



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 06:00 AM GMT

PDB ID : 3A17  
Title : Crystal Structure of Aldoxime Dehydratase (OxdRE) in Complex with Butyraldoxime (Co-crystal)  
Authors : Sawai, H.; Sugimoto, H.; Kato, Y.; Asano, Y.; Shiro, Y.; Aono, S.  
Deposited on : 2009-03-26  
Resolution : 2.50 Å(reported)

This is a full wwPDB validation 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/ValidationPDFNotes.html>

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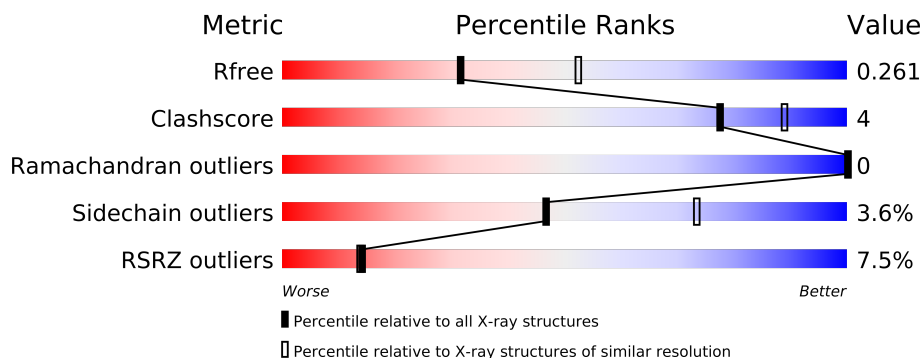
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.50 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))
$R_{free}$	66092	2784 (2.50-2.50)
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)
RSRZ outliers	66119	2785 (2.50-2.50)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	373	
1	B	373	
1	C	373	
1	D	373	
1	E	373	
1	F	373	
1	G	373	
1	H	373	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 23656 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Aldoxime dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	350	Total	C	N	O	S	0	0	0
			2790	1757	492	529	12			
1	B	359	Total	C	N	O	S	0	0	0
			2853	1795	506	540	12			
1	C	352	Total	C	N	O	S	0	1	0
			2812	1770	497	533	12			
1	D	360	Total	C	N	O	S	0	1	0
			2866	1803	510	541	12			
1	E	351	Total	C	N	O	S	0	1	0
			2803	1765	496	530	12			
1	F	359	Total	C	N	O	S	0	0	0
			2853	1795	506	540	12			
1	G	353	Total	C	N	O	S	0	0	0
			2815	1771	497	535	12			
1	H	359	Total	C	N	O	S	0	0	0
			2853	1795	506	540	12			

There are 160 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	INITIATING METHIONINE	UNP Q76K71
A	-18	GLY	-	EXPRESSION TAG	UNP Q76K71
A	-17	SER	-	EXPRESSION TAG	UNP Q76K71
A	-16	SER	-	EXPRESSION TAG	UNP Q76K71
A	-15	HIS	-	EXPRESSION TAG	UNP Q76K71
A	-14	HIS	-	EXPRESSION TAG	UNP Q76K71
A	-13	HIS	-	EXPRESSION TAG	UNP Q76K71
A	-12	HIS	-	EXPRESSION TAG	UNP Q76K71
A	-11	HIS	-	EXPRESSION TAG	UNP Q76K71
A	-10	HIS	-	EXPRESSION TAG	UNP Q76K71
A	-9	SER	-	EXPRESSION TAG	UNP Q76K71
A	-8	SER	-	EXPRESSION TAG	UNP Q76K71
A	-7	GLY	-	EXPRESSION TAG	UNP Q76K71

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	LEU	-	EXPRESSION TAG	UNP Q76K71
A	-5	VAL	-	EXPRESSION TAG	UNP Q76K71
A	-4	PRO	-	EXPRESSION TAG	UNP Q76K71
A	-3	ARG	-	EXPRESSION TAG	UNP Q76K71
A	-2	GLY	-	EXPRESSION TAG	UNP Q76K71
A	-1	SER	-	EXPRESSION TAG	UNP Q76K71
A	0	HIS	-	EXPRESSION TAG	UNP Q76K71
B	-19	MET	-	INITIATING METHIONINE	UNP Q76K71
B	-18	GLY	-	EXPRESSION TAG	UNP Q76K71
B	-17	SER	-	EXPRESSION TAG	UNP Q76K71
B	-16	SER	-	EXPRESSION TAG	UNP Q76K71
B	-15	HIS	-	EXPRESSION TAG	UNP Q76K71
B	-14	HIS	-	EXPRESSION TAG	UNP Q76K71
B	-13	HIS	-	EXPRESSION TAG	UNP Q76K71
B	-12	HIS	-	EXPRESSION TAG	UNP Q76K71
B	-11	HIS	-	EXPRESSION TAG	UNP Q76K71
B	-10	HIS	-	EXPRESSION TAG	UNP Q76K71
B	-9	SER	-	EXPRESSION TAG	UNP Q76K71
B	-8	SER	-	EXPRESSION TAG	UNP Q76K71
B	-7	GLY	-	EXPRESSION TAG	UNP Q76K71
B	-6	LEU	-	EXPRESSION TAG	UNP Q76K71
B	-5	VAL	-	EXPRESSION TAG	UNP Q76K71
B	-4	PRO	-	EXPRESSION TAG	UNP Q76K71
B	-3	ARG	-	EXPRESSION TAG	UNP Q76K71
B	-2	GLY	-	EXPRESSION TAG	UNP Q76K71
B	-1	SER	-	EXPRESSION TAG	UNP Q76K71
B	0	HIS	-	EXPRESSION TAG	UNP Q76K71
C	-19	MET	-	INITIATING METHIONINE	UNP Q76K71
C	-18	GLY	-	EXPRESSION TAG	UNP Q76K71
C	-17	SER	-	EXPRESSION TAG	UNP Q76K71
C	-16	SER	-	EXPRESSION TAG	UNP Q76K71
C	-15	HIS	-	EXPRESSION TAG	UNP Q76K71
C	-14	HIS	-	EXPRESSION TAG	UNP Q76K71
C	-13	HIS	-	EXPRESSION TAG	UNP Q76K71
C	-12	HIS	-	EXPRESSION TAG	UNP Q76K71
C	-11	HIS	-	EXPRESSION TAG	UNP Q76K71
C	-10	HIS	-	EXPRESSION TAG	UNP Q76K71
C	-9	SER	-	EXPRESSION TAG	UNP Q76K71
C	-8	SER	-	EXPRESSION TAG	UNP Q76K71
C	-7	GLY	-	EXPRESSION TAG	UNP Q76K71
C	-6	LEU	-	EXPRESSION TAG	UNP Q76K71
C	-5	VAL	-	EXPRESSION TAG	UNP Q76K71

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-4	PRO	-	EXPRESSION TAG	UNP Q76K71
C	-3	ARG	-	EXPRESSION TAG	UNP Q76K71
C	-2	GLY	-	EXPRESSION TAG	UNP Q76K71
C	-1	SER	-	EXPRESSION TAG	UNP Q76K71
C	0	HIS	-	EXPRESSION TAG	UNP Q76K71
D	-19	MET	-	INITIATING METHIONINE	UNP Q76K71
D	-18	GLY	-	EXPRESSION TAG	UNP Q76K71
D	-17	SER	-	EXPRESSION TAG	UNP Q76K71
D	-16	SER	-	EXPRESSION TAG	UNP Q76K71
D	-15	HIS	-	EXPRESSION TAG	UNP Q76K71
D	-14	HIS	-	EXPRESSION TAG	UNP Q76K71
D	-13	HIS	-	EXPRESSION TAG	UNP Q76K71
D	-12	HIS	-	EXPRESSION TAG	UNP Q76K71
D	-11	HIS	-	EXPRESSION TAG	UNP Q76K71
D	-10	HIS	-	EXPRESSION TAG	UNP Q76K71
D	-9	SER	-	EXPRESSION TAG	UNP Q76K71
D	-8	SER	-	EXPRESSION TAG	UNP Q76K71
D	-7	GLY	-	EXPRESSION TAG	UNP Q76K71
D	-6	LEU	-	EXPRESSION TAG	UNP Q76K71
D	-5	VAL	-	EXPRESSION TAG	UNP Q76K71
D	-4	PRO	-	EXPRESSION TAG	UNP Q76K71
D	-3	ARG	-	EXPRESSION TAG	UNP Q76K71
D	-2	GLY	-	EXPRESSION TAG	UNP Q76K71
D	-1	SER	-	EXPRESSION TAG	UNP Q76K71
D	0	HIS	-	EXPRESSION TAG	UNP Q76K71
E	-19	MET	-	INITIATING METHIONINE	UNP Q76K71
E	-18	GLY	-	EXPRESSION TAG	UNP Q76K71
E	-17	SER	-	EXPRESSION TAG	UNP Q76K71
E	-16	SER	-	EXPRESSION TAG	UNP Q76K71
E	-15	HIS	-	EXPRESSION TAG	UNP Q76K71
E	-14	HIS	-	EXPRESSION TAG	UNP Q76K71
E	-13	HIS	-	EXPRESSION TAG	UNP Q76K71
E	-12	HIS	-	EXPRESSION TAG	UNP Q76K71
E	-11	HIS	-	EXPRESSION TAG	UNP Q76K71
E	-10	HIS	-	EXPRESSION TAG	UNP Q76K71
E	-9	SER	-	EXPRESSION TAG	UNP Q76K71
E	-8	SER	-	EXPRESSION TAG	UNP Q76K71
E	-7	GLY	-	EXPRESSION TAG	UNP Q76K71
E	-6	LEU	-	EXPRESSION TAG	UNP Q76K71
E	-5	VAL	-	EXPRESSION TAG	UNP Q76K71
E	-4	PRO	-	EXPRESSION TAG	UNP Q76K71
E	-3	ARG	-	EXPRESSION TAG	UNP Q76K71

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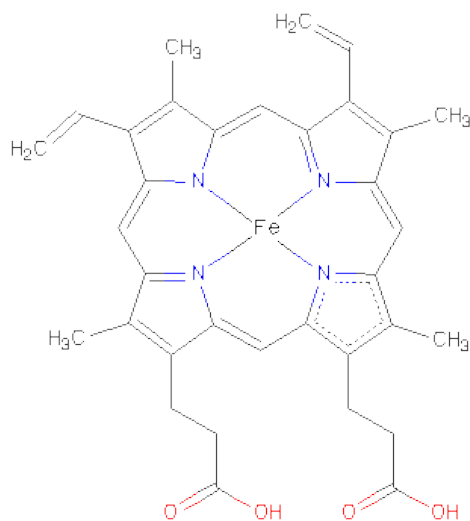
Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	GLY	-	EXPRESSION TAG	UNP Q76K71
E	-1	SER	-	EXPRESSION TAG	UNP Q76K71
E	0	HIS	-	EXPRESSION TAG	UNP Q76K71
F	-19	MET	-	INITIATING METHIONINE	UNP Q76K71
F	-18	GLY	-	EXPRESSION TAG	UNP Q76K71
F	-17	SER	-	EXPRESSION TAG	UNP Q76K71
F	-16	SER	-	EXPRESSION TAG	UNP Q76K71
F	-15	HIS	-	EXPRESSION TAG	UNP Q76K71
F	-14	HIS	-	EXPRESSION TAG	UNP Q76K71
F	-13	HIS	-	EXPRESSION TAG	UNP Q76K71
F	-12	HIS	-	EXPRESSION TAG	UNP Q76K71
F	-11	HIS	-	EXPRESSION TAG	UNP Q76K71
F	-10	HIS	-	EXPRESSION TAG	UNP Q76K71
F	-9	SER	-	EXPRESSION TAG	UNP Q76K71
F	-8	SER	-	EXPRESSION TAG	UNP Q76K71
F	-7	GLY	-	EXPRESSION TAG	UNP Q76K71
F	-6	LEU	-	EXPRESSION TAG	UNP Q76K71
F	-5	VAL	-	EXPRESSION TAG	UNP Q76K71
F	-4	PRO	-	EXPRESSION TAG	UNP Q76K71
F	-3	ARG	-	EXPRESSION TAG	UNP Q76K71
F	-2	GLY	-	EXPRESSION TAG	UNP Q76K71
F	-1	SER	-	EXPRESSION TAG	UNP Q76K71
F	0	HIS	-	EXPRESSION TAG	UNP Q76K71
G	-19	MET	-	INITIATING METHIONINE	UNP Q76K71
G	-18	GLY	-	EXPRESSION TAG	UNP Q76K71
G	-17	SER	-	EXPRESSION TAG	UNP Q76K71
G	-16	SER	-	EXPRESSION TAG	UNP Q76K71
G	-15	HIS	-	EXPRESSION TAG	UNP Q76K71
G	-14	HIS	-	EXPRESSION TAG	UNP Q76K71
G	-13	HIS	-	EXPRESSION TAG	UNP Q76K71
G	-12	HIS	-	EXPRESSION TAG	UNP Q76K71
G	-11	HIS	-	EXPRESSION TAG	UNP Q76K71
G	-10	HIS	-	EXPRESSION TAG	UNP Q76K71
G	-9	SER	-	EXPRESSION TAG	UNP Q76K71
G	-8	SER	-	EXPRESSION TAG	UNP Q76K71
G	-7	GLY	-	EXPRESSION TAG	UNP Q76K71
G	-6	LEU	-	EXPRESSION TAG	UNP Q76K71
G	-5	VAL	-	EXPRESSION TAG	UNP Q76K71
G	-4	PRO	-	EXPRESSION TAG	UNP Q76K71
G	-3	ARG	-	EXPRESSION TAG	UNP Q76K71
G	-2	GLY	-	EXPRESSION TAG	UNP Q76K71
G	-1	SER	-	EXPRESSION TAG	UNP Q76K71

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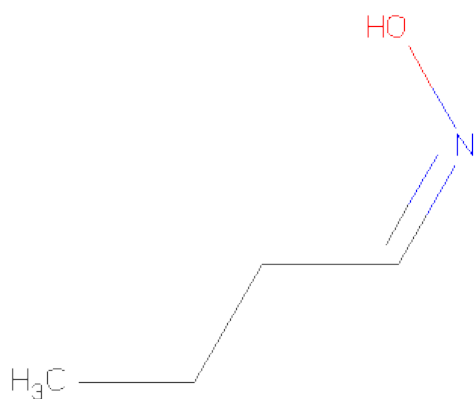
Chain	Residue	Modelled	Actual	Comment	Reference
G	0	HIS	-	EXPRESSION TAG	UNP Q76K71
H	-19	MET	-	INITIATING METHIONINE	UNP Q76K71
H	-18	GLY	-	EXPRESSION TAG	UNP Q76K71
H	-17	SER	-	EXPRESSION TAG	UNP Q76K71
H	-16	SER	-	EXPRESSION TAG	UNP Q76K71
H	-15	HIS	-	EXPRESSION TAG	UNP Q76K71
H	-14	HIS	-	EXPRESSION TAG	UNP Q76K71
H	-13	HIS	-	EXPRESSION TAG	UNP Q76K71
H	-12	HIS	-	EXPRESSION TAG	UNP Q76K71
H	-11	HIS	-	EXPRESSION TAG	UNP Q76K71
H	-10	HIS	-	EXPRESSION TAG	UNP Q76K71
H	-9	SER	-	EXPRESSION TAG	UNP Q76K71
H	-8	SER	-	EXPRESSION TAG	UNP Q76K71
H	-7	GLY	-	EXPRESSION TAG	UNP Q76K71
H	-6	LEU	-	EXPRESSION TAG	UNP Q76K71
H	-5	VAL	-	EXPRESSION TAG	UNP Q76K71
H	-4	PRO	-	EXPRESSION TAG	UNP Q76K71
H	-3	ARG	-	EXPRESSION TAG	UNP Q76K71
H	-2	GLY	-	EXPRESSION TAG	UNP Q76K71
H	-1	SER	-	EXPRESSION TAG	UNP Q76K71
H	0	HIS	-	EXPRESSION TAG	UNP Q76K71

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	D	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	E	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	F	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	G	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	H	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is (1Z)-BUTANAL OXIME (three-letter code: BXO) (formula: C<sub>4</sub>H<sub>9</sub>NO).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			6	4	1	1		
3	B	1	Total	C	N	O	0	0
			6	4	1	1		
3	C	1	Total	C	N	O	0	0
			6	4	1	1		
3	D	1	Total	C	N	O	0	0
			6	4	1	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	E	1	Total	C	N	O	0	0
			6	4	1	1		
3	F	1	Total	C	N	O	0	0
			6	4	1	1		
3	G	1	Total	C	N	O	0	0
			6	4	1	1		
3	H	1	Total	C	N	O	0	0
			6	4	1	1		

- Molecule 4 is water.

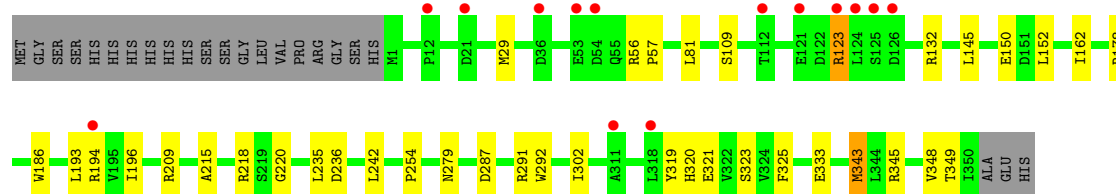
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	105	Total	O	0	0
			105	105		
4	B	60	Total	O	0	0
			60	60		
4	C	93	Total	O	0	0
			93	93		
4	D	77	Total	O	0	0
			77	77		
4	E	83	Total	O	0	0
			83	83		
4	F	58	Total	O	0	0
			58	58		
4	G	84	Total	O	0	0
			84	84		
4	H	59	Total	O	0	0
			59	59		

### 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 errors 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.

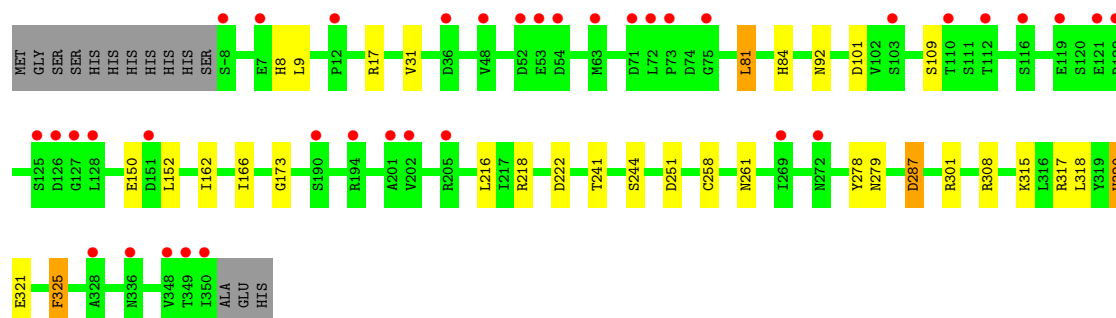
- Molecule 1: Aldoxime dehydratase

Chain A: 



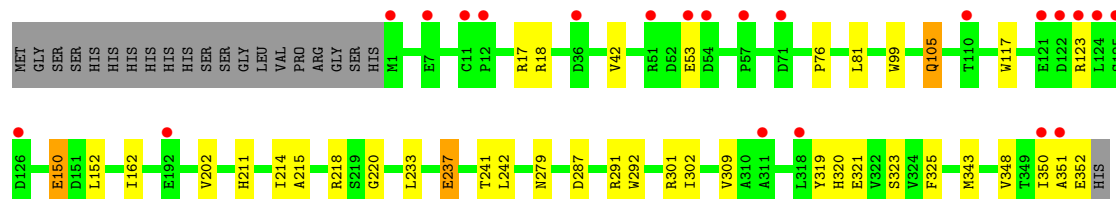
- Molecule 1: Aldoxime dehydratase

Chain B: 



- Molecule 1: Aldoxime dehydratase

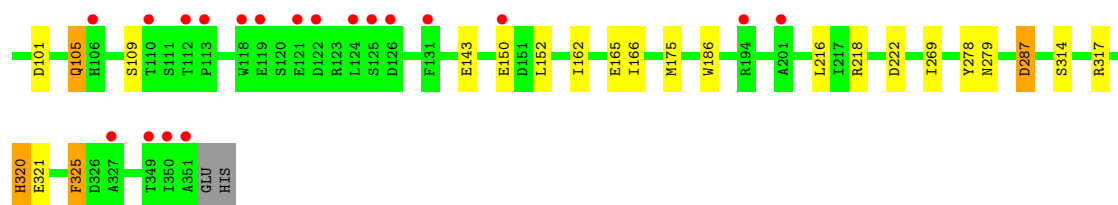
Chain C: 



- Molecule 1: Aldoxime dehydratase

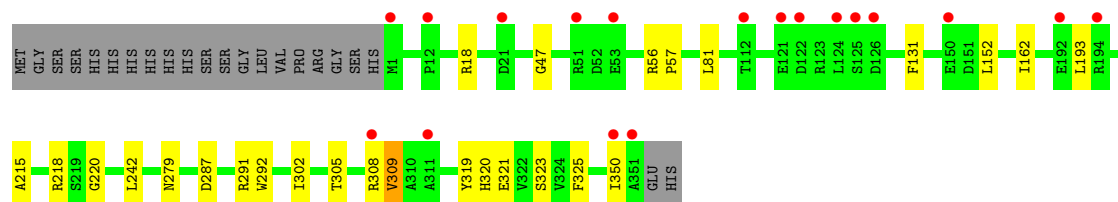
Chain D: 





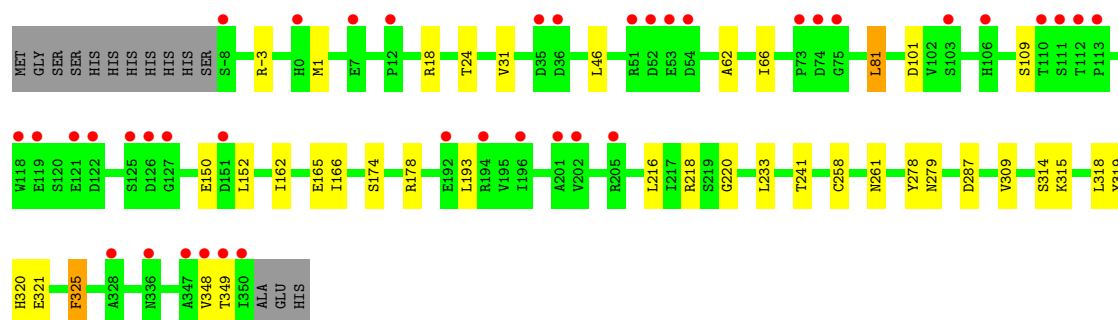
- Molecule 1: Aldoxime dehydratase

Chain E:



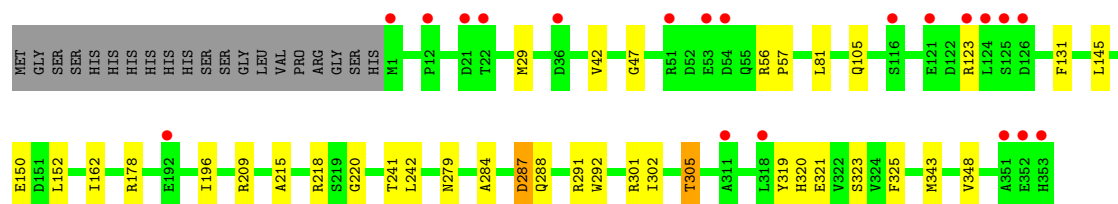
- Molecule 1: Aldoxime dehydratase

Chain F:



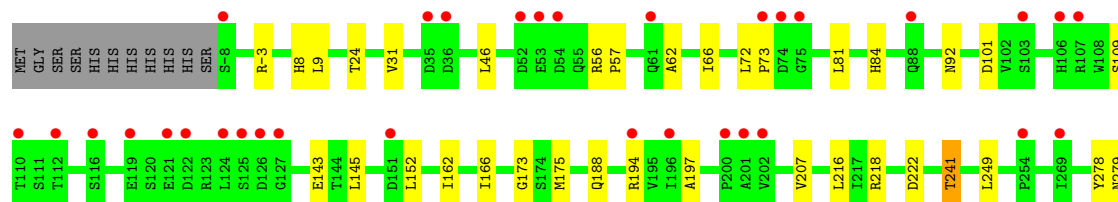
- Molecule 1: Aldoxime dehydratase

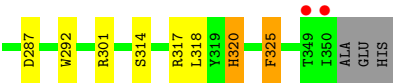
Chain G:



- Molecule 1: Aldoxime dehydratase

Chain H:





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	80.05Å 103.93Å 114.19Å 76.50° 89.46° 87.55°	Depositor
Resolution (Å)	19.90 – 2.50 19.90 – 2.50	Depositor EDS
% Data completeness (in resolution range)	91.2 (19.90-2.50) 91.2 (19.90-2.50)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.30 (at 2.50Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.211 , 0.243 0.225 , 0.261	Depositor DCC
$R_{free}$ test set	5656 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	30.5	Xtriage
Anisotropy	0.447	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 19.6	EDS
Estimated twinning fraction	0.098 for h,-k,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 112560 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	23656	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.34% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

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

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	0.44	0/2866	0.57	0/3900
1	B	0.42	0/2931	0.57	0/3988
1	C	0.44	0/2891	0.57	0/3934
1	D	0.44	0/2947	0.58	0/4009
1	E	0.43	0/2882	0.58	0/3921
1	F	0.42	0/2931	0.57	0/3988
1	G	0.44	0/2892	0.58	0/3934
1	H	0.42	0/2931	0.56	0/3988
All	All	0.43	0/23271	0.57	0/31662

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2790	0	2628	25	0
1	B	2853	0	2691	18	0
1	C	2812	0	2652	21	0
1	D	2866	0	2709	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2803	0	2646	16	0
1	F	2853	0	2691	16	0
1	G	2815	0	2646	21	0
1	H	2853	0	2691	21	0
2	A	43	0	30	4	0
2	B	43	0	30	2	0
2	C	43	0	30	4	0
2	D	43	0	30	1	0
2	E	43	0	30	4	0
2	F	43	0	30	2	0
2	G	43	0	30	4	0
2	H	43	0	30	3	0
3	A	6	0	9	2	0
3	B	6	0	9	1	0
3	C	6	0	9	0	0
3	D	6	0	9	1	0
3	E	6	0	9	0	0
3	F	6	0	9	1	0
3	G	6	0	9	1	0
3	H	6	0	9	3	0
4	A	105	0	0	1	0
4	B	60	0	0	1	0
4	C	93	0	0	0	0
4	D	77	0	0	3	0
4	E	83	0	0	0	0
4	F	58	0	0	0	0
4	G	84	0	0	0	0
4	H	59	0	0	1	0
All	All	23656	0	21666	163	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 4.

All (163) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:-3[A]:ARG:HG2	1:D:-3[A]:ARG:HH11	1.41	0.84
1:G:242:LEU:HD22	2:G:354:HEM:HBB1	1.64	0.80
1:E:242:LEU:HD22	2:E:354:HEM:HBB1	1.69	0.75
1:E:305:THR:O	1:E:309:VAL:HG23	1.88	0.73
1:C:242:LEU:HD22	2:C:354:HEM:HBB1	1.70	0.73
1:D:-8:SER:HA	4:D:516:HOH:O	1.89	0.70
1:F:318:LEU:HD22	2:F:354:HEM:HBB2	1.73	0.70

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:29:MET:HE1	3:A:355:BXO:H4	1.74	0.69
1:G:302:ILE:HA	1:G:305:THR:HG23	1.75	0.67
1:E:308:ARG:HE	1:E:309:VAL:HG22	1.58	0.67
1:G:292:TRP:CZ3	2:G:354:HEM:HBC2	2.29	0.67
1:D:-3[A]:ARG:HG2	1:D:-3[A]:ARG:NH1	2.09	0.64
1:B:31:VAL:HG13	1:B:166:ILE:HG12	1.78	0.64
1:A:235:LEU:HB3	1:D:-3[A]:ARG:HD2	1.79	0.63
1:H:318:LEU:HD22	2:H:354:HEM:HBB2	1.79	0.63
1:C:218:ARG:HB3	1:C:321:GLU:HG2	1.80	0.62
1:A:218:ARG:HB3	1:A:321:GLU:HG2	1.80	0.62
1:G:242:LEU:HD22	2:G:354:HEM:CBB	2.30	0.61
1:F:218:ARG:HG3	1:F:278:TYR:CD2	2.36	0.61
1:A:242:LEU:HD22	2:A:354:HEM:HBB1	1.82	0.60
1:H:218:ARG:HG3	1:H:278:TYR:CD2	2.37	0.59
1:A:292:TRP:CZ3	2:A:354:HEM:HBC2	2.37	0.59
1:G:301:ARG:O	1:G:305:THR:HG22	2.04	0.58
1:C:152:LEU:HD21	1:C:162:ILE:HD13	1.85	0.57
1:E:242:LEU:HD22	2:E:354:HEM:CBB	2.35	0.57
1:B:318:LEU:HD22	2:B:354:HEM:HBB2	1.86	0.57
1:G:218:ARG:HB3	1:G:321:GLU:HG2	1.87	0.56
1:E:218:ARG:HB3	1:E:321:GLU:HG2	1.87	0.56
1:A:235:LEU:HB3	1:D:-3[B]:ARG:HD2	1.86	0.56
1:F:31:VAL:HG13	1:F:166:ILE:HG12	1.87	0.56
1:C:150:GLU:N	1:C:150:GLU:OE1	2.37	0.55
1:B:218:ARG:HG3	1:B:278:TYR:CD2	2.42	0.55
1:C:292:TRP:CZ3	2:C:354:HEM:HBC2	2.42	0.55
1:B:81:LEU:N	1:B:81:LEU:HD23	2.22	0.54
1:G:292:TRP:HZ3	2:G:354:HEM:HBC2	1.70	0.54
1:E:292:TRP:CZ3	2:E:354:HEM:HBC2	2.42	0.54
1:H:216:LEU:HB3	1:H:325:PHE:HZ	1.73	0.54
1:H:31:VAL:HG13	1:H:166:ILE:HG12	1.89	0.54
1:B:251:ASP:HB3	1:C:202:VAL:HG21	1.90	0.53
1:B:244:SER:OG	1:C:351:ALA:HB3	2.07	0.53
1:E:220:GLY:HA3	1:E:319:TYR:CE1	2.44	0.53
1:F:233:LEU:HD21	1:F:309:VAL:HG11	1.91	0.52
1:C:242:LEU:HD22	2:C:354:HEM:CBB	2.38	0.52
1:A:193:LEU:HD11	1:A:343:MET:HG3	1.92	0.52
1:H:145:LEU:HD23	3:H:355:BXO:C4	2.39	0.52
1:A:152:LEU:HD21	1:A:162:ILE:HD13	1.92	0.52
1:C:287:ASP:O	1:C:291:ARG:HG2	2.09	0.51
1:F:220:GLY:HA3	1:F:319:TYR:CE1	2.46	0.51
1:D:42:VAL:HG11	1:D:105:GLN:HG3	1.91	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:287:ASP:O	1:G:291:ARG:HG2	2.12	0.50
1:D:287:ASP:N	1:D:287:ASP:OD1	2.44	0.50
1:G:42:VAL:HG11	1:G:105:GLN:HG3	1.94	0.50
1:A:292:TRP:HZ3	2:A:354:HEM:HBC2	1.77	0.49
1:B:218:ARG:HA	1:B:279:ASN:O	2.13	0.49
1:F:1:MET:HE3	1:F:165:GLU:HB3	1.94	0.49
1:F:218:ARG:HA	1:F:279:ASN:O	2.12	0.49
1:D:218:ARG:HG3	1:D:278:TYR:CD2	2.48	0.49
1:D:31:VAL:HG13	1:D:166:ILE:HG12	1.93	0.49
1:H:72:LEU:HB3	1:H:73:PRO:HD2	1.95	0.49
1:A:302:ILE:HD12	2:A:354:HEM:HBB2	1.94	0.49
1:D:218:ARG:HA	1:D:279:ASN:O	2.13	0.49
1:F:174:SER:O	1:F:178:ARG:HG3	2.13	0.48
1:H:218:ARG:HA	1:H:279:ASN:O	2.13	0.48
1:E:287:ASP:O	1:E:291:ARG:HG2	2.13	0.48
1:H:145:LEU:HD23	3:H:355:BXO:H4A	1.95	0.47
1:C:220:GLY:HA3	1:C:319:TYR:CE1	2.49	0.47
1:C:117:TRP:O	1:C:123:ARG:NH1	2.47	0.47
1:E:215:ALA:HA	1:E:323:SER:O	2.15	0.47
1:E:302:ILE:HD12	2:E:354:HEM:HBB2	1.96	0.47
1:H:222:ASP:HB3	1:H:317:ARG:HB2	1.96	0.47
1:C:215:ALA:HA	1:C:323:SER:O	2.14	0.47
1:H:84:HIS:CE1	1:H:92:ASN:HB2	2.49	0.47
1:B:287:ASP:N	1:B:287:ASP:OD1	2.44	0.47
1:D:218:ARG:HB3	1:D:321:GLU:HG2	1.96	0.47
1:A:220:GLY:HA3	1:A:319:TYR:CE1	2.50	0.46
1:G:29:MET:HE1	3:G:355:BXO:H4	1.98	0.46
1:D:62:ALA:O	1:D:66:ILE:HG13	2.16	0.46
1:B:222:ASP:HB3	1:B:317:ARG:HB2	1.96	0.46
1:G:152:LEU:HD21	1:G:162:ILE:HD13	1.98	0.46
1:B:320:HIS:C	1:B:320:HIS:CD2	2.89	0.46
1:A:215:ALA:HA	1:A:323:SER:O	2.16	0.46
1:D:222:ASP:HB3	1:D:317:ARG:HB2	1.97	0.46
1:H:320:HIS:C	1:H:320:HIS:CD2	2.88	0.46
1:G:215:ALA:HA	1:G:323:SER:O	2.15	0.46
1:E:218:ARG:HA	1:E:279:ASN:O	2.16	0.46
1:F:152:LEU:HD21	1:F:162:ILE:HD13	1.98	0.46
1:A:29:MET:CE	3:A:355:BXO:H4	2.44	0.45
1:C:211:HIS:CE1	1:C:214:ILE:HG13	2.52	0.45
1:F:216:LEU:HB3	1:F:325:PHE:HZ	1.80	0.45
1:A:236:ASP:OD1	1:D:-3[B]:ARG:HD3	2.17	0.45
1:D:84:HIS:CE1	1:D:92:ASN:HB2	2.51	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:193:LEU:O	1:F:348:VAL:HG12	2.17	0.45
1:F:218:ARG:HB3	1:F:321:GLU:HG2	1.99	0.44
1:G:220:GLY:HA3	1:G:319:TYR:CE1	2.53	0.44
1:D:143:GLU:OE2	1:D:175:MET:HG3	2.18	0.44
1:H:249:LEU:HD21	1:H:292:TRP:CE2	2.52	0.44
1:A:287:ASP:O	1:A:291:ARG:HG2	2.16	0.44
1:E:308:ARG:NE	1:E:309:VAL:HG22	2.31	0.44
1:D:11:CYS:HB2	1:D:12:PRO:HD2	1.99	0.44
1:G:218:ARG:HA	1:G:279:ASN:O	2.18	0.44
1:G:150:GLU:OE1	1:G:150:GLU:N	2.49	0.44
1:G:47:GLY:HA3	1:G:131:PHE:CZ	2.53	0.44
1:B:216:LEU:HB3	1:B:325:PHE:HZ	1.83	0.44
1:A:145:LEU:HD22	1:A:178:ARG:NH2	2.33	0.44
1:D:39:GLN:HG2	4:D:514:HOH:O	2.17	0.44
1:A:186:TRP:HH2	1:B:17:ARG:O	1.99	0.43
1:B:218:ARG:HB3	1:B:321:GLU:HG2	2.00	0.43
1:H:173:GLY:HA2	4:H:374:HOH:O	2.18	0.43
1:E:18:ARG:HD2	1:F:18:ARG:HD2	2.00	0.43
1:A:194:ARG:HG2	1:A:348:VAL:CG2	2.49	0.43
1:F:81:LEU:HD23	1:F:81:LEU:N	2.33	0.43
1:C:17:ARG:O	1:D:186:TRP:HH2	2.01	0.43
1:G:196:ILE:HD13	1:G:209:ARG:HB2	2.00	0.43
1:G:145:LEU:HD22	1:G:178:ARG:NH2	2.34	0.43
1:C:233:LEU:HA	1:C:237:GLU:HG3	2.01	0.43
1:G:56:ARG:HB3	1:G:57:PRO:HD3	2.00	0.43
1:A:132:ARG:HG2	1:A:333:GLU:HB2	2.01	0.42
1:B:84:HIS:CE1	1:B:92:ASN:HB2	2.54	0.42
1:D:320:HIS:C	1:D:320:HIS:CD2	2.92	0.42
1:F:62:ALA:O	1:F:66:ILE:HG13	2.19	0.42
1:C:218:ARG:HA	1:C:279:ASN:O	2.19	0.42
1:E:56:ARG:HB3	1:E:57:PRO:HD3	2.00	0.42
1:C:241:THR:HG21	1:C:301:ARG:HB3	2.01	0.42
1:B:173:GLY:HA2	4:B:367:HOH:O	2.20	0.42
1:D:152:LEU:HD21	1:D:162:ILE:HD13	2.01	0.42
1:A:218:ARG:HA	1:A:279:ASN:O	2.19	0.42
1:D:11:CYS:HB2	1:D:12:PRO:CD	2.49	0.41
1:D:24:THR:HA	1:D:25:PRO:HD3	1.95	0.41
1:C:42:VAL:HG11	1:C:105:GLN:HG3	2.02	0.41
1:H:143:GLU:OE2	1:H:175:MET:HG3	2.20	0.41
1:C:18:ARG:HD2	1:D:18:ARG:HD2	2.01	0.41
1:A:56:ARG:HB3	1:A:57:PRO:HD3	2.02	0.41
1:C:302:ILE:HD12	2:C:354:HEM:HBB2	2.02	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:152:LEU:HD21	1:E:162:ILE:HD13	2.02	0.41
1:C:76:PRO:HD3	1:C:99:TRP:CZ2	2.55	0.41
1:D:216:LEU:HB3	1:D:325:PHE:HZ	1.86	0.41
1:H:62:ALA:O	1:H:66:ILE:HG13	2.20	0.41
1:D:218:ARG:HB3	1:D:321:GLU:CG	2.50	0.41
1:A:209:ARG:NE	4:A:588:HOH:O	2.52	0.41
2:D:354:HEM:C4A	3:D:355:BXO:H1	2.56	0.41
1:F:258:CYS:SG	1:F:261:ASN:HB2	2.61	0.41
1:E:47:GLY:HA3	1:E:131:PHE:CZ	2.56	0.41
1:D:1:MET:HE2	1:D:25:PRO:HG2	2.02	0.41
1:A:254:PRO:HD3	1:A:345:ARG:NH2	2.36	0.41
1:H:152:LEU:HD21	1:H:162:ILE:HD13	2.02	0.41
2:B:354:HEM:C4A	3:B:355:BXO:H1	2.56	0.41
1:H:241:THR:HG21	1:H:301:ARG:HD3	2.03	0.41
1:H:56:ARG:HB3	1:H:57:PRO:HD3	2.03	0.41
1:A:123:ARG:HD3	1:A:123:ARG:HA	1.92	0.41
1:B:8:HIS:CE1	1:B:9:LEU:HG	2.56	0.41
1:H:197:ALA:HB3	1:H:207:VAL:HB	2.03	0.41
1:B:258:CYS:SG	1:B:261:ASN:HB2	2.61	0.41
1:D:12:PRO:HD3	4:D:598:HOH:O	2.20	0.41
1:D:1:MET:HE3	1:D:165:GLU:HB3	2.03	0.40
2:H:354:HEM:C4A	3:H:355:BXO:H1	2.56	0.40
1:B:152:LEU:HD21	1:B:162:ILE:HD13	2.02	0.40
1:A:196:ILE:HD13	1:A:209:ARG:HB2	2.02	0.40
1:H:279:ASN:OD1	2:H:354:HEM:HAC	2.22	0.40
1:D:64:ARG:HG2	1:D:269:ILE:HD11	2.04	0.40
1:G:284:ALA:HB3	1:G:288:GLN:HE22	1.87	0.40
2:F:354:HEM:C4A	3:F:355:BXO:H1	2.56	0.40
1:G:241:THR:OG1	1:G:302:ILE:HG23	2.22	0.40
1:H:8:HIS:CE1	1:H:9:LEU:HG	2.56	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

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	348/373 (93%)	346 (99%)	2 (1%)	0	100	100
1	B	357/373 (96%)	354 (99%)	3 (1%)	0	100	100
1	C	351/373 (94%)	346 (99%)	5 (1%)	0	100	100
1	D	359/373 (96%)	357 (99%)	2 (1%)	0	100	100
1	E	350/373 (94%)	346 (99%)	4 (1%)	0	100	100
1	F	357/373 (96%)	354 (99%)	3 (1%)	0	100	100
1	G	351/373 (94%)	347 (99%)	4 (1%)	0	100	100
1	H	357/373 (96%)	357 (100%)	0	0	100	100
All	All	2830/2984 (95%)	2807 (99%)	23 (1%)	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	293/312 (94%)	285 (97%)	8 (3%)	57	83
1	B	300/312 (96%)	289 (96%)	11 (4%)	45	72
1	C	295/312 (95%)	283 (96%)	12 (4%)	41	67
1	D	301/312 (96%)	287 (95%)	14 (5%)	36	61
1	E	294/312 (94%)	288 (98%)	6 (2%)	68	89
1	F	300/312 (96%)	286 (95%)	14 (5%)	36	61
1	G	295/312 (95%)	287 (97%)	8 (3%)	57	83
1	H	300/312 (96%)	287 (96%)	13 (4%)	40	65
All	All	2378/2496 (95%)	2292 (96%)	86 (4%)	47	73

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	81	LEU
1	A	109	SER
1	A	123	ARG

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Mol	Chain	Res	Type
1	A	150	GLU
1	A	320	HIS
1	A	325	PHE
1	A	343	MET
1	A	349	THR
1	B	81	LEU
1	B	101	ASP
1	B	109	SER
1	B	150	GLU
1	B	241	THR
1	B	287	ASP
1	B	301	ARG
1	B	308	ARG
1	B	315	LYS
1	B	320	HIS
1	B	325	PHE
1	C	53	GLU
1	C	81	LEU
1	C	105	GLN
1	C	150	GLU
1	C	237	GLU
1	C	309	VAL
1	C	320	HIS
1	C	325	PHE
1	C	343	MET
1	C	348	VAL
1	C	350	ILE
1	C	352	GLU
1	D	-3[A]	ARG
1	D	-3[B]	ARG
1	D	24	THR
1	D	38	LEU
1	D	46	LEU
1	D	81	LEU
1	D	101	ASP
1	D	105	GLN
1	D	109	SER
1	D	150	GLU
1	D	287	ASP
1	D	314	SER
1	D	320	HIS
1	D	325	PHE

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Mol	Chain	Res	Type
1	E	81	LEU
1	E	193	LEU
1	E	309	VAL
1	E	320	HIS
1	E	325	PHE
1	E	350	ILE
1	F	-3	ARG
1	F	24	THR
1	F	46	LEU
1	F	81	LEU
1	F	101	ASP
1	F	109	SER
1	F	150	GLU
1	F	241	THR
1	F	287	ASP
1	F	314	SER
1	F	315	LYS
1	F	320	HIS
1	F	325	PHE
1	F	349	THR
1	G	81	LEU
1	G	123	ARG
1	G	287	ASP
1	G	305	THR
1	G	320	HIS
1	G	325	PHE
1	G	343	MET
1	G	348	VAL
1	H	-3	ARG
1	H	24	THR
1	H	46	LEU
1	H	81	LEU
1	H	101	ASP
1	H	109	SER
1	H	188	GLN
1	H	194	ARG
1	H	241	THR
1	H	287	ASP
1	H	314	SER
1	H	320	HIS
1	H	325	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	288	GLN
1	B	105	GLN
1	C	288	GLN
1	D	105	GLN
1	E	105	GLN
1	E	288	GLN
1	F	105	GLN
1	G	105	GLN
1	G	288	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

16 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	$\# Z  > 2$
2	HEM	A	354	1,3	49,50,50	2.41	13 (26%)	46,82,82	2.14	9 (19%)
3	BXO	A	355	2	5,5,5	1.16	1 (20%)	4,4,4	5.13	1 (25%)
2	HEM	B	354	1,3	49,50,50	2.71	16 (32%)	46,82,82	2.02	10 (21%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	BXO	B	355	2	5,5,5	1.41	1 (20%)	4,4,4	3.44	2 (50%)
2	HEM	C	354	1,3	49,50,50	2.37	12 (24%)	46,82,82	2.10	9 (19%)
3	BXO	C	355	2	5,5,5	1.81	1 (20%)	4,4,4	4.02	2 (50%)
2	HEM	D	354	1,3	49,50,50	2.62	16 (32%)	46,82,82	1.99	11 (23%)
3	BXO	D	355	2	5,5,5	1.54	1 (20%)	4,4,4	3.33	2 (50%)
2	HEM	E	354	1,3	49,50,50	2.42	13 (26%)	46,82,82	2.12	9 (19%)
3	BXO	E	355	2	5,5,5	1.77	1 (20%)	4,4,4	3.58	2 (50%)
2	HEM	F	354	1,3	49,50,50	2.52	15 (30%)	46,82,82	2.06	10 (21%)
3	BXO	F	355	2	5,5,5	1.41	1 (20%)	4,4,4	3.16	2 (50%)
2	HEM	G	354	1,3	49,50,50	2.50	14 (28%)	46,82,82	2.10	9 (19%)
3	BXO	G	355	2	5,5,5	1.62	1 (20%)	4,4,4	3.51	2 (50%)
2	HEM	H	354	1,3	49,50,50	2.68	13 (26%)	46,82,82	2.01	10 (21%)
3	BXO	H	355	2	5,5,5	1.33	1 (20%)	4,4,4	3.30	2 (50%)

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	HEM	A	354	1,3	-	0/14/114/114	0/0/8/8
3	BXO	A	355	2	-	0/2/3/3	0/0/0/0
2	HEM	B	354	1,3	-	0/14/114/114	0/0/8/8
3	BXO	B	355	2	-	0/2/3/3	0/0/0/0
2	HEM	C	354	1,3	-	0/14/114/114	0/0/8/8
3	BXO	C	355	2	-	0/2/3/3	0/0/0/0
2	HEM	D	354	1,3	-	0/14/114/114	0/0/8/8
3	BXO	D	355	2	-	0/2/3/3	0/0/0/0
2	HEM	E	354	1,3	-	0/14/114/114	0/0/8/8
3	BXO	E	355	2	-	0/2/3/3	0/0/0/0
2	HEM	F	354	1,3	-	0/14/114/114	0/0/8/8
3	BXO	F	355	2	-	0/2/3/3	0/0/0/0
2	HEM	G	354	1,3	-	0/14/114/114	0/0/8/8
3	BXO	G	355	2	-	0/2/3/3	0/0/0/0
2	HEM	H	354	1,3	-	0/14/114/114	0/0/8/8
3	BXO	H	355	2	-	0/2/3/3	0/0/0/0

All (120) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	354	HEM	C2B-C1B	9.04	1.46	1.44
2	E	354	HEM	C2B-C1B	8.44	1.46	1.44
2	B	354	HEM	C2B-C1B	8.31	1.46	1.44
2	C	354	HEM	C2B-C1B	8.00	1.46	1.44
2	G	354	HEM	C2B-C1B	7.93	1.46	1.44
2	H	354	HEM	C2B-C1B	7.79	1.46	1.44
2	F	354	HEM	C2B-C1B	7.45	1.46	1.44
2	H	354	HEM	C3D-C4D	7.04	1.46	1.44
2	A	354	HEM	C2B-C1B	6.96	1.46	1.44
2	B	354	HEM	C2D-C1D	6.29	1.46	1.44
2	B	354	HEM	C3B-C2B	-5.76	1.33	1.43
2	B	354	HEM	C3D-C4D	5.58	1.46	1.44
2	H	354	HEM	C3B-C2B	-5.58	1.34	1.43
2	E	354	HEM	C3C-C2C	-5.55	1.34	1.43
2	F	354	HEM	C3B-C2B	-5.54	1.34	1.43
2	H	354	HEM	C2D-C1D	5.49	1.45	1.44
2	D	354	HEM	C3B-C2B	-5.47	1.34	1.43
2	C	354	HEM	C3C-C2C	-5.44	1.34	1.43
2	G	354	HEM	C3C-C2C	-5.41	1.34	1.43
2	A	354	HEM	C3C-C2C	-5.40	1.34	1.43
2	A	354	HEM	C3D-C4D	5.34	1.45	1.44
2	E	354	HEM	C3B-C2B	-5.30	1.34	1.43
2	E	354	HEM	C3D-C2D	5.27	1.53	1.43
2	C	354	HEM	C3D-C2D	5.25	1.52	1.43
2	D	354	HEM	C3C-C2C	-5.22	1.34	1.43
2	G	354	HEM	C3D-C2D	5.15	1.52	1.43
2	B	354	HEM	C3C-C2C	-5.14	1.34	1.43
2	G	354	HEM	C3B-C2B	-5.14	1.34	1.43
2	A	354	HEM	C3D-C2D	5.06	1.52	1.43
2	F	354	HEM	C3C-C2C	-5.04	1.34	1.43
2	H	354	HEM	C3C-C2C	-5.04	1.35	1.43
2	B	354	HEM	C3D-C2D	4.99	1.52	1.43
2	F	354	HEM	C3D-C2D	4.99	1.52	1.43
2	G	354	HEM	C3D-C4D	4.98	1.45	1.44
2	C	354	HEM	C3B-C2B	-4.98	1.35	1.43
2	A	354	HEM	C3B-C2B	-4.93	1.35	1.43
2	F	354	HEM	C3C-CAC	4.93	1.55	1.40
2	B	354	HEM	C3B-CAB	4.89	1.55	1.40
2	D	354	HEM	C3D-C4D	4.88	1.45	1.44
2	H	354	HEM	C3C-CAC	4.85	1.55	1.40
2	H	354	HEM	C3D-C2D	4.84	1.52	1.43
2	G	354	HEM	C4A-C3A	4.82	1.46	1.40
2	D	354	HEM	C3B-CAB	4.75	1.55	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	354	HEM	C3C-CAC	4.74	1.55	1.40
2	D	354	HEM	C3D-C2D	4.72	1.52	1.43
2	D	354	HEM	C3C-CAC	4.72	1.55	1.40
2	E	354	HEM	C3B-CAB	4.71	1.55	1.40
2	H	354	HEM	C3B-CAB	4.68	1.55	1.40
2	A	354	HEM	C3B-CAB	4.68	1.55	1.40
2	E	354	HEM	C4A-C3A	4.65	1.46	1.40
2	F	354	HEM	C3B-CAB	4.64	1.55	1.40
2	G	354	HEM	C3B-CAB	4.64	1.55	1.40
2	C	354	HEM	C3B-CAB	4.59	1.54	1.40
2	C	354	HEM	C4A-C3A	4.58	1.45	1.40
2	A	354	HEM	C4A-C3A	4.52	1.45	1.40
2	D	354	HEM	C4A-C3A	4.47	1.45	1.40
2	C	354	HEM	C3C-CAC	4.41	1.54	1.40
2	F	354	HEM	C4A-C3A	4.39	1.45	1.40
2	A	354	HEM	C3C-CAC	4.38	1.54	1.40
2	F	354	HEM	C2D-C1D	4.37	1.45	1.44
2	E	354	HEM	C3C-CAC	4.36	1.54	1.40
2	G	354	HEM	C3C-CAC	4.24	1.53	1.40
2	H	354	HEM	C4A-C3A	4.24	1.45	1.40
2	F	354	HEM	C3D-C4D	4.20	1.45	1.44
2	B	354	HEM	C4A-C3A	4.15	1.45	1.40
2	D	354	HEM	C2D-C1D	3.99	1.45	1.44
3	C	355	BXO	C2-C1	3.86	1.53	1.49
2	G	354	HEM	C2D-C1D	3.83	1.45	1.44
3	E	355	BXO	C2-C1	3.78	1.53	1.49
3	G	355	BXO	C2-C1	3.40	1.52	1.49
3	D	355	BXO	C2-C1	3.26	1.52	1.49
2	A	354	HEM	C2D-C1D	3.09	1.45	1.44
2	D	354	HEM	FE-NA	3.09	2.05	1.92
2	C	354	HEM	FE-NA	3.03	2.05	1.92
2	B	354	HEM	CMC-C2C	3.00	1.56	1.47
3	F	355	BXO	C2-C1	2.99	1.52	1.49
3	B	355	BXO	C2-C1	2.92	1.52	1.49
3	H	355	BXO	C2-C1	2.85	1.52	1.49
2	H	354	HEM	CMB-C2B	2.82	1.56	1.47
2	F	354	HEM	CMB-C2B	2.77	1.56	1.47
2	B	354	HEM	CMB-C2B	2.76	1.56	1.47
2	H	354	HEM	FE-NB	2.76	2.07	1.97
2	F	354	HEM	CMC-C2C	2.74	1.55	1.47
2	D	354	HEM	CMC-C2C	2.74	1.55	1.47
2	D	354	HEM	CMB-C2B	2.71	1.55	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	354	HEM	CMC-C2C	2.67	1.55	1.47
2	A	354	HEM	CMB-C2B	2.65	1.55	1.47
2	G	354	HEM	CMB-C2B	2.62	1.55	1.47
2	G	354	HEM	CMC-C2C	2.60	1.55	1.47
2	E	354	HEM	FE-NA	2.58	2.03	1.92
2	F	354	HEM	CMD-C2D	2.57	1.55	1.47
2	E	354	HEM	C2D-C1D	2.54	1.45	1.44
2	A	354	HEM	FE-NA	2.54	2.03	1.92
2	E	354	HEM	CMB-C2B	2.54	1.55	1.47
2	C	354	HEM	CMB-C2B	2.53	1.55	1.47
3	A	355	BXO	C2-C1	2.53	1.51	1.49
2	E	354	HEM	CMC-C2C	2.43	1.55	1.47
2	C	354	HEM	CMC-C2C	2.43	1.54	1.47
2	G	354	HEM	FE-NA	2.41	2.02	1.92
2	F	354	HEM	FE-NB	2.39	2.06	1.97
2	A	354	HEM	CMC-C2C	2.36	1.54	1.47
2	C	354	HEM	CMD-C2D	2.35	1.54	1.47
2	G	354	HEM	CMD-C2D	2.34	1.54	1.47
2	B	354	HEM	FE-NC	2.34	2.06	1.97
2	F	354	HEM	FE-NA	2.33	2.02	1.92
2	E	354	HEM	CMD-C2D	2.32	1.54	1.47
2	A	354	HEM	CMD-C2D	2.26	1.54	1.47
2	C	354	HEM	C3D-C4D	2.23	1.45	1.44
2	H	354	HEM	FE-NC	2.21	2.06	1.97
2	B	354	HEM	FE-ND	2.16	2.05	1.97
2	B	354	HEM	CMD-C2D	2.15	1.54	1.47
2	D	354	HEM	CMD-C2D	2.14	1.54	1.47
2	D	354	HEM	CHA-C4D	2.13	1.38	1.35
2	D	354	HEM	FE-NB	2.14	2.05	1.97
2	B	354	HEM	FE-NA	2.08	2.01	1.92
2	F	354	HEM	CMA-C3A	2.04	1.55	1.51
2	G	354	HEM	FE-NB	2.04	2.05	1.97
2	E	354	HEM	FE-NB	2.03	2.05	1.97
2	B	354	HEM	CHA-C4D	2.02	1.38	1.35
2	D	354	HEM	FE-ND	2.02	2.05	1.97

All (92) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	355	BXO	O1-N1-C1	10.06	121.55	111.66
2	F	354	HEM	C3B-C4B-NB	-7.80	108.42	114.00
2	H	354	HEM	C3B-C4B-NB	-7.77	108.44	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	355	BXO	O1-N1-C1	7.69	119.22	111.66
2	B	354	HEM	C3B-C4B-NB	-7.57	108.58	114.00
2	G	354	HEM	C3B-C4B-NB	-7.51	108.63	114.00
2	A	354	HEM	C3B-C4B-NB	-7.31	108.77	114.00
2	D	354	HEM	C3B-C4B-NB	-7.28	108.79	114.00
2	C	354	HEM	C3B-C4B-NB	-7.21	108.84	114.00
2	E	354	HEM	C3B-C4B-NB	-7.15	108.89	114.00
3	E	355	BXO	O1-N1-C1	6.71	118.26	111.66
2	F	354	HEM	C4D-ND-C1D	6.58	111.89	105.16
3	G	355	BXO	O1-N1-C1	6.57	118.12	111.66
2	D	354	HEM	C4D-ND-C1D	6.23	111.54	105.16
2	B	354	HEM	C4D-ND-C1D	6.19	111.50	105.16
3	B	355	BXO	O1-N1-C1	6.18	117.74	111.66
2	G	354	HEM	C4D-ND-C1D	6.08	111.39	105.16
2	E	354	HEM	C4D-ND-C1D	6.05	111.36	105.16
2	H	354	HEM	C4D-ND-C1D	6.00	111.30	105.16
2	A	354	HEM	C4D-ND-C1D	5.97	111.28	105.16
2	C	354	HEM	C4D-ND-C1D	5.97	111.27	105.16
3	D	355	BXO	O1-N1-C1	5.92	117.48	111.66
3	H	355	BXO	O1-N1-C1	5.83	117.39	111.66
2	C	354	HEM	CBA-CAA-C2A	-5.61	102.81	112.69
2	A	354	HEM	CBA-CAA-C2A	-5.57	102.88	112.69
2	E	354	HEM	CBA-CAA-C2A	-5.50	102.99	112.69
3	F	355	BXO	O1-N1-C1	5.48	117.05	111.66
2	G	354	HEM	CBA-CAA-C2A	-5.46	103.07	112.69
2	F	354	HEM	C2D-C1D-ND	-4.29	107.87	112.93
2	B	354	HEM	C2D-C1D-ND	-4.11	108.08	112.93
2	H	354	HEM	C2D-C1D-ND	-4.00	108.20	112.93
2	D	354	HEM	C2D-C1D-ND	-3.98	108.23	112.93
2	A	354	HEM	C4A-CHB-C1B	-3.59	122.74	127.47
2	C	354	HEM	C4A-CHB-C1B	-3.47	122.91	127.47
2	G	354	HEM	C4A-CHB-C1B	-3.44	122.94	127.47
2	F	354	HEM	CAD-C3D-C4D	3.43	130.69	124.53
2	E	354	HEM	C2D-C1D-ND	-3.41	108.90	112.93
2	G	354	HEM	C2D-C1D-ND	-3.39	108.92	112.93
2	D	354	HEM	CAD-C3D-C4D	3.38	130.60	124.53
2	C	354	HEM	C2D-C1D-ND	-3.35	108.97	112.93
2	A	354	HEM	C2D-C1D-ND	-3.35	108.98	112.93
2	E	354	HEM	C4A-CHB-C1B	-3.30	123.12	127.47
2	H	354	HEM	CAD-C3D-C4D	3.29	130.44	124.53
2	B	354	HEM	CAD-C3D-C4D	3.26	130.40	124.53
2	H	354	HEM	C4A-CHB-C1B	-3.02	123.50	127.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	354	HEM	C4A-CHB-C1B	-2.99	123.53	127.47
2	B	354	HEM	C4C-NC-C1C	2.83	108.47	105.53
2	F	354	HEM	C4A-CHB-C1B	-2.75	123.85	127.47
2	A	354	HEM	C1B-NB-C4B	2.74	107.96	105.16
2	B	354	HEM	C4A-CHB-C1B	-2.73	123.88	127.47
2	E	354	HEM	C3A-C4A-NA	-2.66	107.41	109.41
2	G	354	HEM	C1B-NB-C4B	2.63	107.85	105.16
2	H	354	HEM	CHC-C4B-NB	2.56	126.71	124.58
2	A	354	HEM	C3A-C4A-NA	-2.51	107.51	109.41
2	E	354	HEM	C1B-NB-C4B	2.51	107.72	105.16
2	H	354	HEM	C4C-NC-C1C	2.49	108.12	105.53
2	F	354	HEM	C4C-NC-C1C	2.49	108.12	105.53
3	D	355	BXO	C2-C1-N1	2.45	126.43	120.41
3	F	355	BXO	C2-C1-N1	2.43	126.38	120.41
2	B	354	HEM	C3A-C4A-NA	-2.43	107.58	109.41
2	C	354	HEM	C1B-NB-C4B	2.41	107.63	105.16
2	E	354	HEM	C4C-NC-C1C	2.41	108.04	105.53
2	C	354	HEM	C3A-C4A-NA	-2.40	107.60	109.41
2	F	354	HEM	C1B-NB-C4B	2.40	107.62	105.16
2	F	354	HEM	CHC-C4B-NB	2.39	126.57	124.58
3	H	355	BXO	C2-C1-N1	2.37	126.23	120.41
3	B	355	BXO	C2-C1-N1	2.36	126.22	120.41
2	A	354	HEM	C4C-NC-C1C	2.36	107.98	105.53
2	G	354	HEM	C4C-NC-C1C	2.35	107.98	105.53
2	D	354	HEM	C4C-NC-C1C	2.35	107.97	105.53
2	C	354	HEM	CAD-C3D-C4D	2.35	128.75	124.53
2	E	354	HEM	CAD-C3D-C4D	2.34	128.74	124.53
2	G	354	HEM	C3A-C4A-NA	-2.31	107.67	109.41
2	A	354	HEM	CAD-C3D-C4D	2.26	128.59	124.53
2	D	354	HEM	C3A-C4A-NA	-2.24	107.72	109.41
2	D	354	HEM	C1B-NB-C4B	2.23	107.45	105.16
2	C	354	HEM	C4C-NC-C1C	2.23	107.86	105.53
2	H	354	HEM	C3A-C4A-NA	-2.21	107.74	109.41
3	E	355	BXO	C2-C1-N1	2.18	125.78	120.41
2	H	354	HEM	C1B-NB-C4B	2.18	107.39	105.16
2	B	354	HEM	CMA-C3A-C4A	-2.12	125.36	128.62
2	B	354	HEM	C1B-NB-C4B	2.12	107.33	105.16
2	D	354	HEM	CMA-C3A-C4A	-2.12	125.37	128.62
2	D	354	HEM	CAA-CBA-CGA	-2.11	106.69	113.47
2	G	354	HEM	CAD-C3D-C4D	2.11	128.31	124.53
2	H	354	HEM	CMA-C3A-C4A	-2.10	125.39	128.62
2	B	354	HEM	CHC-C4B-NB	2.08	126.32	124.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	355	BXO	C2-C1-N1	2.08	125.54	120.41
2	D	354	HEM	CHC-C4B-NB	2.04	126.28	124.58
2	F	354	HEM	CMA-C3A-C4A	-2.03	125.49	128.62
3	C	355	BXO	C2-C1-N1	2.01	125.36	120.41
2	F	354	HEM	C3A-C4A-NA	-2.01	107.89	109.41

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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 ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	350/373 (93%)	0.29	14 (4%) 36 37	13, 25, 39, 55	0
1	B	359/373 (96%)	0.59	37 (10%) 7 6	15, 31, 57, 67	0
1	C	352/373 (94%)	0.30	22 (6%) 19 19	14, 25, 40, 55	0
1	D	360/373 (96%)	0.55	30 (8%) 11 10	15, 31, 57, 68	0
1	E	351/373 (94%)	0.30	18 (5%) 27 27	14, 25, 40, 55	0
1	F	359/373 (96%)	0.59	39 (10%) 6 6	15, 31, 57, 68	0
1	G	353/373 (94%)	0.30	20 (5%) 23 23	13, 25, 41, 55	0
1	H	359/373 (96%)	0.54	34 (9%) 8 8	15, 31, 57, 68	0
All	All	2843/2984 (95%)	0.43	214 (7%) 14 14	13, 26, 55, 68	0

All (214) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	353	HIS	7.6
1	H	350	ILE	6.6
1	G	351	ALA	6.5
1	B	126	ASP	5.9
1	F	350	ILE	5.9
1	D	350	ILE	5.8
1	F	125	SER	5.7
1	B	201	ALA	5.7
1	F	201	ALA	5.5
1	B	350	ILE	5.4
1	E	124	LEU	5.4
1	C	351	ALA	5.1
1	F	126	ASP	5.1
1	H	75	GLY	4.9
1	H	201	ALA	4.7
1	G	121	GLU	4.7

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Mol	Chain	Res	Type	RSRZ
1	B	103	SER	4.6
1	F	121	GLU	4.6
1	A	124	LEU	4.6
1	E	12	PRO	4.5
1	H	202	VAL	4.5
1	A	125	SER	4.5
1	E	125	SER	4.4
1	B	125	SER	4.4
1	D	112	THR	4.4
1	G	12	PRO	4.4
1	A	12	PRO	4.2
1	H	121	GLU	4.2
1	F	127	GLY	4.2
1	B	112	THR	4.2
1	F	110	THR	4.0
1	C	12	PRO	4.0
1	B	194	ARG	4.0
1	D	-8	SER	4.0
1	E	351	ALA	3.9
1	F	122	ASP	3.9
1	D	106	HIS	3.9
1	C	125	SER	3.9
1	F	119	GLU	3.8
1	D	125	SER	3.7
1	F	12	PRO	3.7
1	H	122	ASP	3.7
1	D	349	THR	3.7
1	C	311	ALA	3.6
1	H	103	SER	3.6
1	F	106	HIS	3.5
1	C	124	LEU	3.5
1	B	121	GLU	3.4
1	D	119	GLU	3.4
1	D	113	PRO	3.4
1	H	194	ARG	3.4
1	F	74	ASP	3.4
1	H	349	THR	3.4
1	F	52	ASP	3.4
1	H	106	HIS	3.4
1	A	121	GLU	3.4
1	F	349	THR	3.3
1	D	122	ASP	3.3

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Mol	Chain	Res	Type	RSRZ
1	D	110	THR	3.3
1	E	21	ASP	3.3
1	B	122	ASP	3.2
1	H	127	GLY	3.2
1	G	125	SER	3.2
1	E	51	ARG	3.2
1	E	121	GLU	3.2
1	E	311	ALA	3.2
1	G	124	LEU	3.2
1	F	112	THR	3.1
1	H	125	SER	3.1
1	A	123	ARG	3.1
1	H	126	ASP	3.1
1	B	63	MET	3.0
1	A	112	THR	3.0
1	F	196	ILE	3.0
1	F	328	ALA	3.0
1	C	318	LEU	3.0
1	B	119	GLU	3.0
1	B	12	PRO	3.0
1	B	269	ILE	3.0
1	D	201	ALA	2.9
1	D	126	ASP	2.9
1	G	126	ASP	2.9
1	F	336	ASN	2.9
1	H	53	GLU	2.9
1	D	121	GLU	2.9
1	B	73	PRO	2.8
1	D	351	ALA	2.8
1	A	36	ASP	2.8
1	A	311	ALA	2.8
1	B	205	ARG	2.8
1	B	53	GLU	2.8
1	B	72	LEU	2.8
1	F	103	SER	2.8
1	C	192	GLU	2.8
1	F	51	ARG	2.8
1	B	349	THR	2.8
1	A	54	ASP	2.7
1	D	36	ASP	2.7
1	H	151	ASP	2.7
1	H	112	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	36	ASP	2.7
1	E	53	GLU	2.7
1	B	110	THR	2.7
1	B	54	ASP	2.7
1	F	54	ASP	2.7
1	F	-8	SER	2.7
1	D	75	GLY	2.6
1	A	21	ASP	2.6
1	E	350	ILE	2.6
1	D	12	PRO	2.6
1	E	112	THR	2.6
1	C	57	PRO	2.6
1	H	110	THR	2.6
1	G	1	MET	2.5
1	F	151	ASP	2.5
1	H	73	PRO	2.5
1	H	200	PRO	2.5
1	D	73	PRO	2.5
1	B	336	ASN	2.5
1	C	121	GLU	2.5
1	D	150	GLU	2.5
1	E	126	ASP	2.5
1	D	194	ARG	2.5
1	E	1	MET	2.5
1	F	348	VAL	2.4
1	G	311	ALA	2.4
1	F	73	PRO	2.4
1	B	71	ASP	2.4
1	D	35	ASP	2.4
1	E	192	GLU	2.4
1	F	75	GLY	2.4
1	F	113	PRO	2.4
1	H	52	ASP	2.4
1	B	348	VAL	2.4
1	B	127	GLY	2.4
1	B	52	ASP	2.4
1	G	54	ASP	2.4
1	H	61	GLN	2.4
1	D	124	LEU	2.4
1	H	107	ARG	2.4
1	H	119	GLU	2.4
1	B	75	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	G	36	ASP	2.4
1	F	7	GLU	2.4
1	B	116	SER	2.4
1	B	190	SER	2.4
1	H	54	ASP	2.4
1	F	202	VAL	2.3
1	G	123	ARG	2.3
1	G	352	GLU	2.3
1	F	53	GLU	2.3
1	B	-8	SER	2.3
1	D	7	GLU	2.3
1	F	111	SER	2.3
1	B	128	LEU	2.3
1	B	328	ALA	2.3
1	H	36	ASP	2.3
1	A	318	LEU	2.3
1	F	194	ARG	2.3
1	F	118	TRP	2.3
1	H	254	PRO	2.3
1	B	202	VAL	2.3
1	G	22	THR	2.3
1	B	7	GLU	2.3
1	E	150	GLU	2.3
1	H	124	LEU	2.2
1	D	131	PHE	2.2
1	E	194	ARG	2.2
1	F	205	ARG	2.2
1	H	116	SER	2.2
1	D	72	LEU	2.2
1	H	35	ASP	2.2
1	G	51	ARG	2.2
1	G	21	ASP	2.2
1	C	126	ASP	2.2
1	C	1	MET	2.2
1	D	14	THR	2.2
1	C	54	ASP	2.2
1	F	192	GLU	2.2
1	F	0	HIS	2.2
1	C	71	ASP	2.2
1	C	122	ASP	2.2
1	A	53	GLU	2.2
1	D	53	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	G	318	LEU	2.1
1	A	194	ARG	2.1
1	F	347	ALA	2.1
1	F	36	ASP	2.1
1	H	-8	SER	2.1
1	H	88	GLN	2.1
1	C	123	ARG	2.1
1	D	118	TRP	2.1
1	A	126	ASP	2.1
1	B	151	ASP	2.1
1	C	350	ILE	2.1
1	B	272	ASN	2.1
1	G	116	SER	2.1
1	H	196	ILE	2.1
1	C	11	CYS	2.1
1	C	110	THR	2.1
1	E	122	ASP	2.1
1	G	53	GLU	2.1
1	E	308	ARG	2.1
1	B	36	ASP	2.1
1	D	54	ASP	2.1
1	C	51	ARG	2.0
1	C	7	GLU	2.0
1	B	48	VAL	2.0
1	G	192	GLU	2.0
1	D	327	ALA	2.0
1	C	53	GLU	2.0
1	F	35	ASP	2.0
1	H	74	ASP	2.0
1	H	269	ILE	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	BXO	B	355	6/6	0.19	1.48	25,26,27,28	0
3	BXO	D	355	6/6	0.17	0.56	25,26,27,28	0
3	BXO	F	355	6/6	0.16	0.35	25,26,27,28	0
2	HEM	D	354	43/43	0.16	0.05	12,13,17,17	0
2	HEM	H	354	43/43	0.14	-0.05	12,13,17,17	0
3	BXO	G	355	6/6	0.17	-0.07	18,20,21,21	0
3	BXO	A	355	6/6	0.17	-0.22	18,20,21,21	0
2	HEM	F	354	43/43	0.14	-0.29	12,13,17,17	0
2	HEM	E	354	43/43	0.14	-0.39	13,15,18,18	0
3	BXO	H	355	6/6	0.14	-0.49	25,26,27,28	0
2	HEM	B	354	43/43	0.13	-0.51	12,13,17,17	0
3	BXO	E	355	6/6	0.14	-0.54	18,20,21,21	0
3	BXO	C	355	6/6	0.14	-0.69	18,20,21,21	0
2	HEM	A	354	43/43	0.13	-0.86	13,15,18,18	0
2	HEM	G	354	43/43	0.12	-0.86	13,15,18,18	0
2	HEM	C	354	43/43	0.13	-0.95	13,15,18,18	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.