-O3S
2	A	208	EPE	C8-C7-N4-C3
2	A	208	EPE	N4-C7-C8-O8
2	A	208	EPE	C10-C9-N1-C2
2	A	208	EPE	C8-C7-N4-C5
2	A	208	EPE	C9-C10-S-O1S
2	A	208	EPE	C9-C10-S-O2S
2	A	208	EPE	S-C10-C9-N1

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	208	EPE	3	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

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.