



wwPDB X-ray Structure Validation Summary Report

Feb 26, 2014 – 06:00 PM GMT

PDB ID : 3BMO
Title : Structure of Pteridine Reductase 1 (PTR1) from Trypanosoma brucei in ternary complex with cofactor (NADP+) and inhibitor (Compound AX4)
Authors : Martini, V.P.; Iulek, J.; Hunter, W.N.; Tulloch, L.B.
Deposited on : 2007-12-13
Resolution : 1.60 Å(reported)

This is a wwPDB validation summary 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/ValidationPDFNotes.html>

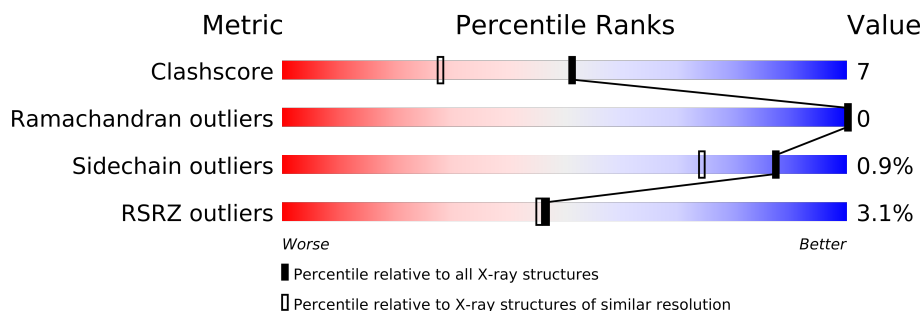
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	79885	2199 (1.60-1.60)
Ramachandran outliers	78287	2126 (1.60-1.60)
Sidechain outliers	78261	2125 (1.60-1.60)
RSRZ outliers	66119	1872 (1.60-1.60)

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. 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	288	
1	C	288	
1	D	288	
2	B	288	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	ACT	A	270	-	X
3	ACT	A	271	-	X
3	ACT	B	269	-	X
3	ACT	C	269	-	X
3	ACT	D	269	-	X
6	AX4	A	274[A]	-	X
6	AX4	D	271[A]	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
7	DTT	A	275	-	X
7	DTT	B	272	-	X
8	D1D	A	276[B]	-	X
8	D1D	B	273[B]	-	X
8	D1D	C	274[B]	-	X
8	D1D	D	273[B]	-	X
9	GOL	B	274[C]	-	X
9	GOL	B	275	-	X

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 9481 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 Pteridine reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	255	Total	C	N	O	S	0	23	0
			2013	1276	346	378	13			
1	C	252	Total	C	N	O	S	0	16	0
			1962	1241	344	365	12			
1	D	251	Total	C	N	O	S	0	21	0
			1981	1266	341	362	12			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP O76290
A	-18	GLY	-	EXPRESSION TAG	UNP O76290
A	-17	SER	-	EXPRESSION TAG	UNP O76290
A	-16	SER	-	EXPRESSION TAG	UNP O76290
A	-15	HIS	-	EXPRESSION TAG	UNP O76290
A	-14	HIS	-	EXPRESSION TAG	UNP O76290
A	-13	HIS	-	EXPRESSION TAG	UNP O76290
A	-12	HIS	-	EXPRESSION TAG	UNP O76290
A	-11	HIS	-	EXPRESSION TAG	UNP O76290
A	-10	HIS	-	EXPRESSION TAG	UNP O76290
A	-9	SER	-	EXPRESSION TAG	UNP O76290
A	-8	SER	-	EXPRESSION TAG	UNP O76290
A	-7	GLY	-	EXPRESSION TAG	UNP O76290
A	-6	LEU	-	EXPRESSION TAG	UNP O76290
A	-5	VAL	-	EXPRESSION TAG	UNP O76290
A	-4	PRO	-	EXPRESSION TAG	UNP O76290
A	-3	ARG	-	EXPRESSION TAG	UNP O76290
A	-2	GLY	-	EXPRESSION TAG	UNP O76290
A	-1	SER	-	EXPRESSION TAG	UNP O76290
A	0	HIS	-	EXPRESSION TAG	UNP O76290
C	-19	MET	-	EXPRESSION TAG	UNP O76290
C	-18	GLY	-	EXPRESSION TAG	UNP O76290
C	-17	SER	-	EXPRESSION TAG	UNP O76290

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-16	SER	-	EXPRESSION TAG	UNP O76290
C	-15	HIS	-	EXPRESSION TAG	UNP O76290
C	-14	HIS	-	EXPRESSION TAG	UNP O76290
C	-13	HIS	-	EXPRESSION TAG	UNP O76290
C	-12	HIS	-	EXPRESSION TAG	UNP O76290
C	-11	HIS	-	EXPRESSION TAG	UNP O76290
C	-10	HIS	-	EXPRESSION TAG	UNP O76290
C	-9	SER	-	EXPRESSION TAG	UNP O76290
C	-8	SER	-	EXPRESSION TAG	UNP O76290
C	-7	GLY	-	EXPRESSION TAG	UNP O76290
C	-6	LEU	-	EXPRESSION TAG	UNP O76290
C	-5	VAL	-	EXPRESSION TAG	UNP O76290
C	-4	PRO	-	EXPRESSION TAG	UNP O76290
C	-3	ARG	-	EXPRESSION TAG	UNP O76290
C	-2	GLY	-	EXPRESSION TAG	UNP O76290
C	-1	SER	-	EXPRESSION TAG	UNP O76290
C	0	HIS	-	EXPRESSION TAG	UNP O76290
D	-19	MET	-	EXPRESSION TAG	UNP O76290
D	-18	GLY	-	EXPRESSION TAG	UNP O76290
D	-17	SER	-	EXPRESSION TAG	UNP O76290
D	-16	SER	-	EXPRESSION TAG	UNP O76290
D	-15	HIS	-	EXPRESSION TAG	UNP O76290
D	-14	HIS	-	EXPRESSION TAG	UNP O76290
D	-13	HIS	-	EXPRESSION TAG	UNP O76290
D	-12	HIS	-	EXPRESSION TAG	UNP O76290
D	-11	HIS	-	EXPRESSION TAG	UNP O76290
D	-10	HIS	-	EXPRESSION TAG	UNP O76290
D	-9	SER	-	EXPRESSION TAG	UNP O76290
D	-8	SER	-	EXPRESSION TAG	UNP O76290
D	-7	GLY	-	EXPRESSION TAG	UNP O76290
D	-6	LEU	-	EXPRESSION TAG	UNP O76290
D	-5	VAL	-	EXPRESSION TAG	UNP O76290
D	-4	PRO	-	EXPRESSION TAG	UNP O76290
D	-3	ARG	-	EXPRESSION TAG	UNP O76290
D	-2	GLY	-	EXPRESSION TAG	UNP O76290
D	-1	SER	-	EXPRESSION TAG	UNP O76290
D	0	HIS	-	EXPRESSION TAG	UNP O76290

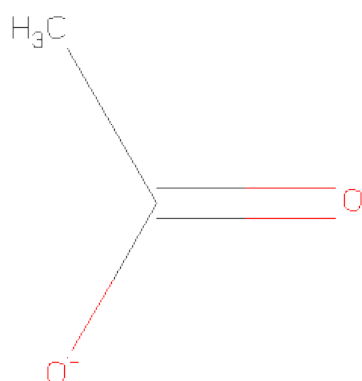
- Molecule 2 is a protein called Pteridine reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	249	Total	C	N	O	S	0	17	0
			1947	1237	336	361	13			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	EXPRESSION TAG	UNP O76290
B	-18	GLY	-	EXPRESSION TAG	UNP O76290
B	-17	SER	-	EXPRESSION TAG	UNP O76290
B	-16	SER	-	EXPRESSION TAG	UNP O76290
B	-15	HIS	-	EXPRESSION TAG	UNP O76290
B	-14	HIS	-	EXPRESSION TAG	UNP O76290
B	-13	HIS	-	EXPRESSION TAG	UNP O76290
B	-12	HIS	-	EXPRESSION TAG	UNP O76290
B	-11	HIS	-	EXPRESSION TAG	UNP O76290
B	-10	HIS	-	EXPRESSION TAG	UNP O76290
B	-9	SER	-	EXPRESSION TAG	UNP O76290
B	-8	SER	-	EXPRESSION TAG	UNP O76290
B	-7	GLY	-	EXPRESSION TAG	UNP O76290
B	-6	LEU	-	EXPRESSION TAG	UNP O76290
B	-5	VAL	-	EXPRESSION TAG	UNP O76290
B	-4	PRO	-	EXPRESSION TAG	UNP O76290
B	-3	ARG	-	EXPRESSION TAG	UNP O76290
B	-2	GLY	-	EXPRESSION TAG	UNP O76290
B	-1	SER	-	EXPRESSION TAG	UNP O76290
B	0	HIS	-	EXPRESSION TAG	UNP O76290

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		

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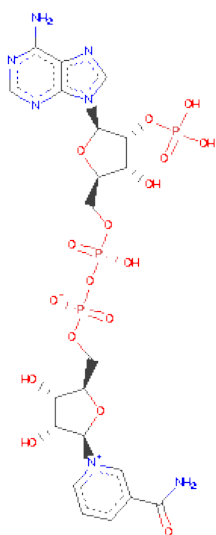
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Na	0	0
			1	1		

- Molecule 5 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDEPHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



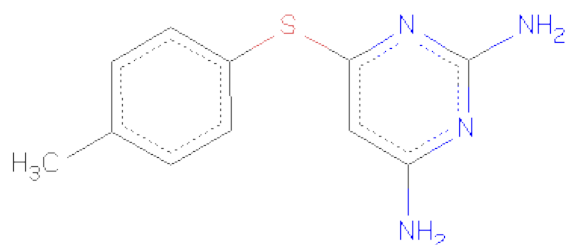
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total 48	C 21	N 7	O 17	P 3	0	0
5	B	1	Total 48	C 21	N 7	O 17	P 3	0	0

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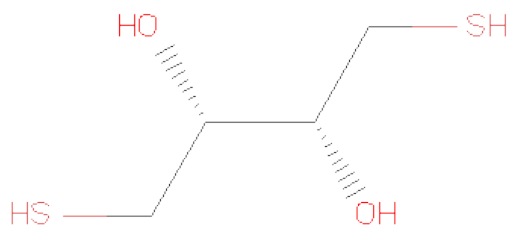
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
5	D	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 6 is 6-[(4-METHYLPHENYL)SULFANYL]PYRIMIDINE-2,4-DIAMINE (three-letter code: AX4) (formula: C₁₁H₁₂N₄S).



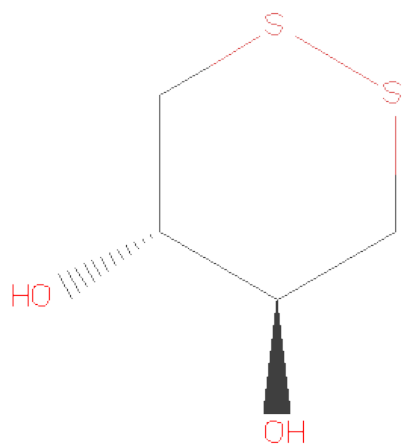
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	S	0	1
			16	11	4	1		
6	B	1	Total	C	N	S	0	1
			16	11	4	1		
6	C	1	Total	C	N	S	0	1
			16	11	4	1		
6	D	1	Total	C	N	S	0	1
			16	11	4	1		

- Molecule 7 is 2,3-DIHYDROXY-1,4-DITHIOBUTANE (three-letter code: DTT) (formula: C₄H₁₀O₂S₂).



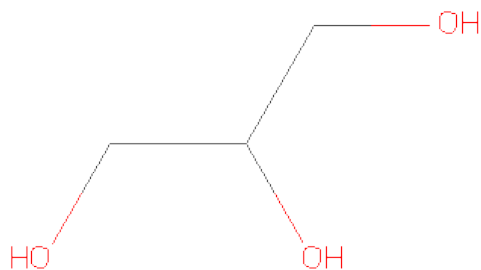
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	O	S	0	0
			8	4	2	2		
7	B	1	Total	C	O	S	0	0
			8	4	2	2		
7	C	1	Total	C	O	S	0	0
			8	4	2	2		
7	D	1	Total	C	O	S	0	0
			8	4	2	2		

- Molecule 8 is (4S,5S)-1,2-DITHIANE-4,5-DIOL (three-letter code: D1D) (formula: C₄H₈O₂S₂).



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

- Molecule 9 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	1	Total C O 6 3 3	0	0
9	B	1	Total C O 6 3 3	0	1
9	B	1	Total C O 6 3 3	0	0

- Molecule 10 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	A	309	Total O 309 309	0	0
10	B	304	Total O 304 304	0	0
10	C	298	Total O 298 298	0	0

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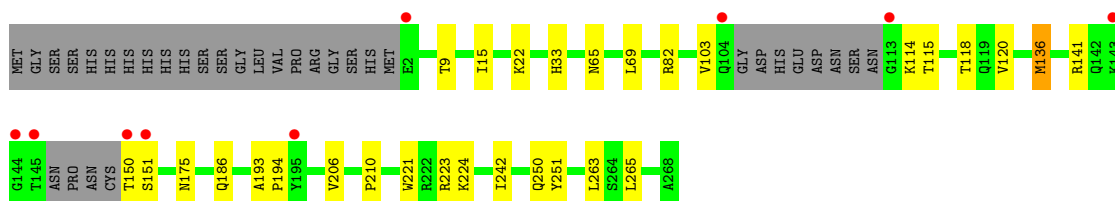
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	D	300	Total 300	O 300	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

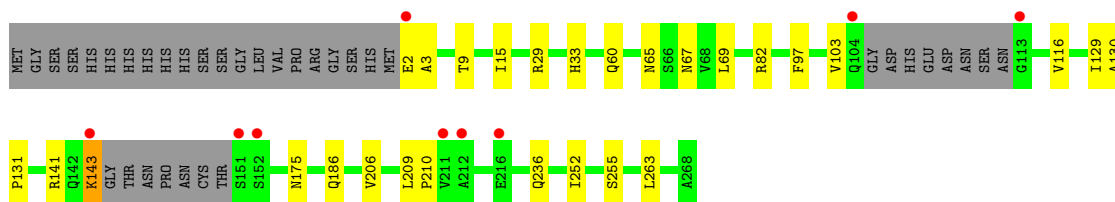
• Molecule 1: Pteridine reductase

Chain A:



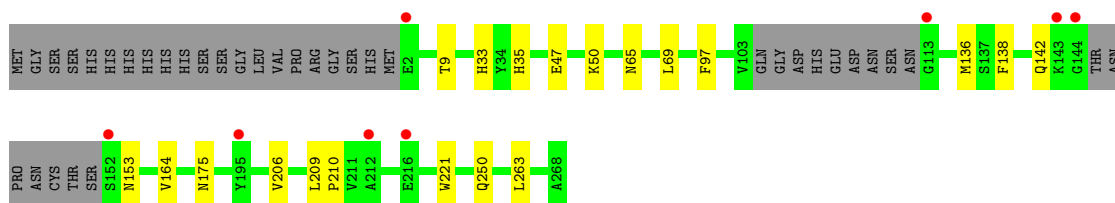
• Molecule 1: Pteridine reductase

Chain C:



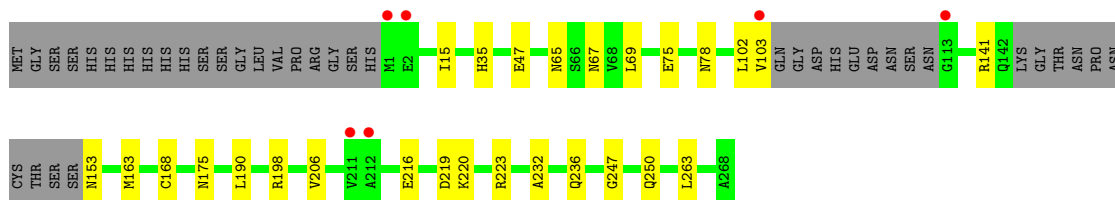
• Molecule 1: Pteridine reductase

Chain D:



• Molecule 2: Pteridine reductase

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.73Å 90.99Å 82.82Å 90.00° 115.79° 90.00°	Depositor
Resolution (Å)	22.38 – 1.60 22.38 – 1.60	Depositor EDS
% Data completeness (in resolution range)	95.7 (22.38-1.60) 95.5 (22.38-1.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.88 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.118 , 0.150 0.136 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	13.6	Xtriage
Anisotropy	0.371	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 52.1	EDS
Estimated twinning fraction	0.013 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	4 of 125368 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9481	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.13 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 9.2899e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities 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: GOL, AX4, NA, D1D, NAP, CSX, ACT, DTT

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	0.72	0/2109	0.81	1/2853 (0.0%)
1	C	0.73	0/2037	0.76	0/2756
1	D	0.69	0/2074	0.77	0/2803
2	B	0.75	0/2017	0.79	2/2728 (0.1%)
All	All	0.72	0/8237	0.78	3/11140 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	82	ARG	NE-CZ-NH1	6.23	123.42	120.30
2	B	198	ARG	NE-CZ-NH1	6.15	123.38	120.30
2	B	198	ARG	NE-CZ-NH2	-5.48	117.56	120.30

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	2013	0	2109	37	0
1	C	1962	0	2045	32	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1981	0	2101	20	0
2	B	1947	0	2040	26	0
3	A	12	0	9	1	0
3	B	4	0	3	0	0
3	C	8	0	6	0	0
3	D	4	0	3	0	0
4	A	1	0	0	0	0
5	A	48	0	25	1	0
5	B	48	0	25	2	0
5	C	48	0	25	1	0
5	D	48	0	25	1	0
6	A	16	0	7	3	0
6	B	16	0	7	0	0
6	C	16	0	7	7	0
6	D	16	0	7	4	0
7	A	8	0	9	2	0
7	B	8	0	9	3	0
7	C	8	0	9	0	0
7	D	8	0	9	0	0
8	A	8	0	8	3	0
8	B	8	0	8	7	0
8	C	8	0	8	2	0
8	D	8	0	8	2	0
9	A	6	0	8	0	0
9	B	12	0	16	6	0
10	A	309	0	0	7	0
10	B	304	0	0	17	0
10	C	298	0	0	8	0
10	D	300	0	0	3	0
All	All	9481	0	8536	123	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 7.

The worst 5 of 123 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:223[B]:ARG:NH1	1:A:224[B]:LYS:HG3	1.43	1.32
1:A:223[B]:ARG:NH1	1:A:224[B]:LYS:CG	1.99	1.26
1:A:223[B]:ARG:HH12	1:A:224[B]:LYS:CG	1.48	1.26
2:B:168:CSX:SG	2:B:168:CSX:OD	2.14	1.04
8:B:273[B]:D1D:H4C2	10:B:433:HOH:O	1.61	1.00

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	271/288 (94%)	261 (96%)	10 (4%)	0	100	100
1	C	262/288 (91%)	251 (96%)	11 (4%)	0	100	100
1	D	266/288 (92%)	256 (96%)	10 (4%)	0	100	100
2	B	259/288 (90%)	251 (97%)	8 (3%)	0	100	100
All	All	1058/1152 (92%)	1019 (96%)	39 (4%)	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	225/231 (97%)	223 (99%)	2 (1%)	87	74
1	C	216/231 (94%)	211 (98%)	5 (2%)	63	32
1	D	220/231 (95%)	220 (100%)	0	100	100
2	B	213/230 (93%)	211 (99%)	2 (1%)	87	74
All	All	874/923 (95%)	865 (99%)	9 (1%)	87	70

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	2	GLU
1	C	143	LYS
1	C	67	ASN

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Mol	Chain	Res	Type
2	B	67[A]	ASN
1	C	60	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	67	ASN
1	C	104	GLN
1	C	236	GLN
1	C	60	GLN
1	C	186	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	CSX	B	168	2	4,6,7	7.46	2 (50%)	3,6,8	3.92	2 (66%)

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	CSX	B	168	2	-	0/2/5/7	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	168	CSX	O-C	14.55	1.21	1.11
2	B	168	CSX	CA-C	2.69	1.53	1.48

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	168	CSX	CA-CB-SG	6.05	119.15	110.82
2	B	168	CSX	C-CA-N	-3.10	110.74	113.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 27 ligands modelled in this entry, 1 is monoatomic - leaving 26 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	ACT	A	269	-	1,3,3	2.24	1 (100%)	0,3,3	0.00	-
3	ACT	A	270	-	1,3,3	0.64	0	0,3,3	0.00	-
3	ACT	A	271	-	1,3,3	0.68	0	0,3,3	0.00	-
5	NAP	A	273	-	52,52,52	1.19	4 (7%)	80,80,80	1.73	13 (16%)
6	AX4	A	274[A]	-	7,7,17	0.37	0	8,8,23	0.72	0
7	DTT	A	275	-	7,7,7	0.63	0	8,8,8	1.07	1 (12%)
8	D1D	A	276[B]	-	8,8,8	1.18	1 (12%)	10,10,10	1.84	2 (20%)
9	GOL	A	277	-	5,5,5	0.59	0	5,5,5	1.04	0
3	ACT	B	269	-	1,3,3	0.80	0	0,3,3	0.00	-
5	NAP	B	270	-	52,52,52	1.30	7 (13%)	80,80,80	1.74	14 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	AX4	B	271[A]	-	7,7,17	0.54	0	8,8,23	0.81	0
7	DTT	B	272	-	7,7,7	0.83	0	8,8,8	1.41	1 (12%)
8	D1D	B	273[B]	-	8,8,8	0.85	0	10,10,10	2.36	1 (10%)
9	GOL	B	274[C]	-	5,5,5	0.60	0	5,5,5	0.19	0
9	GOL	B	275	-	5,5,5	0.34	0	5,5,5	0.62	0
3	ACT	C	269	-	1,3,3	1.48	0	0,3,3	0.00	-
3	ACT	C	270	-	1,3,3	1.95	0	0,3,3	0.00	-
5	NAP	C	271	-	52,52,52	1.22	3 (5%)	80,80,80	1.73	8 (10%)
6	AX4	C	272[A]	-	7,7,17	0.33	0	8,8,23	0.68	0
7	DTT	C	273	-	7,7,7	0.76	0	8,8,8	0.84	0
8	D1D	C	274[B]	-	8,8,8	1.24	1 (12%)	10,10,10	1.76	3 (30%)
3	ACT	D	269	-	1,3,3	0.49	0	0,3,3	0.00	-
5	NAP	D	270	-	52,52,52	1.15	2 (3%)	80,80,80	1.99	14 (17%)
6	AX4	D	271[A]	-	7,7,17	0.35	0	8,8,23	0.68	0
7	DTT	D	272	-	7,7,7	0.67	0	8,8,8	0.88	0
8	D1D	D	273[B]	-	8,8,8	0.87	0	10,10,10	1.54	1 (10%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ACT	A	269	-	-	0/0/0/0	0/0/0/0
3	ACT	A	270	-	-	0/0/0/0	0/0/0/0
3	ACT	A	271	-	-	0/0/0/0	0/0/0/0
5	NAP	A	273	-	-	0/35/67/67	0/3/5/5
6	AX4	A	274[A]	-	-	0/0/0/4	0/1/1/2
7	DTT	A	275	-	2/2/2/2	0/8/8/8	0/0/0/0
8	D1D	A	276[B]	-	-	0/0/11/11	0/1/1/1
9	GOL	A	277	-	-	0/4/4/4	0/0/0/0
3	ACT	B	269	-	-	0/0/0/0	0/0/0/0
5	NAP	B	270	-	-	0/35/67/67	0/3/5/5
6	AX4	B	271[A]	-	-	0/0/0/4	0/1/1/2
7	DTT	B	272	-	-	0/8/8/8	0/0/0/0
8	D1D	B	273[B]	-	-	0/0/11/11	0/1/1/1
9	GOL	B	274[C]	-	-	0/4/4/4	0/0/0/0
9	GOL	B	275	-	-	0/4/4/4	0/0/0/0
3	ACT	C	269	-	-	0/0/0/0	0/0/0/0
3	ACT	C	270	-	-	0/0/0/0	0/0/0/0
5	NAP	C	271	-	-	0/35/67/67	0/3/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	AX4	C	272[A]	-	-	0/0/0/4	0/1/1/2
7	DTT	C	273	-	-	0/8/8/8	0/0/0/0
8	D1D	C	274[B]	-	-	0/0/11/11	0/1/1/1
3	ACT	D	269	-	-	0/0/0/0	0/0/0/0
5	NAP	D	270	-	-	0/35/67/67	0/3/5/5
6	AX4	D	271[A]	-	-	0/0/0/4	0/1/1/2
7	DTT	D	272	-	-	0/8/8/8	0/0/0/0
8	D1D	D	273[B]	-	-	0/0/11/11	0/1/1/1

The worst 5 of 19 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	270	NAP	O7N-C7N	5.40	1.36	1.24
5	C	271	NAP	O7N-C7N	4.87	1.35	1.24
5	A	273	NAP	O7N-C7N	4.16	1.33	1.24
5	B	270	NAP	O7N-C7N	4.10	1.33	1.24
5	C	271	NAP	C2B-C1B	3.45	1.58	1.52

The worst 5 of 58 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	270	NAP	N3A-C2A-N1A	-11.11	119.42	128.71
5	C	271	NAP	N3A-C2A-N1A	-8.39	121.70	128.71
5	A	273	NAP	N3A-C2A-N1A	-8.29	121.78	128.71
5	B	270	NAP	N3A-C2A-N1A	-7.51	122.43	128.71
8	B	273[B]	D1D	C4-S4-S1	6.56	105.42	98.70

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
7	A	275	DTT	C2
7	A	275	DTT	C3

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, 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	255/288 (88%)	-0.35	9 (3%) 42 40	7, 10, 24, 54	0
1	C	252/288 (87%)	-0.33	9 (3%) 41 39	8, 11, 27, 45	0
1	D	251/288 (87%)	-0.35	8 (3%) 45 44	9, 11, 25, 53	0
2	B	249/288 (86%)	-0.51	6 (2%) 56 56	7, 9, 18, 48	0
All	All	1007/1152 (87%)	-0.38	32 (3%) 47 44	7, 11, 24, 54	0

The worst 5 of 32 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	144	GLY	7.3
1	D	113	GLY	7.1
1	C	151	SER	6.8
1	A	150	THR	6.6
1	A	144	GLY	6.5

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

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	CSX	B	168	7/8	0.09	2.71	11,13,26,28	0

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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	ACT	A	271	4/4	0.16	10.80	21,26,28,29	0
3	ACT	A	270	4/4	0.18	9.13	24,24,25,27	0
3	ACT	D	269	4/4	0.17	8.40	23,23,23,26	0
7	DTT	B	272	8/8	0.17	7.83	25,45,53,57	0
3	ACT	C	269	4/4	0.16	6.64	21,21,22,24	0
8	D1D	D	273[B]	8/8	0.19	6.60	26,28,30,31	8
3	ACT	B	269	4/4	0.19	4.83	23,23,24,26	0
9	GOL	B	274[C]	6/6	0.12	4.44	15,23,28,31	6
8	D1D	C	274[B]	8/8	0.14	4.37	17,21,25,27	8
8	D1D	A	276[B]	8/8	0.14	3.75	27,31,34,34	8
9	GOL	B	275	6/6	0.16	3.63	25,29,30,34	0
8	D1D	B	273[B]	8/8	0.13	3.45	16,18,20,21	8
7	DTT	A	275	8/8	0.15	2.65	29,46,52,59	0
6	AX4	A	274[A]	7/16	0.10	1.92	7,9,10,11	7
6	AX4	C	272[A]	7/16	0.08	1.64	13,16,17,18	7
6	AX4	B	271[A]	7/16	0.08	1.53	11,14,15,18	7
3	ACT	C	270	4/4	0.10	1.35	15,17,18,18	0
6	AX4	D	271[A]	7/16	0.10	1.28	10,13,14,15	7
7	DTT	C	273	8/8	0.11	1.23	27,42,49,56	0
7	DTT	D	272	8/8	0.11	0.86	27,46,52,58	0
9	GOL	A	277	6/6	0.07	0.08	15,18,19,23	0
5	NAP	A	273	48/48	0.06	-0.25	6,9,10,13	0
5	NAP	D	270	48/48	0.05	-0.46	8,10,12,13	0
5	NAP	B	270	48/48	0.05	-0.75	5,8,10,11	0
3	ACT	A	269	4/4	0.04	-0.86	11,12,12,16	0
5	NAP	C	271	48/48	0.05	-1.02	7,10,11,12	0
4	NA	A	272	1/1	0.04	-2.43	28,28,28,28	0

6.5 Other polymers ⓘ

There are no such residues in this entry.