



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 02:11 am GMT

PDB ID : 1ATJ  
Title : RECOMBINANT HORSERADISH PEROXIDASE C1A  
Authors : Gajhede, M.; Schuller, D.J.; Henriksen, A.; Smith, A.T.; Poulos, T.L.  
Deposited on : 1997-08-14  
Resolution : 2.15 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](mailto: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.

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

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

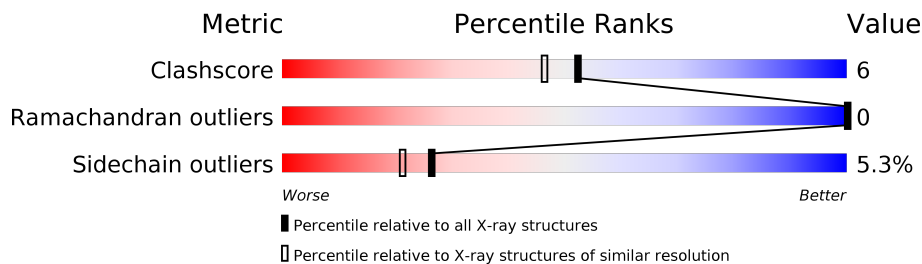
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.15 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	112137	1278 (2.16-2.16)
Ramachandran outliers	110173	1256 (2.16-2.16)
Sidechain outliers	110143	1255 (2.16-2.16)

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  $\geq 3$ , 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	306	
1	B	306	
1	C	306	
1	D	306	
1	E	306	
1	F	306	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 18528 atoms, of which 3156 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.

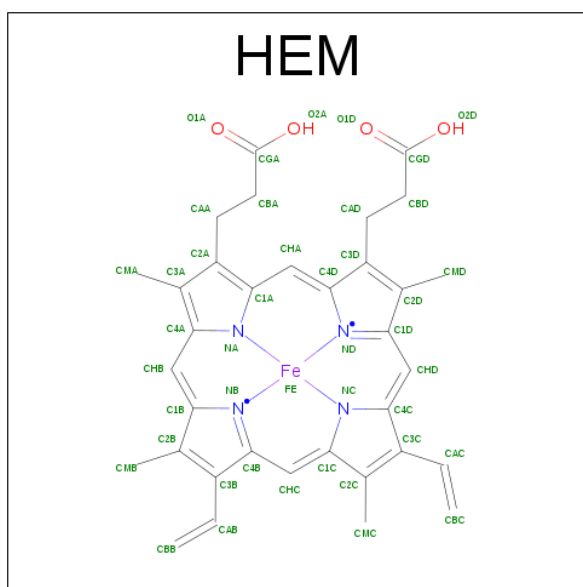
- Molecule 1 is a protein called PEROXIDASE C1A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	306	Total	C	H	N	O	S	0	0	0
			2895	1481	526	421	455	12			
1	B	306	Total	C	H	N	O	S	0	0	0
			2895	1481	526	421	455	12			
1	C	306	Total	C	H	N	O	S	0	0	0
			2895	1481	526	421	455	12			
1	D	306	Total	C	H	N	O	S	0	0	0
			2895	1481	526	421	455	12			
1	E	306	Total	C	H	N	O	S	0	0	0
			2895	1481	526	421	455	12			
1	F	306	Total	C	H	N	O	S	0	0	0
			2895	1481	526	421	455	12			

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	2	Total	Ca	0	0
			2	2		
2	E	2	Total	Ca	0	0
			2	2		
2	B	2	Total	Ca	0	0
			2	2		
2	C	2	Total	Ca	0	0
			2	2		
2	A	2	Total	Ca	0	0
			2	2		
2	F	2	Total	Ca	0	0
			2	2		

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: C<sub>34</sub>H<sub>32</sub>FeN<sub>4</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	
			43	34	1	4	4	
3	B	1	Total	C	Fe	N	O	
			43	34	1	4	4	
3	C	1	Total	C	Fe	N	O	
			43	34	1	4	4	
3	D	1	Total	C	Fe	N	O	
			43	34	1	4	4	
3	E	1	Total	C	Fe	N	O	
			43	34	1	4	4	
3	F	1	Total	C	Fe	N	O	
			43	34	1	4	4	

- Molecule 4 is water.

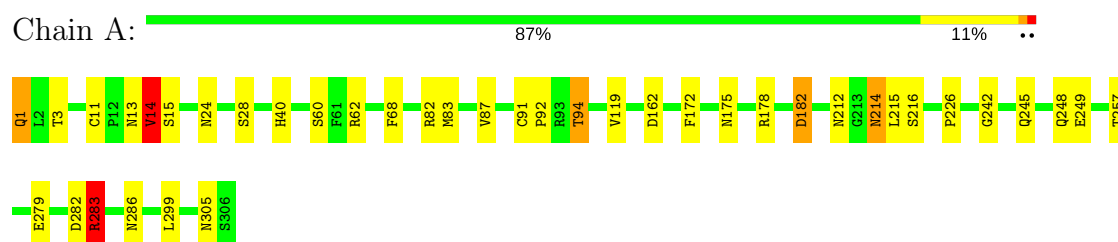
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	147	Total	O		
			147	147	0	0
4	B	150	Total	O		
			150	150	0	0
4	C	150	Total	O		
			150	150	0	0
4	D	148	Total	O		
			148	148	0	0
4	E	146	Total	O		
			146	146	0	0
4	F	147	Total	O		
			147	147	0	0

### 3 Residue-property plots

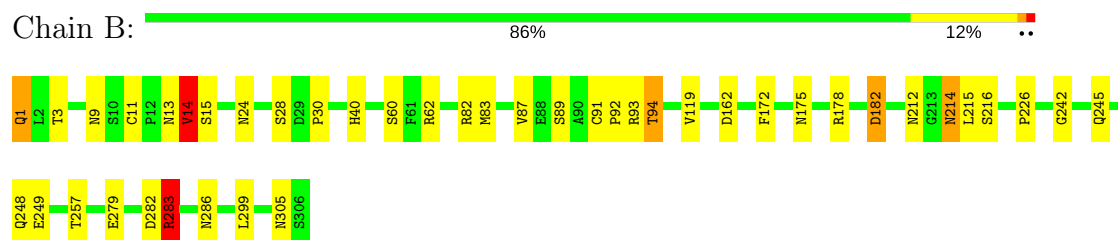
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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.

Note EDS was not executed.

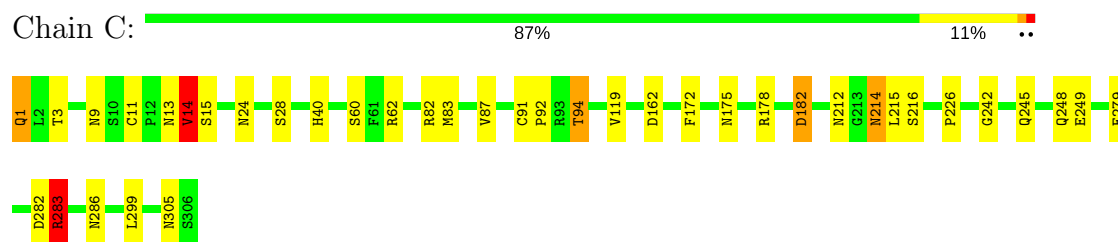
#### • Molecule 1: PEROXIDASE C1A



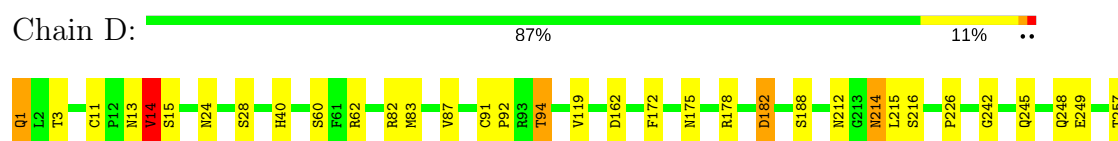
#### • Molecule 1: PEROXIDASE C1A



#### • Molecule 1: PEROXIDASE C1A

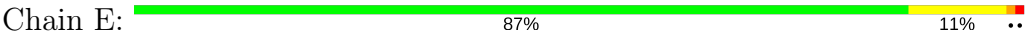


#### • Molecule 1: PEROXIDASE C1A

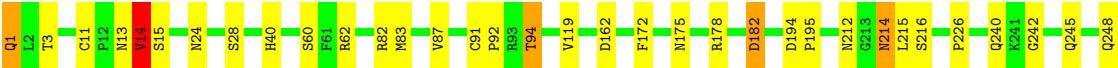
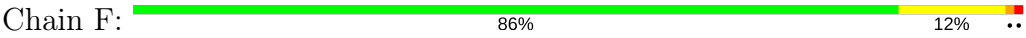




● Molecule 1: PEROXIDASE C1A



● Molecule 1: PEROXIDASE C1A



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 31 1 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	159.17Å 159.17Å 114.27Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	100.00 – 2.15	Depositor
% Data completeness (in resolution range)	83.7 (100.00-2.15)	Depositor
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, $R_{free}$	0.222 , 0.242	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	18528	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	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: HEM, CA

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	0.74	3/2415 (0.1%)	0.94	7/3286 (0.2%)
1	B	0.74	3/2415 (0.1%)	0.94	7/3286 (0.2%)
1	C	0.74	3/2415 (0.1%)	0.94	7/3286 (0.2%)
1	D	0.74	3/2415 (0.1%)	0.94	7/3286 (0.2%)
1	E	0.74	3/2415 (0.1%)	0.94	7/3286 (0.2%)
1	F	0.74	3/2415 (0.1%)	0.94	7/3286 (0.2%)
All	All	0.74	18/14490 (0.1%)	0.94	42/19716 (0.2%)

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	182	ASP	CB-CG	11.30	1.75	1.51
1	B	182	ASP	CB-CG	11.29	1.75	1.51
1	A	182	ASP	CB-CG	11.29	1.75	1.51
1	C	182	ASP	CB-CG	11.29	1.75	1.51
1	D	182	ASP	CB-CG	11.28	1.75	1.51

The worst 5 of 42 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	182	ASP	CB-CG-OD2	19.79	136.11	118.30
1	E	182	ASP	CB-CG-OD2	19.77	136.09	118.30
1	C	182	ASP	CB-CG-OD2	19.75	136.07	118.30
1	D	182	ASP	CB-CG-OD2	19.73	136.06	118.30
1	A	182	ASP	CB-CG-OD2	19.72	136.05	118.30

There are no chirality outliers.

There are no planarity outliers.



## 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	2369	526	2321	31	0
1	B	2369	526	2321	29	3
1	C	2369	526	2321	30	0
1	D	2369	526	2321	34	0
1	E	2369	526	2321	28	2
1	F	2369	526	2321	30	1
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
3	A	43	0	30	1	0
3	B	43	0	30	1	0
3	C	43	0	30	1	0
3	D	43	0	30	1	0
3	E	43	0	30	1	0
3	F	43	0	30	1	0
4	A	147	0	0	8	0
4	B	150	0	0	10	0
4	C	150	0	0	10	3
4	D	148	0	0	13	1
4	E	146	0	0	9	0
4	F	147	0	0	8	3
All	All	15372	3156	14106	179	7

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 6.

The worst 5 of 179 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:C:182:ASP:CB	1:C:182:ASP:CG	1.75	1.55
1:D:182:ASP:CB	1:D:182:ASP:CG	1.75	1.54
1:B:182:ASP:CG	1:B:182:ASP:CB	1.75	1.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:182:ASP:CG	1:E:182:ASP:CB	1.75	1.53
1:A:182:ASP:CG	1:A:182:ASP:CB	1.75	1.51

The worst 5 of 7 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:B:257:THR:CG2	4:C:457:HOH:O[2_565]	1.05	1.15
1:B:89:SER:HG	1:B:93:ARG:HH21[6_566]	0.82	0.78
1:E:257:THR:CB	4:F:456:HOH:O[4_555]	1.77	0.43
1:B:257:THR:CB	4:C:457:HOH:O[2_565]	1.89	0.31
4:C:484:HOH:O	4:F:480:HOH:O[4_555]	2.03	0.17

## 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	304/306 (99%)	297 (98%)	7 (2%)	0	100	100
1	B	304/306 (99%)	297 (98%)	7 (2%)	0	100	100
1	C	304/306 (99%)	297 (98%)	7 (2%)	0	100	100
1	D	304/306 (99%)	297 (98%)	7 (2%)	0	100	100
1	E	304/306 (99%)	297 (98%)	7 (2%)	0	100	100
1	F	304/306 (99%)	297 (98%)	7 (2%)	0	100	100
All	All	1824/1836 (99%)	1782 (98%)	42 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	266/266 (100%)	252 (95%)	14 (5%)	26	21
1	B	266/266 (100%)	252 (95%)	14 (5%)	26	21
1	C	266/266 (100%)	252 (95%)	14 (5%)	26	21
1	D	266/266 (100%)	252 (95%)	14 (5%)	26	21
1	E	266/266 (100%)	252 (95%)	14 (5%)	26	21
1	F	266/266 (100%)	252 (95%)	14 (5%)	26	21
All	All	1596/1596 (100%)	1512 (95%)	84 (5%)	26	21

5 of 84 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	172	PHE
1	D	24	ASN
1	F	60	SER
1	C	214	ASN
1	D	1	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 79 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	255	ASN
1	D	214	ASN
1	F	245	GLN
1	C	271	GLN
1	D	16	ASN

### 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 ⓘ

Of 18 ligands modelled in this entry, 12 are monoatomic - leaving 6 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$
3	HEM	A	350	1	28,50,50	2.00	9 (32%)	17,82,82	2.29	5 (29%)
3	HEM	B	350	1	28,50,50	2.00	9 (32%)	17,82,82	2.30	5 (29%)
3	HEM	C	350	1	28,50,50	1.99	9 (32%)	17,82,82	2.29	5 (29%)
3	HEM	D	350	1	28,50,50	2.00	9 (32%)	17,82,82	2.29	5 (29%)
3	HEM	E	350	1	28,50,50	1.99	9 (32%)	17,82,82	2.28	5 (29%)
3	HEM	F	350	1	28,50,50	1.99	9 (32%)	17,82,82	2.28	5 (29%)

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
3	HEM	A	350	1	-	0/6/54/54	0/0/8/8
3	HEM	B	350	1	-	0/6/54/54	0/0/8/8
3	HEM	C	350	1	-	0/6/54/54	0/0/8/8
3	HEM	D	350	1	-	0/6/54/54	0/0/8/8
3	HEM	E	350	1	-	0/6/54/54	0/0/8/8
3	HEM	F	350	1	-	0/6/54/54	0/0/8/8

The worst 5 of 54 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	350	HEM	C3C-CAC	-3.18	1.41	1.47
3	C	350	HEM	C3C-CAC	-3.18	1.41	1.47
3	B	350	HEM	C3C-CAC	-3.17	1.41	1.47
3	F	350	HEM	C3C-CAC	-3.16	1.41	1.47
3	D	350	HEM	C3C-CAC	-3.16	1.41	1.47

The worst 5 of 30 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	350	HEM	CMA-C3A-C4A	-4.66	121.31	128.46
3	C	350	HEM	CMA-C3A-C4A	-4.63	121.35	128.46
3	A	350	HEM	CMA-C3A-C4A	-4.62	121.36	128.46
3	F	350	HEM	CMA-C3A-C4A	-4.62	121.37	128.46
3	D	350	HEM	CMA-C3A-C4A	-4.60	121.39	128.46

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	350	HEM	1	0
3	B	350	HEM	1	0
3	C	350	HEM	1	0
3	D	350	HEM	1	0
3	E	350	HEM	1	0
3	F	350	HEM	1	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.