



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:14 PM GMT

PDB ID : 1O2D
Title : Crystal structure of Alcohol dehydrogenase, iron-containing (TM0920) from *Thermotoga maritima* at 1.30 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2003-02-27
Resolution : 1.30 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

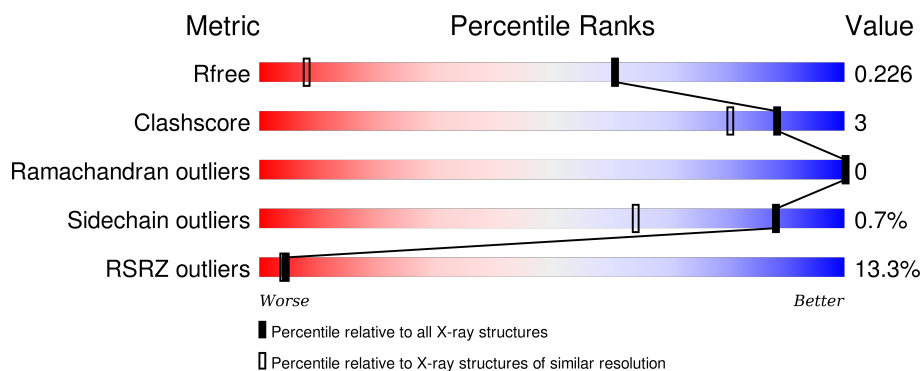
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.30 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	1475 (1.34-1.26)
Clashscore	102246	1031 (1.32-1.28)
Ramachandran outliers	100387	1504 (1.34-1.26)
Sidechain outliers	100360	1503 (1.34-1.26)
RSRZ outliers	91569	1476 (1.34-1.26)

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	371	<div> <div>11%</div> <div>91%</div> <div>5% . .</div> </div>
1	B	371	<div> <div>14%</div> <div>92%</div> <div>5% . .</div> </div>

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
4	TRS	A	2801	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6604 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 Alcohol dehydrogenase, iron-containing.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	359	Total	C	N	O	S	Se	0	12	0
			2805	1797	461	533	2	12			
1	B	360	Total	C	N	O	S	Se	0	10	0
			2791	1789	460	527	2	13			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	LEADER SEQUENCE	UNP Q9X022
A	-10	GLY	-	LEADER SEQUENCE	UNP Q9X022
A	-9	SER	-	LEADER SEQUENCE	UNP Q9X022
A	-8	ASP	-	LEADER SEQUENCE	UNP Q9X022
A	-7	LYS	-	LEADER SEQUENCE	UNP Q9X022
A	-6	ILE	-	LEADER SEQUENCE	UNP Q9X022
A	-5	HIS	-	LEADER SEQUENCE	UNP Q9X022
A	-4	HIS	-	LEADER SEQUENCE	UNP Q9X022
A	-3	HIS	-	LEADER SEQUENCE	UNP Q9X022
A	-2	HIS	-	LEADER SEQUENCE	UNP Q9X022
A	-1	HIS	-	LEADER SEQUENCE	UNP Q9X022
A	0	HIS	-	LEADER SEQUENCE	UNP Q9X022
A	1	VAL	-	LEADER SEQUENCE	UNP Q9X022
A	6	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
A	76	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
A	99	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
A	163	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
A	178	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
A	215	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
A	236	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
A	245	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
A	276	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
A	282	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
A	285	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
B	-11	MET	-	LEADER SEQUENCE	UNP Q9X022

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	GLY	-	LEADER SEQUENCE	UNP Q9X022
B	-9	SER	-	LEADER SEQUENCE	UNP Q9X022
B	-8	ASP	-	LEADER SEQUENCE	UNP Q9X022
B	-7	LYS	-	LEADER SEQUENCE	UNP Q9X022
B	-6	ILE	-	LEADER SEQUENCE	UNP Q9X022
B	-5	HIS	-	LEADER SEQUENCE	UNP Q9X022
B	-4	HIS	-	LEADER SEQUENCE	UNP Q9X022
B	-3	HIS	-	LEADER SEQUENCE	UNP Q9X022
B	-2	HIS	-	LEADER SEQUENCE	UNP Q9X022
B	-1	HIS	-	LEADER SEQUENCE	UNP Q9X022
B	0	HIS	-	LEADER SEQUENCE	UNP Q9X022
B	1	VAL	-	LEADER SEQUENCE	UNP Q9X022
B	6	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
B	76	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
B	99	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
B	163	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
B	178	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
B	215	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
B	236	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
B	245	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
B	276	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
B	282	MSE	MET	MODIFIED RESIDUE	UNP Q9X022
B	285	MSE	MET	MODIFIED RESIDUE	UNP Q9X022

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

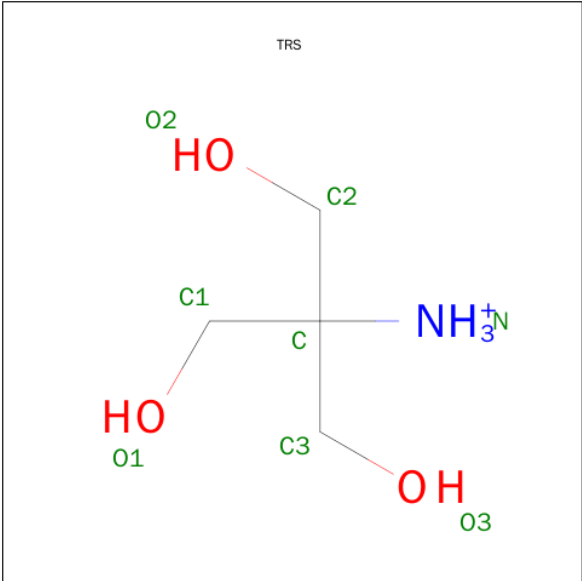
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Fe 1 1	0	0
2	A	1	Total Fe 1 1	0	0

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



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

- Molecule 4 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: C₄H₁₂NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	1
			10	5	1	4		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			8	4	1	3		

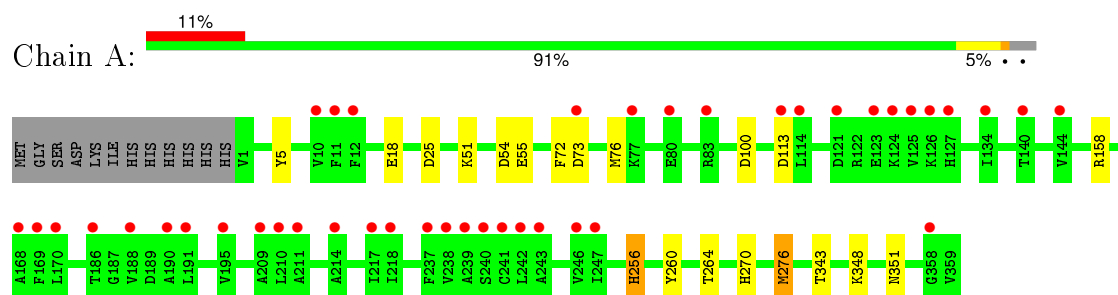
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	479	Total	O	0	0
			479	479		
5	B	413	Total	O	0	0
			413	413		

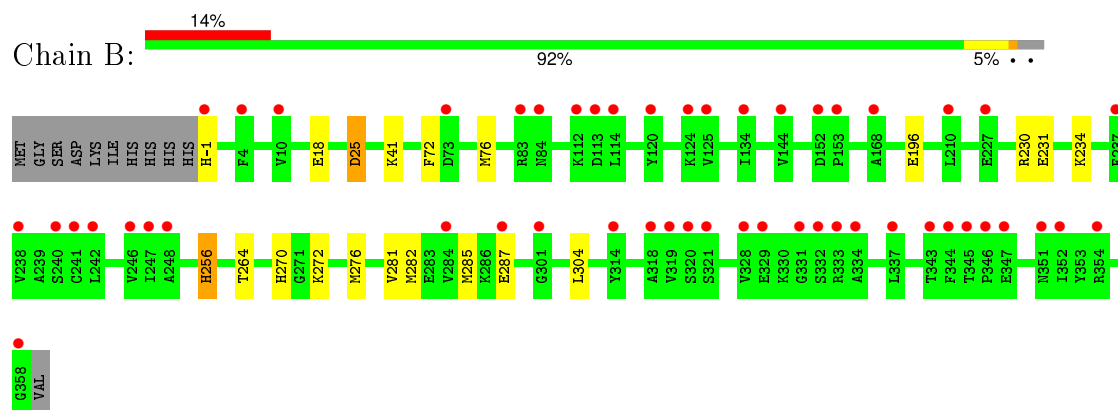
3 Residue-property plots [i](#)

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: Alcohol dehydrogenase, iron-containing



- Molecule 1: Alcohol dehydrogenase, iron-containing



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.05Å 85.28Å 72.08Å 90.00° 96.29° 90.00°	Depositor
Resolution (Å)	38.19 – 1.30 38.19 – 1.30	Depositor EDS
% Data completeness (in resolution range)	97.8 (38.19-1.30) 97.8 (38.19-1.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 1.30Å)	Xtriage
Refinement program	REFMAC 5.1.955	Depositor
R, R_{free}	0.137 , 0.170 0.204 , 0.226	Depositor DCC
R_{free} test set	8371 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	13.1	Xtriage
Anisotropy	0.179	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 50.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 167310 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6604	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.41% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.63	1/2903 (0.0%)	0.76	3/3906 (0.1%)
1	B	0.58	0/2886	0.74	1/3887 (0.0%)
All	All	0.61	1/5789 (0.0%)	0.75	4/7793 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	276	MSE	SE-CE	-7.78	1.49	1.95

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	25	ASP	CB-CG-OD2	6.19	123.87	118.30
1	A	25	ASP	CB-CG-OD1	5.88	123.59	118.30
1	A	113	ASP	CB-CG-OD2	5.86	123.57	118.30
1	A	158	ARG	NE-CZ-NH2	-5.68	117.46	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2805	0	2819	14	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2791	0	2763	14	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	48	0	24	2	0
3	B	48	0	25	2	0
4	A	8	0	12	0	0
4	B	10	0	6	0	0
5	A	479	0	0	7	1
5	B	413	0	0	6	1
All	All	6604	0	5649	29	2

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 3.

All (29) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:THR:HG22	5:A:3015:HOH:O	1.64	0.96
1:B:287:GLU:O	5:B:2833:HOH:O	1.92	0.86
1:A:351[B]:ASN:ND2	5:A:2920:HOH:O	2.15	0.77
1:A:351[B]:ASN:OD1	5:A:3041:HOH:O	2.11	0.68
1:A:256[B]:HIS:NE2	1:A:260:TYR:OH	2.27	0.66
1:A:18:GLU:OE2	5:A:3037:HOH:O	2.16	0.61
1:B:281:VAL:HG12	1:B:285:MSE:HE3	1.84	0.60
1:B:264:THR:HG22	5:B:3205:HOH:O	2.02	0.59
1:A:54:ASP:CG	5:A:3128:HOH:O	2.40	0.59
1:B:270:HIS:CD2	3:B:1801:NAP:H6N	2.40	0.56
1:A:270:HIS:CD2	3:A:1800:NAP:H6N	2.40	0.56
1:B:41:LYS:C	5:B:3046:HOH:O	2.44	0.55
1:A:256[B]:HIS:CD2	1:A:260:TYR:OH	2.61	0.54
1:B:231:GLU:OE2	1:B:234:LYS:NZ	2.39	0.54
1:B:76:MSE:HE2	5:B:3007:HOH:O	2.10	0.51
1:B:282[A]:MSE:SE	1:B:304:LEU:HD23	2.61	0.50
1:A:72:PHE:HB3	1:A:76:MSE:HE3	1.95	0.47
1:B:18:GLU:HG3	5:B:3120:HOH:O	2.14	0.47
1:B:281:VAL:HG12	1:B:285:MSE:CE	2.47	0.45
1:B:196:GLU:OE2	1:B:256[A]:HIS:ND1	2.43	0.45
1:B:272:LYS:O	1:B:276:MSE:HG3	2.17	0.44
1:A:51:LYS:HG2	5:A:3247:HOH:O	2.17	0.44
1:B:72:PHE:HB3	1:B:76:MSE:HE3	1.99	0.44
1:A:55:GLU:HB2	5:A:3247:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:ASP:OD2	3:A:1800:NAP:H2N	2.19	0.42
3:B:1801:NAP:C5B	5:B:3208:HOH:O	2.68	0.42
1:A:343:THR:O	1:A:348:LYS:HD2	2.21	0.41

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:-1:HIS:N	5:A:2941:HOH:O[2_647]	2.15	0.05
1:A:5:TYR:OH	5:B:3213:HOH:O[2_657]	2.18	0.02

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	369/371 (100%)	361 (98%)	8 (2%)	0	100	100
1	B	368/371 (99%)	361 (98%)	7 (2%)	0	100	100
All	All	737/742 (99%)	722 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	311/308 (101%)	308 (99%)	3 (1%)	82	51
1	B	302/308 (98%)	299 (99%)	3 (1%)	82	51
All	All	613/616 (100%)	607 (99%)	6 (1%)	88	51

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	ASP
1	A	256[A]	HIS
1	A	256[B]	HIS
1	B	25	ASP
1	B	256[A]	HIS
1	B	256[B]	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	-1	HIS

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](#)

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	NAP	A	1800	-	42,52,52	1.42	3 (7%)	54,80,80	2.00	7 (12%)
4	TRS	A	2801	-	7,7,7	0.77	0	9,9,9	1.27	1 (11%)
3	NAP	B	1801	-	42,52,52	1.52	3 (7%)	54,80,80	1.55	6 (11%)
4	TRS	B	2800[A]	-	7,7,7	0.51	0	9,9,9	0.56	0
4	TRS	B	2800[B]	-	7,7,7	0.56	0	9,9,9	0.92	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
3	NAP	A	1800	-	-	0/27/67/67	0/5/5/5
4	TRS	A	2801	-	-	0/9/9/9	0/0/0/0
3	NAP	B	1801	-	-	0/27/67/67	0/5/5/5
4	TRS	B	2800[A]	-	-	0/9/9/9	0/0/0/0
4	TRS	B	2800[B]	-	-	0/9/9/9	0/0/0/0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1800	NAP	C2A-N1A	2.05	1.37	1.33
3	B	1801	NAP	C2A-N1A	2.16	1.38	1.33
3	A	1800	NAP	C2A-N3A	2.91	1.37	1.32
3	B	1801	NAP	C2A-N3A	3.40	1.38	1.32
3	A	1800	NAP	O7N-C7N	7.30	1.39	1.24
3	B	1801	NAP	O7N-C7N	7.56	1.40	1.24

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1800	NAP	N3A-C2A-N1A	-11.57	120.04	128.89
3	B	1801	NAP	N3A-C2A-N1A	-7.34	123.28	128.89
3	B	1801	NAP	O7N-C7N-N7N	-3.70	117.39	122.59
3	A	1800	NAP	O7N-C7N-N7N	-3.58	117.55	122.59
3	A	1800	NAP	C1B-N9A-C4A	-2.29	123.48	126.94
3	A	1800	NAP	O3-PA-O5B	-2.12	97.32	102.94
3	B	1801	NAP	C3N-C7N-N7N	2.10	120.12	117.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1801	NAP	O2A-PA-O3	2.26	115.36	105.09
3	A	1800	NAP	C2A-N1A-C6A	2.33	122.93	118.77
4	A	2801	TRS	C2-C-N	2.62	112.85	108.09
3	B	1801	NAP	O7N-C7N-C3N	2.67	122.50	119.59
3	A	1800	NAP	O7N-C7N-C3N	3.15	123.02	119.59
3	A	1800	NAP	O4D-C1D-N1N	3.62	112.10	108.13
3	B	1801	NAP	O4D-C1D-N1N	3.66	112.16	108.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1800	NAP	2	0
3	B	1801	NAP	2	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	348/371 (93%)	0.64	42 (12%) 6 5	5, 9, 19, 26	0
1	B	349/371 (94%)	0.73	51 (14%) 3 3	5, 10, 19, 25	0
All	All	697/742 (93%)	0.69	93 (13%) 4 4	5, 10, 19, 26	0

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	113	ASP	5.2
1	A	125	VAL	4.9
1	A	73	ASP	4.3
1	A	241	CYS	4.0
1	B	344	PHE	4.0
1	B	343	THR	3.9
1	B	345	THR	3.9
1	A	113	ASP	3.8
1	A	80	GLU	3.7
1	A	242	LEU	3.6
1	A	77	LYS	3.6
1	A	114	LEU	3.5
1	A	238	VAL	3.4
1	B	346	PRO	3.4
1	B	318	ALA	3.4
1	B	358	GLY	3.4
1	A	123	GLU	3.3
1	A	247	ILE	3.3
1	A	246	VAL	3.2
1	B	114	LEU	3.2
1	A	218	ILE	3.2
1	B	83	ARG	3.1
1	B	144	VAL	3.1
1	B	319	VAL	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	210	LEU	3.0
1	B	347	GLU	3.0
1	A	144	VAL	3.0
1	B	125	VAL	3.0
1	B	351	ASN	2.9
1	B	112	LYS	2.9
1	B	321	SER	2.9
1	B	246	VAL	2.9
1	A	239	ALA	2.8
1	A	358	GLY	2.8
1	B	248	ALA	2.8
1	A	237	PHE	2.8
1	A	190	ALA	2.7
1	B	242	LEU	2.7
1	B	153	PRO	2.7
1	B	-1	HIS	2.7
1	B	331	GLY	2.7
1	A	170	LEU	2.6
1	B	134	ILE	2.6
1	B	10	VAL	2.6
1	B	237	PHE	2.6
1	B	241	CYS	2.6
1	A	126	LYS	2.6
1	A	214	ALA	2.6
1	B	354	ARG	2.5
1	A	243	ALA	2.5
1	B	287	GLU	2.5
1	B	332	SER	2.5
1	A	124	LYS	2.4
1	A	10	VAL	2.4
1	A	11	PHE	2.4
1	A	12	PHE	2.4
1	A	217	ILE	2.4
1	B	329	GLU	2.4
1	B	333	ARG	2.4
1	B	247	ILE	2.4
1	B	168	ALA	2.4
1	A	195	VAL	2.4
1	B	210	LEU	2.3
1	B	84	ASN	2.3
1	A	211	ALA	2.3
1	A	121	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	134	ILE	2.3
1	B	120	TYR	2.3
1	A	191	LEU	2.3
1	B	4	PHE	2.3
1	B	320	SER	2.3
1	B	238	VAL	2.3
1	A	168	ALA	2.3
1	A	140	THR	2.2
1	B	73	ASP	2.2
1	B	124	LYS	2.2
1	A	169	PHE	2.2
1	B	284	VAL	2.2
1	B	352	ILE	2.2
1	B	240[A]	SER	2.1
1	A	188	VAL	2.1
1	B	314	TYR	2.1
1	A	186	THR	2.1
1	A	127	HIS	2.1
1	B	227	GLU	2.1
1	B	328	VAL	2.1
1	B	301	GLY	2.1
1	B	337	LEU	2.1
1	B	152	ASP	2.1
1	A	209	ALA	2.0
1	B	334	ALA	2.0
1	A	240[A]	SER	2.0
1	A	83	ARG	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
4	TRS	A	2801	8/8	0.73	0.24	3.80	29,30,30,35	0
3	NAP	B	1801	48/48	0.95	0.08	-0.75	9,14,23,27	0
3	NAP	A	1800	48/48	0.97	0.07	-1.16	8,12,22,24	0
2	FE	A	900	1/1	0.99	0.10	-1.17	12,12,12,12	1
4	TRS	B	2800[A]	8/8	0.92	0.10	-	14,17,19,22	2
2	FE	B	900	1/1	0.99	0.07	-	10,10,10,10	1
4	TRS	B	2800[B]	8/8	0.92	0.10	-	15,17,19,22	2

6.5 Other polymers

There are no such residues in this entry.