



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:56 PM GMT

PDB ID : 4AUB
Title : the complex Structure of the bacterial aldo-keto reductase AKR14A1 with NADP and citrate
Authors : Zhu, X.; Ellis, E.M.; Lapthorn, A.
Deposited on : 2012-05-16
Resolution : 2.05 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

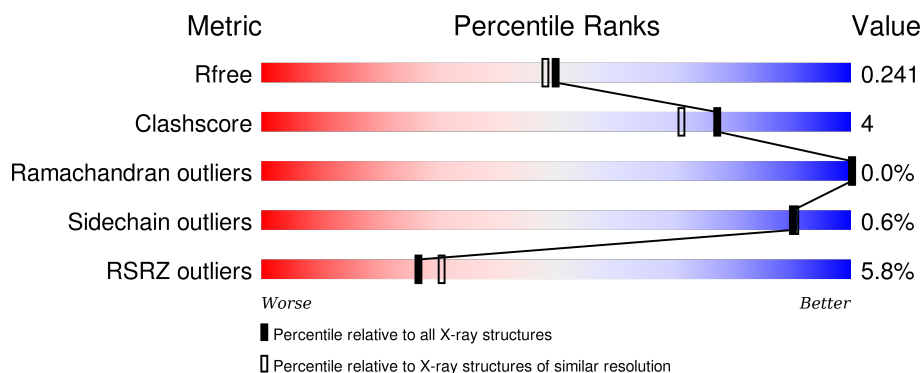
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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}	91344	1192 (2.04-2.04)
Clashscore	102246	1269 (2.04-2.04)
Ramachandran outliers	100387	1258 (2.04-2.04)
Sidechain outliers	100360	1258 (2.04-2.04)
RSRZ outliers	91569	1194 (2.04-2.04)

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 ≥ 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.

Mol	Chain	Length	Quality of chain
1	A	366	<div> <div>3%</div> <div>78%</div> <div>7%</div> <div>15%</div> </div>
1	B	366	<div> <div>9%</div> <div>82%</div> <div>9%</div> <div>8%</div> </div>
1	C	366	<div> <div>7%</div> <div>83%</div> <div>7%</div> <div>10%</div> </div>
1	D	366	<div> <div>81%</div> <div>16%</div> </div>
1	E	366	<div> <div>8%</div> <div>75%</div> <div>5%</div> <div>19%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	366	
1	G	366	
1	H	366	

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
2	NAP	A	500	-	-	-	X
2	NAP	F	500	-	-	-	X
3	FLC	A	501	-	-	-	X
3	FLC	C	501	-	-	-	X
3	FLC	D	501	-	-	-	X
3	FLC	F	501	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 21483 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 ALDO-KETO REDUCTASE AKR14A1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	312	Total	C	N	O	S	0	2	0
			2472	1561	436	465	10			
1	B	335	Total	C	N	O	S	0	1	0
			2655	1673	470	500	12			
1	C	330	Total	C	N	O	S	0	1	0
			2608	1646	457	493	12			
1	D	308	Total	C	N	O	S	0	0	0
			2430	1535	427	458	10			
1	E	297	Total	C	N	O	S	0	0	0
			2345	1484	411	440	10			
1	F	298	Total	C	N	O	S	0	0	0
			2359	1493	416	440	10			
1	G	335	Total	C	N	O	S	0	3	0
			2665	1681	471	501	12			
1	H	297	Total	C	N	O	S	0	0	0
			2348	1485	411	442	10			

There are 160 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP Q46851
A	-18	GLY	-	EXPRESSION TAG	UNP Q46851
A	-17	SER	-	EXPRESSION TAG	UNP Q46851
A	-16	SER	-	EXPRESSION TAG	UNP Q46851
A	-15	HIS	-	EXPRESSION TAG	UNP Q46851
A	-14	HIS	-	EXPRESSION TAG	UNP Q46851
A	-13	HIS	-	EXPRESSION TAG	UNP Q46851
A	-12	HIS	-	EXPRESSION TAG	UNP Q46851
A	-11	HIS	-	EXPRESSION TAG	UNP Q46851
A	-10	HIS	-	EXPRESSION TAG	UNP Q46851
A	-9	SER	-	EXPRESSION TAG	UNP Q46851
A	-8	SER	-	EXPRESSION TAG	UNP Q46851
A	-7	GLY	-	EXPRESSION TAG	UNP Q46851

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	LEU	-	EXPRESSION TAG	UNP Q46851
A	-5	VAL	-	EXPRESSION TAG	UNP Q46851
A	-4	PRO	-	EXPRESSION TAG	UNP Q46851
A	-3	ARG	-	EXPRESSION TAG	UNP Q46851
A	-2	GLY	-	EXPRESSION TAG	UNP Q46851
A	-1	SER	-	EXPRESSION TAG	UNP Q46851
A	0	HIS	-	EXPRESSION TAG	UNP Q46851
B	-19	MET	-	EXPRESSION TAG	UNP Q46851
B	-18	GLY	-	EXPRESSION TAG	UNP Q46851
B	-17	SER	-	EXPRESSION TAG	UNP Q46851
B	-16	SER	-	EXPRESSION TAG	UNP Q46851
B	-15	HIS	-	EXPRESSION TAG	UNP Q46851
B	-14	HIS	-	EXPRESSION TAG	UNP Q46851
B	-13	HIS	-	EXPRESSION TAG	UNP Q46851
B	-12	HIS	-	EXPRESSION TAG	UNP Q46851
B	-11	HIS	-	EXPRESSION TAG	UNP Q46851
B	-10	HIS	-	EXPRESSION TAG	UNP Q46851
B	-9	SER	-	EXPRESSION TAG	UNP Q46851
B	-8	SER	-	EXPRESSION TAG	UNP Q46851
B	-7	GLY	-	EXPRESSION TAG	UNP Q46851
B	-6	LEU	-	EXPRESSION TAG	UNP Q46851
B	-5	VAL	-	EXPRESSION TAG	UNP Q46851
B	-4	PRO	-	EXPRESSION TAG	UNP Q46851
B	-3	ARG	-	EXPRESSION TAG	UNP Q46851
B	-2	GLY	-	EXPRESSION TAG	UNP Q46851
B	-1	SER	-	EXPRESSION TAG	UNP Q46851
B	0	HIS	-	EXPRESSION TAG	UNP Q46851
C	-19	MET	-	EXPRESSION TAG	UNP Q46851
C	-18	GLY	-	EXPRESSION TAG	UNP Q46851
C	-17	SER	-	EXPRESSION TAG	UNP Q46851
C	-16	SER	-	EXPRESSION TAG	UNP Q46851
C	-15	HIS	-	EXPRESSION TAG	UNP Q46851
C	-14	HIS	-	EXPRESSION TAG	UNP Q46851
C	-13	HIS	-	EXPRESSION TAG	UNP Q46851
C	-12	HIS	-	EXPRESSION TAG	UNP Q46851
C	-11	HIS	-	EXPRESSION TAG	UNP Q46851
C	-10	HIS	-	EXPRESSION TAG	UNP Q46851
C	-9	SER	-	EXPRESSION TAG	UNP Q46851
C	-8	SER	-	EXPRESSION TAG	UNP Q46851
C	-7	GLY	-	EXPRESSION TAG	UNP Q46851
C	-6	LEU	-	EXPRESSION TAG	UNP Q46851
C	-5	VAL	-	EXPRESSION TAG	UNP Q46851

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-4	PRO	-	EXPRESSION TAG	UNP Q46851
C	-3	ARG	-	EXPRESSION TAG	UNP Q46851
C	-2	GLY	-	EXPRESSION TAG	UNP Q46851
C	-1	SER	-	EXPRESSION TAG	UNP Q46851
C	0	HIS	-	EXPRESSION TAG	UNP Q46851
D	-19	MET	-	EXPRESSION TAG	UNP Q46851
D	-18	GLY	-	EXPRESSION TAG	UNP Q46851
D	-17	SER	-	EXPRESSION TAG	UNP Q46851
D	-16	SER	-	EXPRESSION TAG	UNP Q46851
D	-15	HIS	-	EXPRESSION TAG	UNP Q46851
D	-14	HIS	-	EXPRESSION TAG	UNP Q46851
D	-13	HIS	-	EXPRESSION TAG	UNP Q46851
D	-12	HIS	-	EXPRESSION TAG	UNP Q46851
D	-11	HIS	-	EXPRESSION TAG	UNP Q46851
D	-10	HIS	-	EXPRESSION TAG	UNP Q46851
D	-9	SER	-	EXPRESSION TAG	UNP Q46851
D	-8	SER	-	EXPRESSION TAG	UNP Q46851
D	-7	GLY	-	EXPRESSION TAG	UNP Q46851
D	-6	LEU	-	EXPRESSION TAG	UNP Q46851
D	-5	VAL	-	EXPRESSION TAG	UNP Q46851
D	-4	PRO	-	EXPRESSION TAG	UNP Q46851
D	-3	ARG	-	EXPRESSION TAG	UNP Q46851
D	-2	GLY	-	EXPRESSION TAG	UNP Q46851
D	-1	SER	-	EXPRESSION TAG	UNP Q46851
D	0	HIS	-	EXPRESSION TAG	UNP Q46851
E	-19	MET	-	EXPRESSION TAG	UNP Q46851
E	-18	GLY	-	EXPRESSION TAG	UNP Q46851
E	-17	SER	-	EXPRESSION TAG	UNP Q46851
E	-16	SER	-	EXPRESSION TAG	UNP Q46851
E	-15	HIS	-	EXPRESSION TAG	UNP Q46851
E	-14	HIS	-	EXPRESSION TAG	UNP Q46851
E	-13	HIS	-	EXPRESSION TAG	UNP Q46851
E	-12	HIS	-	EXPRESSION TAG	UNP Q46851
E	-11	HIS	-	EXPRESSION TAG	UNP Q46851
E	-10	HIS	-	EXPRESSION TAG	UNP Q46851
E	-9	SER	-	EXPRESSION TAG	UNP Q46851
E	-8	SER	-	EXPRESSION TAG	UNP Q46851
E	-7	GLY	-	EXPRESSION TAG	UNP Q46851
E	-6	LEU	-	EXPRESSION TAG	UNP Q46851
E	-5	VAL	-	EXPRESSION TAG	UNP Q46851
E	-4	PRO	-	EXPRESSION TAG	UNP Q46851
E	-3	ARG	-	EXPRESSION TAG	UNP Q46851

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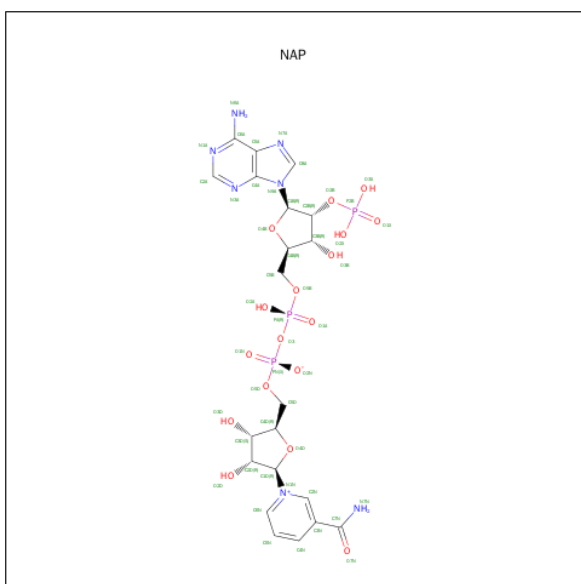
Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	GLY	-	EXPRESSION TAG	UNP Q46851
E	-1	SER	-	EXPRESSION TAG	UNP Q46851
E	0	HIS	-	EXPRESSION TAG	UNP Q46851
F	-19	MET	-	EXPRESSION TAG	UNP Q46851
F	-18	GLY	-	EXPRESSION TAG	UNP Q46851
F	-17	SER	-	EXPRESSION TAG	UNP Q46851
F	-16	SER	-	EXPRESSION TAG	UNP Q46851
F	-15	HIS	-	EXPRESSION TAG	UNP Q46851
F	-14	HIS	-	EXPRESSION TAG	UNP Q46851
F	-13	HIS	-	EXPRESSION TAG	UNP Q46851
F	-12	HIS	-	EXPRESSION TAG	UNP Q46851
F	-11	HIS	-	EXPRESSION TAG	UNP Q46851
F	-10	HIS	-	EXPRESSION TAG	UNP Q46851
F	-9	SER	-	EXPRESSION TAG	UNP Q46851
F	-8	SER	-	EXPRESSION TAG	UNP Q46851
F	-7	GLY	-	EXPRESSION TAG	UNP Q46851
F	-6	LEU	-	EXPRESSION TAG	UNP Q46851
F	-5	VAL	-	EXPRESSION TAG	UNP Q46851
F	-4	PRO	-	EXPRESSION TAG	UNP Q46851
F	-3	ARG	-	EXPRESSION TAG	UNP Q46851
F	-2	GLY	-	EXPRESSION TAG	UNP Q46851
F	-1	SER	-	EXPRESSION TAG	UNP Q46851
F	0	HIS	-	EXPRESSION TAG	UNP Q46851
G	-19	MET	-	EXPRESSION TAG	UNP Q46851
G	-18	GLY	-	EXPRESSION TAG	UNP Q46851
G	-17	SER	-	EXPRESSION TAG	UNP Q46851
G	-16	SER	-	EXPRESSION TAG	UNP Q46851
G	-15	HIS	-	EXPRESSION TAG	UNP Q46851
G	-14	HIS	-	EXPRESSION TAG	UNP Q46851
G	-13	HIS	-	EXPRESSION TAG	UNP Q46851
G	-12	HIS	-	EXPRESSION TAG	UNP Q46851
G	-11	HIS	-	EXPRESSION TAG	UNP Q46851
G	-10	HIS	-	EXPRESSION TAG	UNP Q46851
G	-9	SER	-	EXPRESSION TAG	UNP Q46851
G	-8	SER	-	EXPRESSION TAG	UNP Q46851
G	-7	GLY	-	EXPRESSION TAG	UNP Q46851
G	-6	LEU	-	EXPRESSION TAG	UNP Q46851
G	-5	VAL	-	EXPRESSION TAG	UNP Q46851
G	-4	PRO	-	EXPRESSION TAG	UNP Q46851
G	-3	ARG	-	EXPRESSION TAG	UNP Q46851
G	-2	GLY	-	EXPRESSION TAG	UNP Q46851
G	-1	SER	-	EXPRESSION TAG	UNP Q46851

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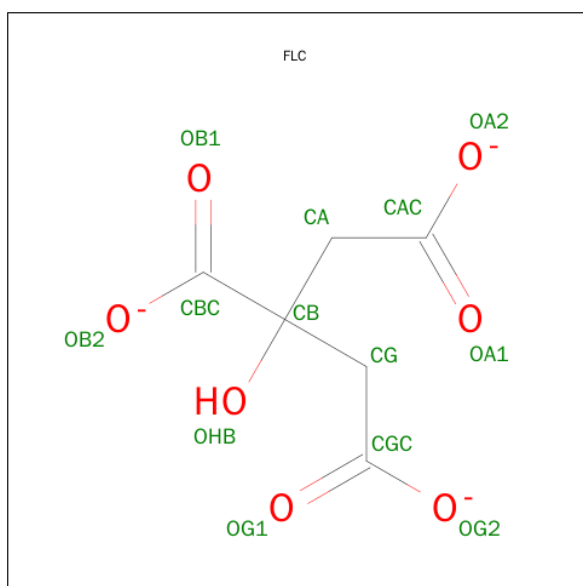
Chain	Residue	Modelled	Actual	Comment	Reference
G	0	HIS	-	EXPRESSION TAG	UNP Q46851
H	-19	MET	-	EXPRESSION TAG	UNP Q46851
H	-18	GLY	-	EXPRESSION TAG	UNP Q46851
H	-17	SER	-	EXPRESSION TAG	UNP Q46851
H	-16	SER	-	EXPRESSION TAG	UNP Q46851
H	-15	HIS	-	EXPRESSION TAG	UNP Q46851
H	-14	HIS	-	EXPRESSION TAG	UNP Q46851
H	-13	HIS	-	EXPRESSION TAG	UNP Q46851
H	-12	HIS	-	EXPRESSION TAG	UNP Q46851
H	-11	HIS	-	EXPRESSION TAG	UNP Q46851
H	-10	HIS	-	EXPRESSION TAG	UNP Q46851
H	-9	SER	-	EXPRESSION TAG	UNP Q46851
H	-8	SER	-	EXPRESSION TAG	UNP Q46851
H	-7	GLY	-	EXPRESSION TAG	UNP Q46851
H	-6	LEU	-	EXPRESSION TAG	UNP Q46851
H	-5	VAL	-	EXPRESSION TAG	UNP Q46851
H	-4	PRO	-	EXPRESSION TAG	UNP Q46851
H	-3	ARG	-	EXPRESSION TAG	UNP Q46851
H	-2	GLY	-	EXPRESSION TAG	UNP Q46851
H	-1	SER	-	EXPRESSION TAG	UNP Q46851
H	0	HIS	-	EXPRESSION TAG	UNP Q46851

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	E	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	F	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	G	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	H	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is CITRATE ANION (three-letter code: FLC) (formula: $C_6H_5O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			13	6	7		
3	B	1	Total	C	O	0	0
			13	6	7		
3	C	1	Total	C	O	0	0
			13	6	7		
3	D	1	Total	C	O	0	0
			13	6	7		
3	E	1	Total	C	O	0	0
			13	6	7		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	F	1	Total	C	O	0	0
			13	6	7		
3	G	1	Total	C	O	0	0
			13	6	7		
3	H	1	Total	C	O	0	0
			13	6	7		

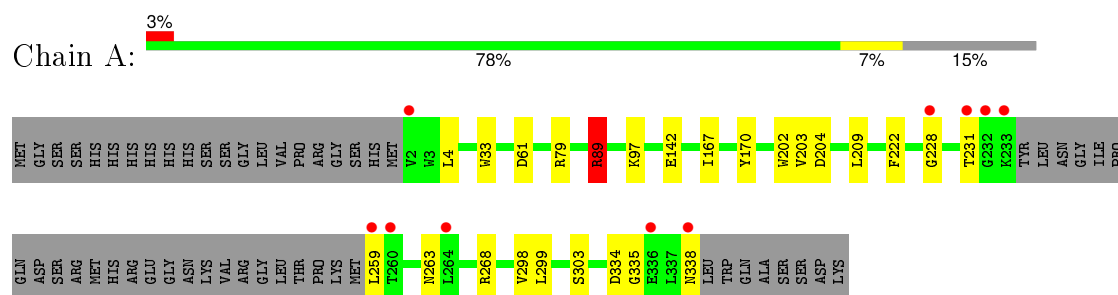
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	181	Total	O	0	0
			181	181		
4	B	155	Total	O	0	0
			155	155		
4	C	142	Total	O	0	0
			142	142		
4	D	161	Total	O	0	0
			161	161		
4	E	121	Total	O	0	0
			121	121		
4	F	156	Total	O	0	0
			156	156		
4	G	132	Total	O	0	0
			132	132		
4	H	113	Total	O	0	0
			113	113		

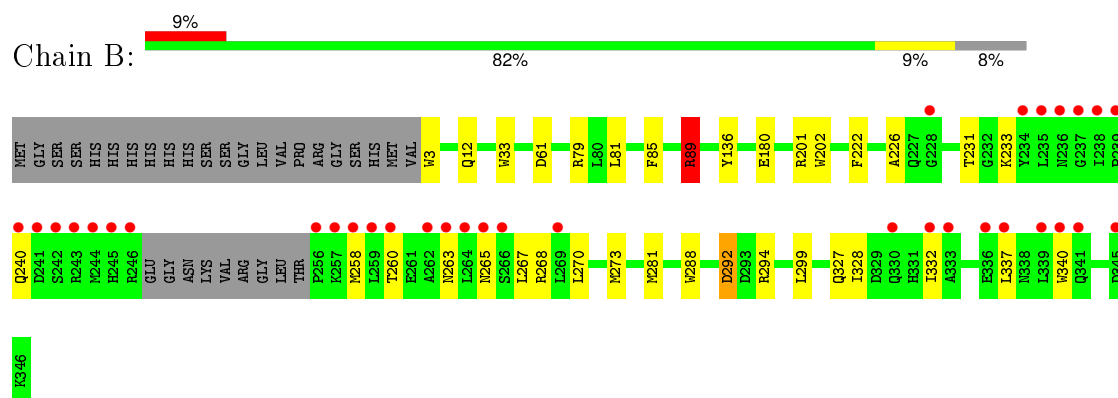
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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.

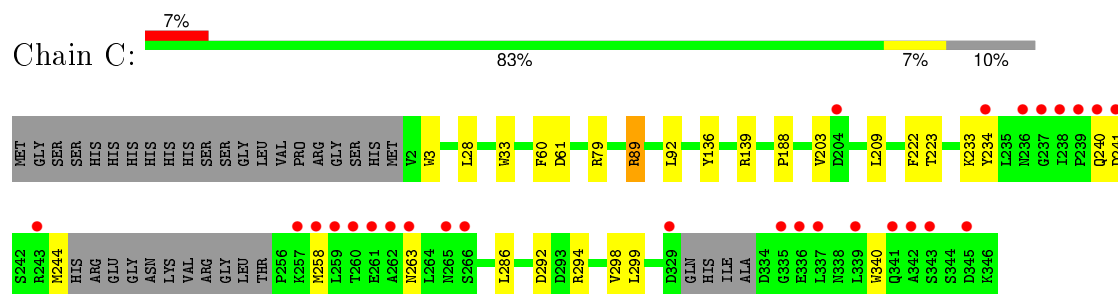
- Molecule 1: ALDO-KETO REDUCTASE AKR14A1



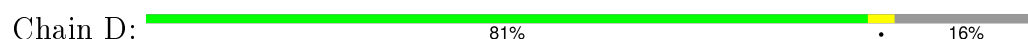
- Molecule 1: ALDO-KETO REDUCTASE AKR14A1

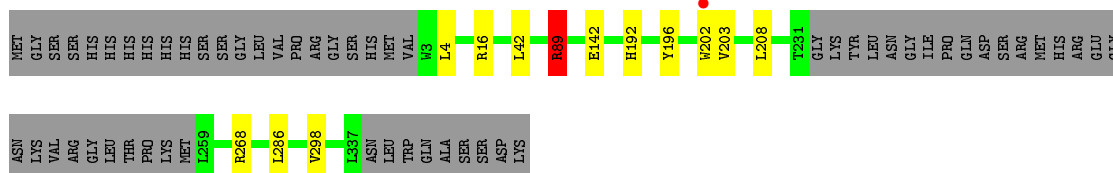


- Molecule 1: ALDO-KETO REDUCTASE AKR14A1

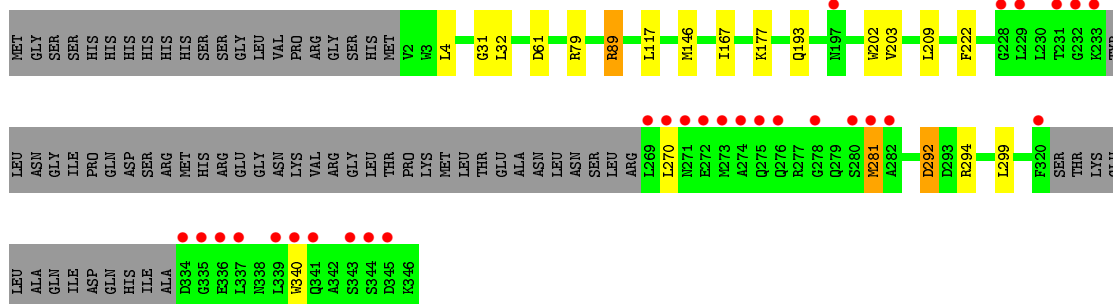


- Molecule 1: ALDO-KETO REDUCTASE AKR14A1

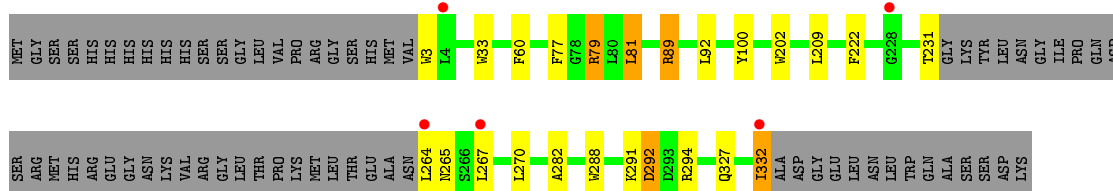




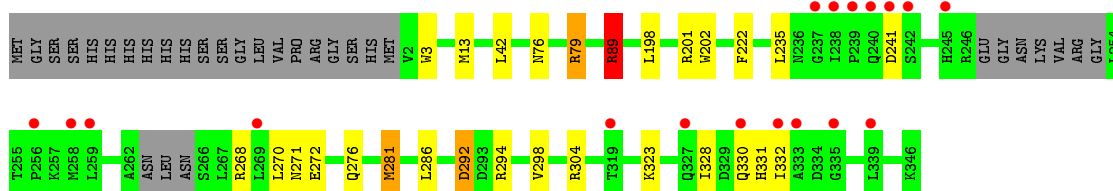
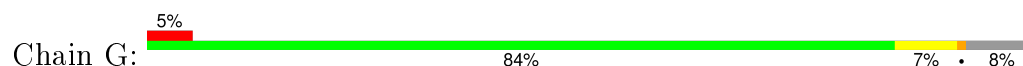
• Molecule 1: ALDO-KETO REDUCTASE AKR14A1



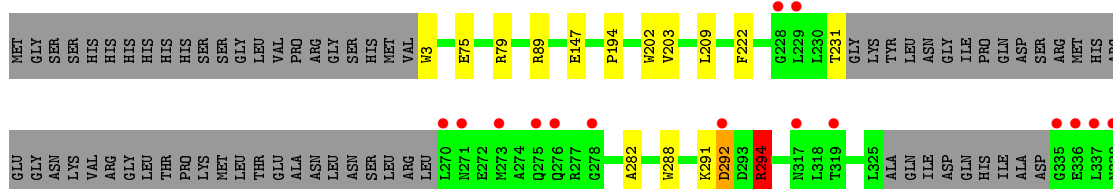
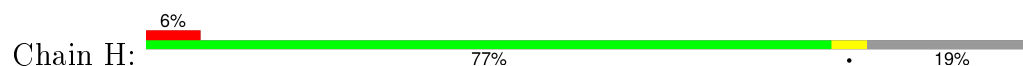
• Molecule 1: ALDO-KETO REDUCTASE AKR14A1

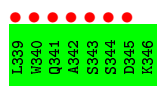


• Molecule 1: ALDO-KETO REDUCTASE AKR14A1



• Molecule 1: ALDO-KETO REDUCTASE AKR14A1





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	91.81Å 191.09Å 97.30Å 90.00° 105.97° 90.00°	Depositor
Resolution (Å)	95.35 – 2.05 48.87 – 2.05	Depositor EDS
% Data completeness (in resolution range)	92.2 (95.35-2.05) 92.2 (48.87-2.05)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.34 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.202 , 0.238 0.203 , 0.241	Depositor DCC
R_{free} test set	10115 reflections (5.76%)	DCC
Wilson B-factor (Å ²)	25.0	Xtriage
Anisotropy	0.080	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 185058 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	21483	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.57% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.97	2/2528 (0.1%)	0.88	2/3421 (0.1%)
1	B	0.93	5/2716 (0.2%)	0.84	2/3672 (0.1%)
1	C	0.94	2/2664 (0.1%)	0.85	2/3601 (0.1%)
1	D	1.01	1/2480 (0.0%)	0.90	4/3358 (0.1%)
1	E	0.87	2/2395 (0.1%)	0.86	3/3239 (0.1%)
1	F	0.92	4/2409 (0.2%)	0.87	3/3261 (0.1%)
1	G	0.93	2/2733 (0.1%)	0.92	7/3694 (0.2%)
1	H	0.89	4/2398 (0.2%)	0.85	3/3243 (0.1%)
All	All	0.93	22/20323 (0.1%)	0.87	26/27489 (0.1%)

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
1	B	0	2
1	C	0	2
1	D	0	1
1	E	0	1
1	F	0	1
1	G	0	1
All	All	0	9

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	202	TRP	CD2-CE2	8.72	1.51	1.41
1	B	202	TRP	CD2-CE2	6.96	1.49	1.41
1	H	3	TRP	CD2-CE2	6.76	1.49	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	202	TRP	CD2-CE2	6.49	1.49	1.41
1	B	33	TRP	CD2-CE2	6.39	1.49	1.41
1	B	3	TRP	CD2-CE2	6.32	1.49	1.41
1	E	202	TRP	CD2-CE2	6.20	1.48	1.41
1	H	202	TRP	CD2-CE2	6.05	1.48	1.41
1	F	33	TRP	CD2-CE2	6.05	1.48	1.41
1	E	340	TRP	CD2-CE2	6.03	1.48	1.41
1	G	3	TRP	CD2-CE2	6.00	1.48	1.41
1	B	340	TRP	CD2-CE2	5.83	1.48	1.41
1	F	288	TRP	CD2-CE2	5.77	1.48	1.41
1	G	202	TRP	CD2-CE2	5.75	1.48	1.41
1	F	3	TRP	CD2-CE2	5.75	1.48	1.41
1	A	202	TRP	CG-CD1	5.68	1.44	1.36
1	C	340	TRP	CD2-CE2	5.48	1.48	1.41
1	C	3	TRP	CD2-CE2	5.45	1.47	1.41
1	B	288	TRP	CD2-CE2	5.39	1.47	1.41
1	F	202	TRP	CD2-CE2	5.33	1.47	1.41
1	H	288	TRP	CD2-CE2	5.32	1.47	1.41
1	H	147	GLU	CG-CD	5.13	1.59	1.51

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	292	ASP	CB-CG-OD1	9.55	126.90	118.30
1	B	292	ASP	CB-CG-OD1	8.81	126.23	118.30
1	E	292	ASP	CB-CG-OD1	8.73	126.16	118.30
1	G	79	ARG	NE-CZ-NH1	8.16	124.38	120.30
1	E	89	ARG	CG-CD-NE	7.62	127.80	111.80
1	E	292	ASP	CB-CG-OD2	-7.42	111.62	118.30
1	H	294	ARG	CG-CD-NE	7.38	127.31	111.80
1	B	89	ARG	CG-CD-NE	7.31	127.16	111.80
1	C	89	ARG	CG-CD-NE	7.23	126.97	111.80
1	H	89	ARG	CG-CD-NE	7.21	126.95	111.80
1	D	89	ARG	CG-CD-NE	7.12	126.75	111.80
1	A	89	ARG	CG-CD-NE	7.11	126.74	111.80
1	F	292	ASP	CB-CA-C	-6.95	96.51	110.40
1	H	292	ASP	CB-CA-C	-6.93	96.55	110.40
1	G	89	ARG	CG-CD-NE	6.90	126.29	111.80
1	F	89	ARG	CG-CD-NE	6.82	126.13	111.80
1	A	334	ASP	CB-CG-OD1	6.26	123.93	118.30
1	G	79	ARG	NE-CZ-NH2	-6.02	117.29	120.30
1	D	16	ARG	NE-CZ-NH2	-5.68	117.46	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	79	ARG	NE-CZ-NH2	-5.67	117.47	120.30
1	G	276	GLN	CA-CB-CG	5.58	125.67	113.40
1	C	292	ASP	CB-CG-OD1	-5.35	113.49	118.30
1	D	268	ARG	NE-CZ-NH1	5.28	122.94	120.30
1	D	4	LEU	CB-CG-CD1	-5.12	102.30	111.00
1	G	272	GLU	CB-CA-C	-5.07	100.26	110.40
1	G	13	MET	CG-SD-CE	5.05	108.27	100.20

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	89	ARG	Sidechain
1	B	136	TYR	Peptide
1	B	89	ARG	Sidechain
1	C	136	TYR	Peptide
1	C	89	ARG	Sidechain
1	D	89	ARG	Sidechain
1	E	89	ARG	Sidechain
1	F	89	ARG	Sidechain
1	G	89	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	2472	0	2446	24	0
1	B	2655	0	2617	28	0
1	C	2608	0	2577	21	0
1	D	2430	0	2396	7	0
1	E	2345	0	2305	19	0
1	F	2359	0	2332	17	0
1	G	2665	0	2634	31	0
1	H	2348	0	2307	7	0
2	A	48	0	25	6	0
2	B	48	0	25	1	0
2	C	48	0	25	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	48	0	25	2	0
2	F	48	0	25	2	0
2	G	48	0	25	1	0
2	H	48	0	25	1	0
3	A	13	0	5	0	0
3	B	13	0	5	0	0
3	C	13	0	5	0	0
3	D	13	0	5	0	0
3	E	13	0	5	1	0
3	F	13	0	5	3	0
3	G	13	0	5	1	0
3	H	13	0	5	1	0
4	A	181	0	0	3	0
4	B	155	0	0	2	0
4	C	142	0	0	3	0
4	D	161	0	0	2	0
4	E	121	0	0	0	0
4	F	156	0	0	1	0
4	G	132	0	0	1	0
4	H	113	0	0	0	0
All	All	21483	0	19829	147	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 4.

All (147) 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:B:180:GLU:HG2	4:B:2124:HOH:O	1.59	1.02
1:B:328:ILE:HG23	1:B:332:ILE:HD12	1.40	1.02
1:D:142:GLU:HB3	4:D:2095:HOH:O	1.63	0.97
1:A:79:ARG:NH1	1:G:79:ARG:CZ	2.27	0.97
1:F:292:ASP:HB3	1:F:294:ARG:HG2	1.50	0.93
1:F:332:ILE:HD12	1:F:332:ILE:O	1.67	0.93
1:F:270:LEU:HD11	1:F:332:ILE:HG21	1.51	0.90
1:B:328:ILE:HG23	1:B:332:ILE:CD1	2.04	0.86
1:G:76:ASN:OD1	1:G:79:ARG:NH1	2.08	0.85
1:G:271:ASN:OD1	1:G:281:MET:HE2	1.77	0.85
1:G:201:ARG:HH22	1:G:328:ILE:HG22	1.45	0.81
1:A:222:PHE:CE1	2:A:500:NAP:C4N	2.67	0.76
1:B:328:ILE:CG2	1:B:332:ILE:HD12	2.17	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:267:LEU:HD22	1:B:281:MET:CE	2.17	0.73
1:F:270:LEU:HD11	1:F:332:ILE:CG2	2.17	0.73
1:B:267:LEU:CD2	1:B:281:MET:HE1	2.20	0.72
1:H:231:THR:HG22	1:H:282:ALA:HB2	1.71	0.71
3:F:501:FLC:OA2	3:F:501:FLC:CBC	2.39	0.70
1:B:201:ARG:HH22	1:B:328:ILE:HG22	1.56	0.70
1:E:292:ASP:HB3	1:E:294:ARG:HG3	1.73	0.70
1:A:79:ARG:NH1	1:G:79:ARG:NH2	2.40	0.69
1:C:233:LYS:HE3	1:C:234:TYR:CZ	2.28	0.69
1:C:222:PHE:CE1	2:C:500:NAP:C4N	2.75	0.69
1:B:265:ASN:O	1:B:268:ARG:HG2	1.94	0.68
1:A:79:ARG:CZ	1:G:79:ARG:NE	2.57	0.67
1:B:270:LEU:HD21	1:B:332:ILE:HD11	1.77	0.67
1:F:292:ASP:CB	1:F:294:ARG:HG2	2.23	0.67
1:C:258:MET:O	1:C:263:ASN:ND2	2.28	0.66
1:C:244:MET:C	4:C:2117:HOH:O	2.34	0.66
1:B:267:LEU:CD2	1:B:281:MET:CE	2.75	0.64
1:B:222:PHE:CE1	2:B:500:NAP:C4N	2.80	0.64
1:A:79:ARG:CZ	1:G:79:ARG:CZ	2.79	0.61
1:E:146:MET:HE3	1:E:177:LYS:CG	2.31	0.61
1:B:267:LEU:HD23	1:B:281:MET:HE1	1.83	0.61
3:E:501:FLC:OA2	3:E:501:FLC:CBC	2.49	0.60
1:G:201:ARG:HH21	1:G:332:ILE:HD11	1.67	0.60
1:C:233:LYS:NZ	2:C:500:NAP:O1X	2.31	0.60
1:A:79:ARG:HH11	1:G:79:ARG:NH2	1.99	0.59
1:E:222:PHE:CE1	2:E:500:NAP:C4N	2.84	0.59
1:F:222:PHE:CE1	2:F:500:NAP:C4N	2.86	0.59
1:H:292:ASP:HB3	1:H:294:ARG:HG3	1.85	0.58
1:B:270:LEU:HD21	1:B:332:ILE:CD1	2.33	0.58
1:F:231:THR:HG22	1:F:282:ALA:HB2	1.85	0.57
1:F:332:ILE:O	1:F:332:ILE:CD1	2.48	0.57
1:B:292:ASP:OD1	1:B:294:ARG:HB2	2.05	0.57
1:D:142:GLU:HG2	4:D:2093:HOH:O	2.05	0.56
1:A:142:GLU:HG2	4:A:2117:HOH:O	2.06	0.56
1:G:330:GLN:HG3	1:G:331:HIS:CE1	2.41	0.56
1:G:222:PHE:CE1	2:G:500:NAP:C4N	2.89	0.56
1:C:79:ARG:CD	1:E:79:ARG:CZ	2.84	0.56
1:G:271:ASN:CG	1:G:281:MET:HE2	2.26	0.55
1:G:235:LEU:HD11	1:G:281:MET:HE1	1.89	0.54
1:B:267:LEU:HD22	1:B:281:MET:HE2	1.88	0.54
1:G:42:LEU:HD13	1:G:79:ARG:NH1	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:196:TYR:CG	1:D:203:VAL:HG21	2.44	0.53
1:G:292:ASP:OD1	1:G:294:ARG:HB2	2.07	0.53
1:F:294:ARG:HH11	1:F:294:ARG:HG3	1.72	0.53
1:A:222:PHE:CD1	2:A:500:NAP:C4N	2.91	0.53
1:D:286:LEU:HD22	1:D:298:VAL:HG21	1.90	0.52
1:G:201:ARG:HH22	1:G:328:ILE:CG2	2.17	0.52
1:E:61:ASP:HB3	1:E:299:LEU:HD11	1.90	0.52
1:B:267:LEU:HD22	1:B:281:MET:HE1	1.83	0.52
1:F:264:LEU:HA	1:F:267:LEU:HD12	1.91	0.52
1:D:42:LEU:HD22	1:F:79:ARG:HH22	1.74	0.52
1:A:259:LEU:N	4:A:2157:HOH:O	2.42	0.52
1:C:233:LYS:HE3	1:C:234:TYR:CE2	2.46	0.51
1:F:327:GLN:HG3	4:F:2154:HOH:O	2.11	0.51
1:G:330:GLN:HB3	1:G:331:HIS:CD2	2.45	0.51
1:B:79:ARG:CD	1:H:79:ARG:CZ	2.88	0.51
1:G:241:ASP:OD1	1:G:304[A]:ARG:NH2	2.45	0.50
1:B:201:ARG:NH2	1:B:328:ILE:HG22	2.24	0.50
1:B:81:LEU:HD12	1:B:85:PHE:HB2	1.94	0.50
1:A:268:ARG:HD3	4:A:2159:HOH:O	2.12	0.50
1:C:28:LEU:HD12	1:C:28:LEU:N	2.26	0.49
1:A:33:TRP:HB2	2:A:500:NAP:H2D	1.93	0.49
1:H:222:PHE:CE1	2:H:500:NAP:C4N	2.95	0.49
1:C:139:ARG:NE	4:C:2060:HOH:O	2.46	0.49
1:E:146:MET:CE	1:E:177:LYS:HG2	2.43	0.49
1:A:61:ASP:HB3	1:A:299:LEU:HD11	1.94	0.48
1:D:196:TYR:CD1	1:D:203:VAL:CG2	2.96	0.48
1:A:263:ASN:ND2	1:A:335:GLY:O	2.45	0.48
1:A:203:VAL:HG12	1:A:209:LEU:HG	1.96	0.48
1:G:76:ASN:OD1	1:G:79:ARG:CZ	2.61	0.48
1:C:209:LEU:HD13	1:C:294:ARG:HB3	1.96	0.48
1:A:299:LEU:O	2:A:500:NAP:H51N	2.13	0.48
1:C:60:PHE:HE2	1:C:92:LEU:HD22	1.79	0.48
1:B:258:MET:O	1:B:263:ASN:ND2	2.47	0.48
1:G:330:GLN:CG	1:G:331:HIS:NE2	2.77	0.47
1:G:330:GLN:HG2	1:G:331:HIS:NE2	2.29	0.47
1:E:203:VAL:HG12	1:E:209:LEU:HG	1.97	0.47
1:A:303:SER:OG	2:A:500:NAP:O3X	2.25	0.47
1:E:32:LEU:N	1:E:32:LEU:HD12	2.30	0.47
1:G:271:ASN:CB	1:G:281:MET:HE2	2.45	0.47
1:B:273:MET:CE	1:B:327:GLN:HB3	2.45	0.46
1:A:204:ASP:OD1	1:A:209:LEU:HD11	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:270:LEU:HB3	1:E:281:MET:HE1	1.97	0.45
1:E:146:MET:CE	1:E:177:LYS:CG	2.94	0.45
1:E:146:MET:HE3	1:E:177:LYS:HG2	1.98	0.45
1:G:328:ILE:HG23	1:G:332:ILE:HG12	1.97	0.45
1:G:198:LEU:HD22	1:G:332:ILE:CD1	2.48	0.44
3:G:501:FLC:CBC	3:G:501:FLC:OA2	2.65	0.44
2:F:500:NAP:C4N	3:F:501:FLC:CGC	2.95	0.44
1:C:79:ARG:HD2	1:E:79:ARG:CZ	2.47	0.44
1:E:31:GLY:C	1:E:32:LEU:HD12	2.38	0.44
1:C:188:PRO:HG2	4:C:2105:HOH:O	2.17	0.44
1:E:146:MET:HE3	1:E:177:LYS:HG3	2.00	0.43
1:C:286:LEU:HD22	1:C:298:VAL:HG21	2.00	0.43
1:G:270:LEU:HD21	1:G:332:ILE:HG21	1.99	0.43
1:H:291:LYS:HG2	1:H:292:ASP:OD1	2.18	0.43
1:A:298:VAL:O	1:A:298:VAL:HG23	2.18	0.43
1:C:61:ASP:HB3	1:C:299:LEU:HD11	2.00	0.43
1:B:226:ALA:HB1	1:B:337:LEU:HD22	2.01	0.43
1:C:240:GLN:O	1:C:241:ASP:HB2	2.19	0.43
1:G:201:ARG:NH2	1:G:328:ILE:CG2	2.81	0.43
1:G:323:LYS:HG3	4:G:2128:HOH:O	2.17	0.43
1:E:117:LEU:HA	1:E:117:LEU:HD23	1.86	0.43
1:C:79:ARG:HD2	1:E:79:ARG:NE	2.34	0.43
1:A:167:ILE:HD12	1:A:170:TYR:CG	2.54	0.42
1:G:76:ASN:OD1	1:G:79:ARG:NH2	2.52	0.42
1:E:193:GLN:OE1	2:E:500:NAP:H2N	2.19	0.42
1:B:231:THR:OG1	1:B:233:LYS:HB2	2.19	0.42
1:A:61:ASP:CB	1:A:299:LEU:HD11	2.49	0.42
1:C:203:VAL:CG1	1:C:209:LEU:HG	2.49	0.42
1:H:203:VAL:CG1	1:H:209:LEU:HG	2.50	0.42
1:G:286:LEU:HD22	1:G:298:VAL:HG21	2.01	0.42
1:A:222:PHE:CD1	2:A:500:NAP:C5N	3.02	0.41
1:B:79:ARG:NH2	1:H:75:GLU:OE1	2.52	0.41
1:B:12:GLN:HG2	4:B:2004:HOH:O	2.20	0.41
1:F:291:LYS:HG2	1:F:292:ASP:OD1	2.19	0.41
1:B:273:MET:HE2	1:B:327:GLN:HB3	2.01	0.41
1:B:61:ASP:HB3	1:B:299:LEU:HD11	2.02	0.41
1:F:77:PHE:CE2	1:F:81:LEU:HD12	2.56	0.41
1:G:330:GLN:CG	1:G:331:HIS:CD2	3.04	0.41
1:B:260:THR:H	1:B:263:ASN:HD22	1.68	0.41
1:C:222:PHE:CD1	2:C:500:NAP:C5N	3.04	0.41
3:H:501:FLC:CBC	3:H:501:FLC:OA2	2.69	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:100:TYR:CG	3:F:501:FLC:HA1	2.56	0.40
1:C:33:TRP:HB2	2:C:500:NAP:H2D	2.02	0.40
1:G:281:MET:HB2	1:G:281:MET:HE2	1.87	0.40
1:C:223:THR:HG22	2:C:500:NAP:O2N	2.20	0.40
1:D:192:HIS:CE1	1:D:208:LEU:HD21	2.57	0.40
1:F:60:PHE:HE2	1:F:92:LEU:HD22	1.86	0.40
1:F:209:LEU:HD13	1:F:294:ARG:HB3	2.04	0.40
1:A:61:ASP:OD2	1:A:97:LYS:NZ	2.46	0.40
1:A:228:GLY:HA2	1:A:231:THR:HG23	2.03	0.40
1:A:4:LEU:HD12	1:E:4:LEU:HD12	2.03	0.40
1:E:167:ILE:HD13	1:E:167:ILE:HG21	1.88	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	310/366 (85%)	300 (97%)	10 (3%)	0	100	100
1	B	332/366 (91%)	322 (97%)	10 (3%)	0	100	100
1	C	325/366 (89%)	314 (97%)	11 (3%)	0	100	100
1	D	304/366 (83%)	297 (98%)	7 (2%)	0	100	100
1	E	291/366 (80%)	282 (97%)	9 (3%)	0	100	100
1	F	294/366 (80%)	287 (98%)	7 (2%)	0	100	100
1	G	332/366 (91%)	322 (97%)	10 (3%)	0	100	100
1	H	291/366 (80%)	283 (97%)	7 (2%)	1 (0%)	46	36
All	All	2479/2928 (85%)	2407 (97%)	71 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	194	PRO

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	262/307 (85%)	260 (99%)	2 (1%)	86	86
1	B	282/307 (92%)	280 (99%)	2 (1%)	88	88
1	C	278/307 (91%)	278 (100%)	0	100	100
1	D	257/307 (84%)	256 (100%)	1 (0%)	93	94
1	E	247/307 (80%)	246 (100%)	1 (0%)	93	94
1	F	250/307 (81%)	247 (99%)	3 (1%)	78	76
1	G	284/307 (92%)	281 (99%)	3 (1%)	80	79
1	H	248/307 (81%)	247 (100%)	1 (0%)	93	94
All	All	2108/2456 (86%)	2095 (99%)	13 (1%)	90	90

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	89	ARG
1	A	338	ASN
1	B	89	ARG
1	B	240	GLN
1	D	89	ARG
1	E	281	MET
1	F	81	LEU
1	F	265	ASN
1	F	332	ILE
1	G	89	ARG
1	G	268	ARG
1	G	281	MET
1	H	294	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such

sidechains are listed below:

Mol	Chain	Res	Type
1	B	240	GLN
1	B	245	HIS
1	B	263	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

15 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	NAP	A	500	-	42,52,52	1.23	7 (16%)	54,80,80	2.07	15 (27%)
3	FLC	A	501	-	3,12,12	1.48	0	3,17,17	2.28	2 (66%)
2	NAP	B	500	-	42,52,52	1.32	6 (14%)	54,80,80	2.16	10 (18%)
3	FLC	B	501	-	3,12,12	1.10	0	3,17,17	1.80	1 (33%)
2	NAP	C	500	-	42,52,52	1.20	4 (9%)	54,80,80	2.01	13 (24%)
3	FLC	C	501	-	3,12,12	1.19	0	3,17,17	1.18	0
3	FLC	D	501	-	3,12,12	1.29	1 (33%)	3,17,17	0.81	0
2	NAP	E	500	-	42,52,52	0.95	3 (7%)	54,80,80	2.10	10 (18%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	FLC	E	501	-	3,12,12	1.58	1 (33%)	3,17,17	2.58	1 (33%)
2	NAP	F	500	-	42,52,52	1.18	5 (11%)	54,80,80	1.76	7 (12%)
3	FLC	F	501	-	3,12,12	1.33	0	3,17,17	3.90	2 (66%)
2	NAP	G	500	-	42,52,52	1.18	4 (9%)	54,80,80	1.84	7 (12%)
3	FLC	G	501	-	3,12,12	0.95	0	3,17,17	1.14	0
2	NAP	H	500	-	42,52,52	1.08	4 (9%)	54,80,80	2.08	9 (16%)
3	FLC	H	501	-	3,12,12	0.43	0	3,17,17	1.15	0

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	NAP	A	500	-	-	0/27/67/67	0/5/5/5
3	FLC	A	501	-	-	0/6/16/16	0/0/0/0
2	NAP	B	500	-	-	0/27/67/67	0/5/5/5
3	FLC	B	501	-	-	0/6/16/16	0/0/0/0
2	NAP	C	500	-	-	0/27/67/67	0/5/5/5
3	FLC	C	501	-	-	0/6/16/16	0/0/0/0
3	FLC	D	501	-	-	0/6/16/16	0/0/0/0
2	NAP	E	500	-	-	0/27/67/67	0/5/5/5
3	FLC	E	501	-	-	0/6/16/16	0/0/0/0
2	NAP	F	500	-	-	0/27/67/67	0/5/5/5
3	FLC	F	501	-	-	0/6/16/16	0/0/0/0
2	NAP	G	500	-	-	0/27/67/67	0/5/5/5
3	FLC	G	501	-	-	0/6/16/16	0/0/0/0
2	NAP	H	500	-	-	0/27/67/67	0/5/5/5
3	FLC	H	501	-	-	0/6/16/16	0/0/0/0

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	501	FLC	CG-CB	-2.00	1.51	1.54
2	B	500	NAP	C2A-N3A	2.01	1.35	1.32
2	A	500	NAP	C2A-N3A	2.02	1.35	1.32
2	C	500	NAP	C2A-N3A	2.06	1.35	1.32
3	D	501	FLC	OHB-CB	2.07	1.46	1.43
2	F	500	NAP	C4A-N3A	2.09	1.38	1.35
2	F	500	NAP	C2A-N3A	2.12	1.35	1.32
2	C	500	NAP	O4B-C1B	2.14	1.43	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	NAP	C4N-C3N	2.15	1.43	1.39
2	F	500	NAP	O4B-C1B	2.21	1.44	1.41
2	E	500	NAP	C4A-N3A	2.23	1.38	1.35
2	E	500	NAP	O4B-C1B	2.28	1.44	1.41
2	A	500	NAP	O2D-C2D	2.33	1.48	1.43
2	A	500	NAP	O4D-C1D	2.37	1.44	1.41
2	B	500	NAP	C4N-C3N	2.37	1.43	1.39
2	A	500	NAP	O4B-C1B	2.40	1.44	1.41
2	G	500	NAP	C5A-C4A	2.46	1.46	1.40
2	H	500	NAP	O4D-C1D	2.50	1.44	1.41
2	H	500	NAP	C2A-N3A	2.52	1.36	1.32
2	G	500	NAP	C2A-N3A	2.61	1.36	1.32
2	F	500	NAP	O4D-C1D	2.67	1.44	1.41
2	A	500	NAP	PA-O1A	2.72	1.61	1.51
2	B	500	NAP	C3N-C7N	2.72	1.54	1.50
2	G	500	NAP	O4B-C1B	2.75	1.44	1.41
2	B	500	NAP	O4D-C1D	2.83	1.44	1.41
2	H	500	NAP	C5A-C4A	3.01	1.47	1.40
2	H	500	NAP	O4B-C1B	3.04	1.45	1.41
2	A	500	NAP	C5A-C4A	3.11	1.47	1.40
2	B	500	NAP	O4B-C1B	3.23	1.45	1.41
2	E	500	NAP	C5A-C4A	3.27	1.47	1.40
2	F	500	NAP	C5A-C4A	3.45	1.48	1.40
2	C	500	NAP	C5A-C4A	3.50	1.48	1.40
2	G	500	NAP	O4D-C1D	3.79	1.46	1.41
2	B	500	NAP	C5A-C4A	3.89	1.49	1.40
2	C	500	NAP	O4D-C1D	4.04	1.46	1.41

All (77) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	500	NAP	N3A-C2A-N1A	-9.63	121.52	128.89
2	E	500	NAP	N3A-C2A-N1A	-9.22	121.83	128.89
2	F	500	NAP	N3A-C2A-N1A	-8.70	122.24	128.89
2	C	500	NAP	N3A-C2A-N1A	-8.64	122.28	128.89
2	B	500	NAP	N3A-C2A-N1A	-8.61	122.30	128.89
2	A	500	NAP	N3A-C2A-N1A	-8.39	122.47	128.89
2	G	500	NAP	N3A-C2A-N1A	-8.08	122.71	128.89
3	F	501	FLC	CB-CA-CAC	-5.79	105.70	114.96
2	A	500	NAP	O3-PN-O5D	-4.12	92.00	102.94
3	E	501	FLC	CB-CA-CAC	-4.09	108.41	114.96
2	C	500	NAP	O3-PN-O5D	-4.05	92.19	102.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	NAP	C4A-C5A-N7A	-3.56	106.20	109.48
2	G	500	NAP	C4A-C5A-N7A	-3.55	106.21	109.48
3	F	501	FLC	CB-CG-CGC	-3.48	109.39	114.96
2	F	500	NAP	O3-PN-O5D	-3.40	93.92	102.94
2	E	500	NAP	O2B-P2B-O1X	-3.37	98.70	107.11
2	H	500	NAP	O7N-C7N-C3N	-3.35	115.93	119.59
2	E	500	NAP	O3-PN-O5D	-3.32	94.12	102.94
2	B	500	NAP	O3-PN-O5D	-3.30	94.18	102.94
2	G	500	NAP	O7N-C7N-N7N	-3.13	118.19	122.59
3	A	501	FLC	CB-CG-CGC	-3.11	109.99	114.96
2	F	500	NAP	C1B-N9A-C4A	-2.85	122.64	126.94
2	H	500	NAP	O3-PN-O5D	-2.67	95.85	102.94
2	A	500	NAP	O2B-P2B-O1X	-2.65	100.50	107.11
2	H	500	NAP	C1B-N9A-C4A	-2.57	123.06	126.94
2	B	500	NAP	C4A-C5A-N7A	-2.54	107.15	109.48
2	C	500	NAP	O2B-P2B-O1X	-2.50	100.87	107.11
2	B	500	NAP	O7N-C7N-N7N	-2.48	119.11	122.59
2	C	500	NAP	O7N-C7N-N7N	-2.39	119.23	122.59
2	E	500	NAP	C4A-C5A-N7A	-2.34	107.32	109.48
2	A	500	NAP	O3-PA-O5B	-2.34	96.73	102.94
3	B	501	FLC	CB-CG-CGC	-2.32	111.25	114.96
2	B	500	NAP	C1B-N9A-C4A	-2.24	123.56	126.94
2	C	500	NAP	C2N-C3N-C4N	-2.21	115.82	118.29
3	A	501	FLC	CB-CA-CAC	-2.20	111.43	114.96
2	A	500	NAP	O4D-C4D-C5D	-2.20	101.44	109.32
2	A	500	NAP	C1B-N9A-C4A	-2.17	123.67	126.94
2	B	500	NAP	C2N-C3N-C4N	-2.16	115.89	118.29
2	G	500	NAP	O3-PN-O5D	-2.14	97.26	102.94
2	C	500	NAP	O5B-PA-O1A	-2.05	101.67	109.62
2	F	500	NAP	C4A-C5A-N7A	-2.02	107.62	109.48
2	A	500	NAP	C3N-C2N-N1N	2.05	122.72	120.36
2	H	500	NAP	O2N-PN-O3	2.11	114.64	105.09
2	G	500	NAP	O3X-P2B-O2X	2.12	115.46	107.38
2	E	500	NAP	O2A-PA-O1A	2.19	124.38	112.53
2	H	500	NAP	O2X-P2B-O1X	2.23	117.77	110.58
2	C	500	NAP	O2X-P2B-O1X	2.29	117.96	110.58
2	E	500	NAP	C2A-N1A-C6A	2.29	122.87	118.77
2	A	500	NAP	C3N-C7N-N7N	2.32	120.35	117.82
2	A	500	NAP	O2N-PN-O1N	2.33	125.17	112.53
2	G	500	NAP	C3N-C2N-N1N	2.33	123.05	120.36
2	E	500	NAP	C3N-C2N-N1N	2.40	123.13	120.36
2	F	500	NAP	C2A-N1A-C6A	2.45	123.15	118.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	500	NAP	O2X-P2B-O1X	2.48	118.55	110.58
2	A	500	NAP	O2D-C2D-C3D	2.51	119.98	111.83
2	C	500	NAP	P2B-O2B-C2B	2.63	127.87	121.56
2	A	500	NAP	C4D-O4D-C1D	2.63	112.61	109.72
2	B	500	NAP	C2A-N1A-C6A	2.63	123.47	118.77
2	H	500	NAP	O2A-PA-O1A	2.65	126.88	112.53
2	C	500	NAP	C2A-N1A-C6A	2.65	123.50	118.77
2	F	500	NAP	O2A-PA-O1A	2.66	126.97	112.53
2	E	500	NAP	P2B-O2B-C2B	2.68	127.99	121.56
2	C	500	NAP	O2A-PA-O1A	2.71	127.23	112.53
2	E	500	NAP	O2X-P2B-O1X	2.81	119.62	110.58
2	B	500	NAP	O2A-PA-O1A	3.20	129.89	112.53
2	A	500	NAP	P2B-O2B-C2B	3.25	129.35	121.56
2	H	500	NAP	C3N-C7N-N7N	3.48	121.63	117.82
2	C	500	NAP	C4D-O4D-C1D	3.62	113.69	109.72
2	A	500	NAP	O2A-PA-O1A	3.67	132.40	112.53
2	C	500	NAP	O7N-C7N-C3N	3.67	123.60	119.59
2	C	500	NAP	O4D-C1D-N1N	4.19	112.73	108.13
2	B	500	NAP	C3N-C7N-N7N	4.64	122.90	117.82
2	A	500	NAP	O4D-C1D-N1N	4.66	113.25	108.13
2	G	500	NAP	C3N-C7N-N7N	5.11	123.41	117.82
2	H	500	NAP	O4D-C1D-N1N	6.94	115.76	108.13
2	E	500	NAP	O4D-C1D-N1N	7.22	116.06	108.13
2	B	500	NAP	O4D-C1D-N1N	7.84	116.75	108.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

11 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	NAP	6	0
2	B	500	NAP	1	0
2	C	500	NAP	5	0
2	E	500	NAP	2	0
3	E	501	FLC	1	0
2	F	500	NAP	2	0
3	F	501	FLC	3	0
2	G	500	NAP	1	0
3	G	501	FLC	1	0
2	H	500	NAP	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	501	FLC	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 ⓘ

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, 95th 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 ²)	Q<0.9
1	A	312/366 (85%)	0.14	10 (3%) 51 58	12, 21, 47, 77	1 (0%)
1	B	335/366 (91%)	0.39	34 (10%) 9 10	14, 26, 73, 97	1 (0%)
1	C	330/366 (90%)	0.36	27 (8%) 14 16	13, 24, 67, 98	1 (0%)
1	D	308/366 (84%)	-0.09	1 (0%) 94 95	12, 20, 41, 55	1 (0%)
1	E	297/366 (81%)	0.49	29 (9%) 10 10	16, 28, 65, 81	0
1	F	298/366 (81%)	0.00	5 (1%) 73 78	13, 22, 48, 78	1 (0%)
1	G	335/366 (91%)	0.24	18 (5%) 29 34	14, 24, 59, 84	0
1	H	297/366 (81%)	0.36	22 (7%) 17 20	15, 26, 59, 79	2 (0%)
All	All	2512/2928 (85%)	0.24	146 (5%) 26 30	12, 24, 59, 98	7 (0%)

All (146) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	228	GLY	7.8
1	C	237	GLY	7.7
1	C	262	ALA	7.3
1	B	263	ASN	6.8
1	C	260	THR	6.8
1	F	264	LEU	6.6
1	C	263	ASN	5.7
1	E	339	LEU	5.6
1	C	265	ASN	5.5
1	C	241	ASP	5.5
1	B	246	ARG	5.4
1	B	260	THR	5.3
1	C	239	PRO	5.3
1	B	262	ALA	5.3
1	B	264	LEU	5.2
1	A	228	GLY	5.2

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Mol	Chain	Res	Type	RSRZ
1	B	235	LEU	5.0
1	B	337	LEU	4.8
1	H	345	ASP	4.7
1	B	258	MET	4.7
1	A	259	LEU	4.7
1	B	266	SER	4.6
1	E	271	ASN	4.6
1	B	236	ASN	4.6
1	H	337	LEU	4.4
1	B	333	ALA	4.4
1	E	275	GLN	4.4
1	H	341	GLN	4.3
1	B	241	ASP	4.3
1	B	228	GLY	4.2
1	E	340	TRP	4.2
1	E	345	ASP	4.1
1	G	330	GLN	4.1
1	E	276	GLN	4.0
1	H	340	TRP	4.0
1	G	259	LEU	3.9
1	E	272	GLU	3.9
1	H	339	LEU	3.9
1	B	265	ASN	3.9
1	E	320	PHE	3.9
1	B	340	TRP	3.9
1	G	333	ALA	3.9
1	C	339	LEU	3.8
1	C	240	GLN	3.8
1	G	240	GLN	3.7
1	C	342	ALA	3.7
1	E	233	LYS	3.6
1	C	345	ASP	3.6
1	A	260	THR	3.5
1	B	339	LEU	3.5
1	E	269	LEU	3.5
1	G	241	ASP	3.5
1	H	342	ALA	3.5
1	G	256	PRO	3.5
1	B	244	MET	3.5
1	D	202	TRP	3.4
1	E	270	LEU	3.4
1	B	240	GLN	3.4

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Mol	Chain	Res	Type	RSRZ
1	E	231	THR	3.4
1	E	334	ASP	3.3
1	E	232	GLY	3.3
1	C	343	SER	3.3
1	C	259	LEU	3.3
1	C	238	ILE	3.3
1	H	344	SER	3.3
1	E	337	LEU	3.2
1	G	245	HIS	3.2
1	A	264	LEU	3.1
1	E	336	GLU	3.1
1	G	332	ILE	3.1
1	E	273	MET	3.1
1	H	319	THR	3.1
1	C	336	GLU	3.1
1	C	266	SER	3.1
1	B	256	PRO	3.1
1	H	343	SER	3.0
1	B	245	HIS	3.0
1	A	232	GLY	3.0
1	C	335	GLY	3.0
1	C	337	LEU	3.0
1	E	278	GLY	3.0
1	C	261	GLU	2.9
1	F	332	ILE	2.9
1	B	259	LEU	2.9
1	E	344	SER	2.9
1	H	276	GLN	2.9
1	E	281	MET	2.9
1	B	341	GLN	2.8
1	E	335	GLY	2.8
1	C	234	TYR	2.8
1	G	238	ILE	2.8
1	A	2	VAL	2.8
1	B	239	PRO	2.8
1	B	269	LEU	2.8
1	H	228	GLY	2.8
1	B	234	TYR	2.7
1	H	335	GLY	2.7
1	G	339	LEU	2.7
1	G	319	THR	2.7
1	C	341	GLN	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	233	LYS	2.7
1	B	238	ILE	2.7
1	H	229	LEU	2.7
1	B	332	ILE	2.7
1	B	242	SER	2.7
1	F	228	GLY	2.7
1	G	269	LEU	2.6
1	C	236	ASN	2.6
1	A	231	THR	2.6
1	G	242	SER	2.5
1	G	335	GLY	2.5
1	B	345	ASP	2.5
1	H	273	MET	2.5
1	G	239	PRO	2.5
1	H	292	ASP	2.5
1	C	257	LYS	2.5
1	H	317	ASN	2.5
1	E	229	LEU	2.5
1	H	278	GLY	2.5
1	B	243	ARG	2.4
1	F	4	LEU	2.4
1	H	336	GLU	2.4
1	C	204	ASP	2.4
1	A	338	ASN	2.4
1	B	257	LYS	2.4
1	E	197	ASN	2.4
1	G	237	GLY	2.4
1	G	258	MET	2.4
1	C	258	MET	2.3
1	B	330	GLN	2.3
1	H	271	ASN	2.3
1	H	270	LEU	2.3
1	E	343	SER	2.2
1	B	237	GLY	2.2
1	E	282	ALA	2.2
1	E	280	SER	2.2
1	F	267	LEU	2.2
1	C	243	ARG	2.2
1	E	341	GLN	2.1
1	E	274	ALA	2.1
1	H	275	GLN	2.1
1	C	329	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	336	GLU	2.0
1	B	336	GLU	2.0
1	G	327	GLN	2.0
1	H	338	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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, 95th 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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	FLC	A	501	13/13	0.82	0.37	21.01	46,57,63,65	0
3	FLC	F	501	13/13	0.86	0.19	8.60	32,43,49,50	0
3	FLC	D	501	13/13	0.79	0.21	5.52	42,51,56,59	0
2	NAP	A	500	48/48	0.73	0.38	3.14	28,55,79,82	0
2	NAP	F	500	48/48	0.86	0.22	2.28	22,43,56,60	0
3	FLC	C	501	13/13	0.89	0.19	2.27	33,42,48,58	0
3	FLC	G	501	13/13	0.91	0.18	1.86	34,39,46,51	0
2	NAP	H	500	48/48	0.90	0.20	1.41	33,44,50,52	0
3	FLC	E	501	13/13	0.91	0.22	1.36	38,42,48,50	0
3	FLC	B	501	13/13	0.89	0.19	1.16	34,37,45,51	0
3	FLC	H	501	13/13	0.89	0.19	1.09	38,46,49,49	0
2	NAP	B	500	48/48	0.92	0.21	0.49	27,43,58,61	0
2	NAP	C	500	48/48	0.91	0.17	0.44	19,40,51,53	0
2	NAP	E	500	48/48	0.92	0.21	0.39	28,43,50,57	0
2	NAP	G	500	48/48	0.96	0.14	0.07	21,30,37,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.