



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:56 PM GMT

PDB ID : 4GO2
Title : Crystal structure of the c-terminal domain of 10'formyltetrahydrofolate dehydrogenase in complex with Thio-NADP
Authors : Tsybovsky, Y.
Deposited on : 2012-08-18
Resolution : 2.28 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

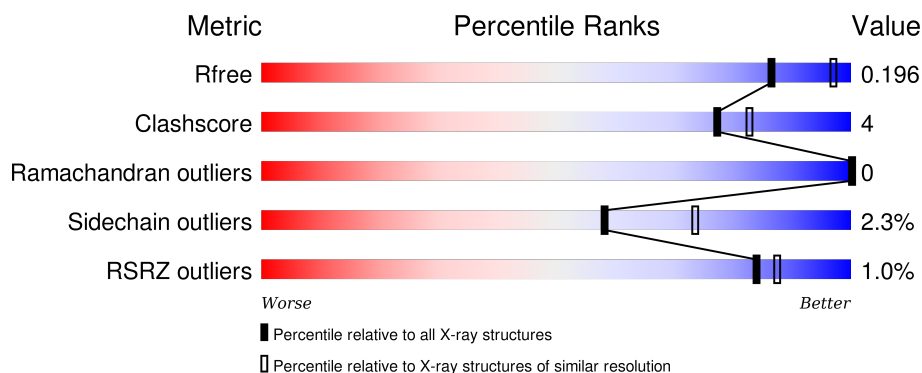
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.28 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	5193 (2.30-2.26)
Clashscore	102246	5929 (2.30-2.26)
Ramachandran outliers	100387	5851 (2.30-2.26)
Sidechain outliers	100360	5850 (2.30-2.26)
RSRZ outliers	91569	5204 (2.30-2.26)

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	517	<div> <div>%</div> <div> <div></div> <div>88%</div> <div>8%</div> <div>• •</div> </div> </div>
1	B	517	<div> <div>2%</div> <div> <div></div> <div>90%</div> <div>6%</div> <div>• •</div> </div> </div>
1	C	517	<div> <div></div> <div> <div></div> <div>87%</div> <div>9%</div> <div>• •</div> </div> </div>
1	D	517	<div> <div>%</div> <div> <div></div> <div>87%</div> <div>9%</div> <div>•</div> </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
2	SO4	A	3005	-	-	-	X
2	SO4	A	3006	-	-	-	X
2	SO4	B	1006	-	-	-	X
2	SO4	D	1004	-	-	-	X
3	GOL	A	3008	-	-	-	X
3	GOL	C	1008	-	-	-	X
3	GOL	D	1007	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 17137 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 Cytosolic 10-formyltetrahydrofolate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	498	Total	C	N	O	S	26	22	0
			3930	2497	674	737	22			
1	B	498	Total	C	N	O	S	16	23	0
			3928	2494	675	737	22			
1	C	498	Total	C	N	O	S	17	20	0
			3914	2486	676	730	22			
1	D	498	Total	C	N	O	S	15	21	0
			3922	2492	672	736	22			

There are 44 discrepancies between the modelled and reference sequences:

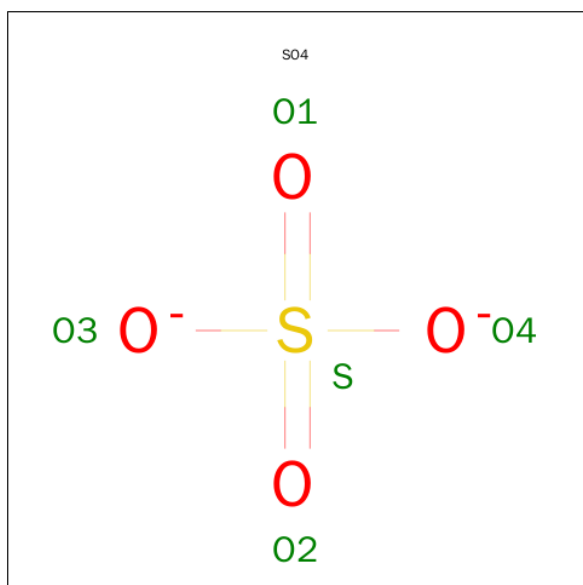
Chain	Residue	Modelled	Actual	Comment	Reference
A	386	MET	-	EXPRESSION TAG	UNP P28037
A	387	ARG	-	EXPRESSION TAG	UNP P28037
A	388	GLY	-	EXPRESSION TAG	UNP P28037
A	389	SER	-	EXPRESSION TAG	UNP P28037
A	390	HIS	-	EXPRESSION TAG	UNP P28037
A	391	HIS	-	EXPRESSION TAG	UNP P28037
A	392	HIS	-	EXPRESSION TAG	UNP P28037
A	393	HIS	-	EXPRESSION TAG	UNP P28037
A	394	HIS	-	EXPRESSION TAG	UNP P28037
A	395	THR	-	EXPRESSION TAG	UNP P28037
A	396	THR	-	EXPRESSION TAG	UNP P28037
B	386	MET	-	EXPRESSION TAG	UNP P28037
B	387	ARG	-	EXPRESSION TAG	UNP P28037
B	388	GLY	-	EXPRESSION TAG	UNP P28037
B	389	SER	-	EXPRESSION TAG	UNP P28037
B	390	HIS	-	EXPRESSION TAG	UNP P28037
B	391	HIS	-	EXPRESSION TAG	UNP P28037
B	392	HIS	-	EXPRESSION TAG	UNP P28037
B	393	HIS	-	EXPRESSION TAG	UNP P28037
B	394	HIS	-	EXPRESSION TAG	UNP P28037
B	395	THR	-	EXPRESSION TAG	UNP P28037

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Chain	Residue	Modelled	Actual	Comment	Reference
B	396	THR	-	EXPRESSION TAG	UNP P28037
C	386	MET	-	EXPRESSION TAG	UNP P28037
C	387	ARG	-	EXPRESSION TAG	UNP P28037
C	388	GLY	-	EXPRESSION TAG	UNP P28037
C	389	SER	-	EXPRESSION TAG	UNP P28037
C	390	HIS	-	EXPRESSION TAG	UNP P28037
C	391	HIS	-	EXPRESSION TAG	UNP P28037
C	392	HIS	-	EXPRESSION TAG	UNP P28037
C	393	HIS	-	EXPRESSION TAG	UNP P28037
C	394	HIS	-	EXPRESSION TAG	UNP P28037
C	395	THR	-	EXPRESSION TAG	UNP P28037
C	396	THR	-	EXPRESSION TAG	UNP P28037
D	386	MET	-	EXPRESSION TAG	UNP P28037
D	387	ARG	-	EXPRESSION TAG	UNP P28037
D	388	GLY	-	EXPRESSION TAG	UNP P28037
D	389	SER	-	EXPRESSION TAG	UNP P28037
D	390	HIS	-	EXPRESSION TAG	UNP P28037
D	391	HIS	-	EXPRESSION TAG	UNP P28037
D	392	HIS	-	EXPRESSION TAG	UNP P28037
D	393	HIS	-	EXPRESSION TAG	UNP P28037
D	394	HIS	-	EXPRESSION TAG	UNP P28037
D	395	THR	-	EXPRESSION TAG	UNP P28037
D	396	THR	-	EXPRESSION TAG	UNP P28037

- Molecule 2 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



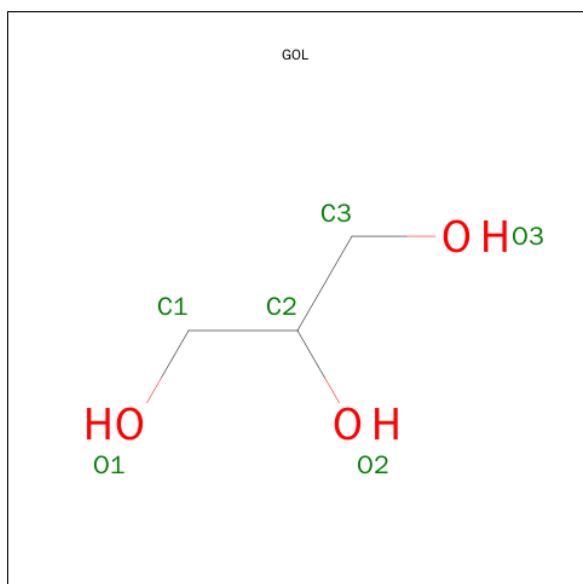
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0

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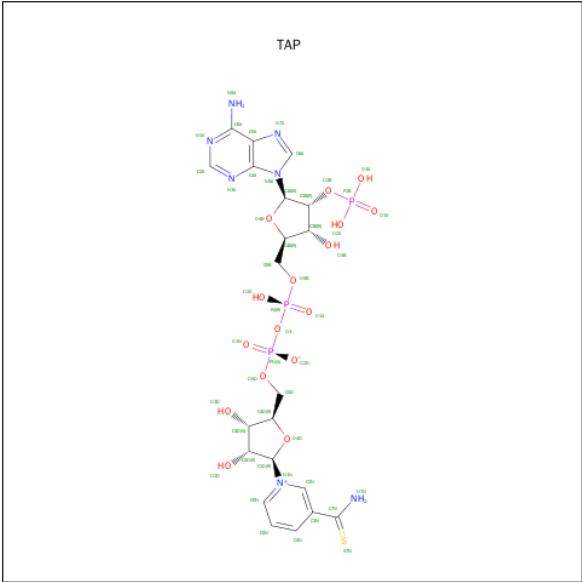
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 4 is 7-THIONICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: TAP) (formula: $C_{21}H_{28}N_7O_{16}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	S	0	1
			64	26	7	25	5	1		
4	B	1	Total	C	N	O	P	S	0	1
			64	26	7	25	5	1		
4	C	1	Total	C	N	O	P	S	0	1
			64	26	7	25	5	1		
4	D	1	Total	C	N	O	P	S	0	1
			64	26	7	25	5	1		

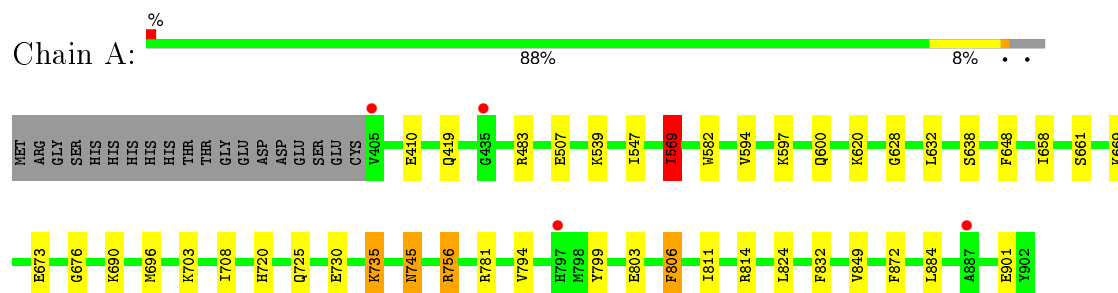
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	248	Total	O	0	0
			248	248		
5	B	247	Total	O	0	0
			247	247		
5	C	283	Total	O	0	0
			283	283		
5	D	245	Total	O	0	0
			245	245		

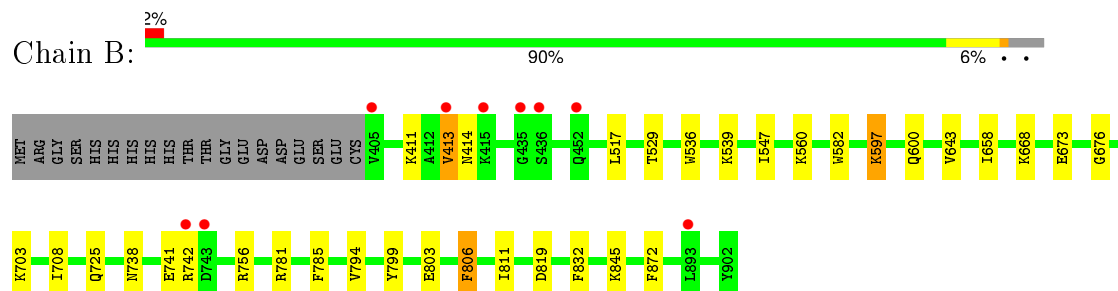
3 Residue-property plots [i](#)

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 ($RSRZ > 2$). 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.

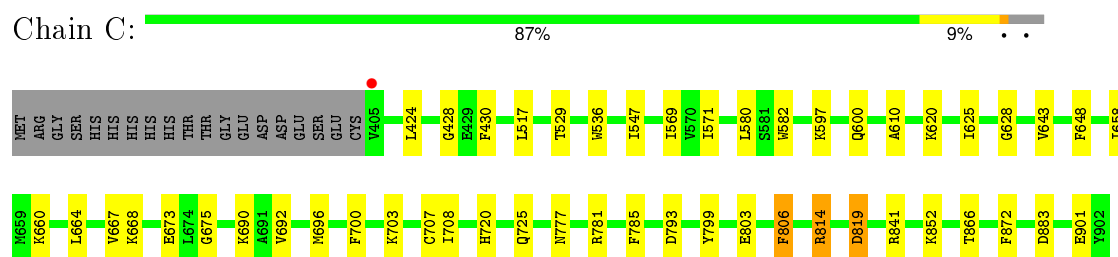
- Molecule 1: Cytosolic 10-formyltetrahydrofolate dehydrogenase



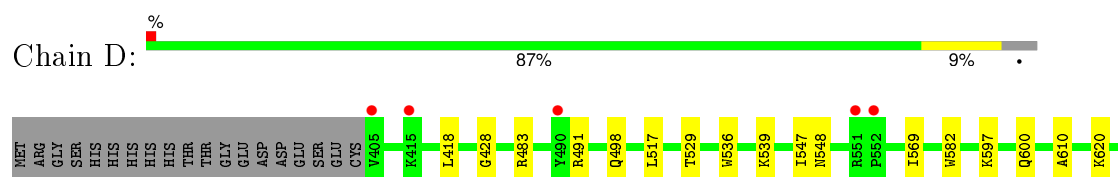
- Molecule 1: Cytosolic 10-formyltetrahydrofolate dehydrogenase

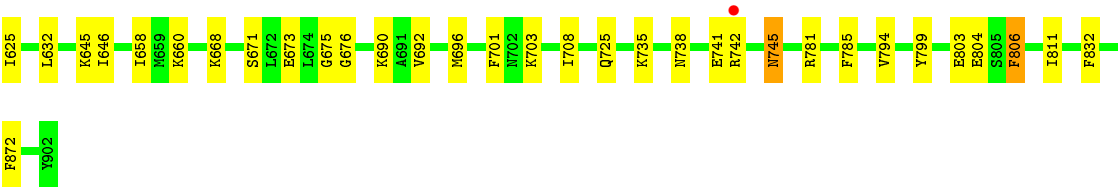


- Molecule 1: Cytosolic 10-formyltetrahydrofolate dehydrogenase



- Molecule 1: Cytosolic 10-formyltetrahydrofolate dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	258.94Å 194.39Å 97.11Å 90.00° 109.02° 90.00°	Depositor
Resolution (Å)	49.88 – 2.28 48.80 – 2.28	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.88-2.28) 99.9 (48.80-2.28)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.89 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.177 , 0.196 0.176 , 0.196	Depositor DCC
R_{free} test set	10330 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	29.2	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 44.2	EDS
Estimated twinning fraction	0.010 for -h-2*k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 205441 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	17137	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.79% 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: GOL, TAP, SO4

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	0.71	4/4093 (0.1%)	0.76	5/5532 (0.1%)
1	B	0.69	4/4102 (0.1%)	0.65	4/5544 (0.1%)
1	C	0.62	0/4076	0.68	3/5509 (0.1%)
1	D	0.62	3/4083 (0.1%)	0.67	5/5519 (0.1%)
All	All	0.66	11/16354 (0.1%)	0.69	17/22104 (0.1%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	725[A]	GLN	CG-CD	-16.23	1.13	1.51
1	B	725[B]	GLN	CG-CD	-16.23	1.13	1.51
1	A	756[A]	ARG	CD-NE	15.36	1.72	1.46
1	A	756[B]	ARG	CD-NE	15.36	1.72	1.46
1	A	725	GLN	CG-CD	-10.80	1.26	1.51
1	A	690	LYS	CG-CD	-8.19	1.24	1.52
1	D	725[A]	GLN	CG-CD	-7.20	1.34	1.51
1	D	725[B]	GLN	CG-CD	-7.20	1.34	1.51
1	D	690	LYS	CG-CD	-7.03	1.28	1.52
1	B	560[A]	LYS	CG-CD	-5.62	1.33	1.52
1	B	560[B]	LYS	CG-CD	-5.62	1.33	1.52

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	756[A]	ARG	CD-NE-CZ	17.79	148.50	123.60
1	A	756[B]	ARG	CD-NE-CZ	17.79	148.50	123.60
1	A	756[A]	ARG	CG-CD-NE	11.94	136.87	111.80
1	A	756[B]	ARG	CG-CD-NE	11.94	136.87	111.80
1	B	725[A]	GLN	CB-CG-CD	7.90	132.14	111.60
1	B	725[B]	GLN	CB-CG-CD	7.90	132.14	111.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	725[A]	GLN	CG-CD-OE1	7.71	137.02	121.60
1	D	725[B]	GLN	CG-CD-OE1	7.71	137.02	121.60
1	C	690	LYS	CB-CG-CD	6.92	129.59	111.60
1	D	725[A]	GLN	CG-CD-NE2	-6.89	100.16	116.70
1	D	725[B]	GLN	CG-CD-NE2	-6.89	100.16	116.70
1	C	814	ARG	NE-CZ-NH2	-6.51	117.05	120.30
1	B	725[A]	GLN	CG-CD-OE1	-6.33	108.95	121.60
1	B	725[B]	GLN	CG-CD-OE1	-6.33	108.95	121.60
1	A	569	ILE	CG1-CB-CG2	-5.47	99.36	111.40
1	C	690	LYS	CG-CD-CE	-5.42	95.64	111.90
1	D	690	LYS	CG-CD-CE	-5.01	96.86	111.90

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3930	0	3941	26	0
1	B	3928	0	3938	21	0
1	C	3914	0	3927	40	0
1	D	3922	0	3933	32	0
2	A	35	0	0	1	0
2	B	40	0	0	0	0
2	C	35	0	0	1	0
2	D	30	0	0	0	0
3	A	6	0	8	0	0
3	B	6	0	8	0	0
3	C	6	0	8	1	0
3	D	6	0	8	0	0
4	A	64	0	22	2	0
4	B	64	0	22	4	0
4	C	64	0	22	6	0
4	D	64	0	22	2	0
5	A	248	0	0	3	0
5	B	247	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	283	0	0	4	0
5	D	245	0	0	1	0
All	All	17137	0	15859	118	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 (118) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:841[B]:ARG:HH11	1:C:841[B]:ARG:HG2	1.24	1.03
1:A:597:LYS:HE2	1:A:628:GLY:O	1.58	1.01
1:C:597:LYS:HE2	1:C:628:GLY:O	1.69	0.92
4:B:1010[B]:TAP:H6N	4:B:1010[B]:TAP:H52N	1.55	0.87
1:C:692:VAL:O	1:C:696:MET:HG3	1.75	0.85
1:A:806:PHE:CE1	4:A:3009[A]:TAP:H2D	2.22	0.74
1:D:708:ILE:CG2	1:D:872:PHE:HE1	2.00	0.74
1:C:841[B]:ARG:NH1	1:C:841[B]:ARG:HG2	2.00	0.73
1:B:799:TYR:CE2	1:B:803[B]:GLU:HG3	2.25	0.72
1:A:799:TYR:CE2	1:A:803[A]:GLU:HG3	2.27	0.68
1:C:675:GLY:HA2	4:C:1009[B]:TAP:N7N	2.09	0.67
1:D:428:GLY:HA3	1:D:620[B]:LYS:HD2	1.76	0.66
1:A:735:LYS:H	1:A:745:ASN:HD21	1.44	0.66
1:D:735:LYS:H	1:D:745:ASN:HD21	1.43	0.65
1:D:708:ILE:HG21	1:D:872:PHE:HE1	1.61	0.65
1:A:600:GLN:NE2	5:A:3264:HOH:O	2.24	0.65
1:D:708:ILE:CG2	1:D:872:PHE:CE1	2.81	0.63
1:C:806:PHE:CE1	4:C:1009[A]:TAP:H2D	2.34	0.62
1:D:806:PHE:CE1	4:D:1008[A]:TAP:H2D	2.34	0.62
1:C:707[B]:CYS:HB3	4:C:1009[B]:TAP:S7N	2.39	0.62
1:C:610:ALA:HB2	1:C:625:ILE:HD12	1.81	0.61
1:D:676:GLY:HA2	1:D:832[A]:PHE:CD1	2.35	0.61
1:B:597:LYS:O	1:B:597:LYS:HG3	2.00	0.61
1:C:799:TYR:CE2	1:C:803[A]:GLU:HG3	2.34	0.61
4:B:1010[B]:TAP:H6N	4:B:1010[B]:TAP:C5D	2.28	0.60
1:C:428:GLY:HA3	1:C:620[B]:LYS:HG3	1.82	0.60
1:B:806:PHE:CE1	4:B:1010[A]:TAP:H2D	2.37	0.59
1:C:708:ILE:CG2	1:C:872:PHE:HE1	2.16	0.59
1:C:571:ILE:HD12	1:C:580:LEU:HA	1.84	0.59
1:B:756[B]:ARG:HG3	1:B:756[B]:ARG:HH11	1.67	0.59
1:A:410:GLU:OE1	1:A:419[B]:GLN:NE2	2.30	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:696:MET:CE	1:C:700:PHE:HE2	2.18	0.57
1:C:696:MET:HE2	1:C:700:PHE:HE2	1.68	0.57
1:B:600:GLN:NE2	5:B:1234:HOH:O	2.38	0.56
1:C:696:MET:HE2	1:C:700:PHE:CE2	2.40	0.56
1:C:675:GLY:CA	4:C:1009[B]:TAP:N7N	2.70	0.55
1:D:675:GLY:HA2	4:D:1008[B]:TAP:N7N	2.22	0.55
1:C:852:LYS:NZ	2:C:1001:SO4:O3	2.34	0.55
1:D:610:ALA:HB2	1:D:625:ILE:HD12	1.89	0.55
1:D:735:LYS:N	1:D:745:ASN:HD21	2.05	0.55
1:C:600:GLN:NE2	5:C:1237:HOH:O	2.40	0.54
1:D:738:ASN:HB3	1:D:741[A]:GLU:HG3	1.89	0.54
1:D:645:LYS:NZ	1:D:671[B]:SER:OG	2.36	0.54
1:A:735:LYS:N	1:A:745:ASN:HD21	2.07	0.53
1:A:735:LYS:H	1:A:745:ASN:ND2	2.07	0.53
1:D:781[B]:ARG:HD2	1:D:785:PHE:CD2	2.44	0.53
1:C:692:VAL:O	1:C:696:MET:CG	2.54	0.52
1:B:414:ASN:OD1	1:B:742:ARG:NH2	2.39	0.52
1:D:708:ILE:HG21	1:D:872:PHE:CE1	2.42	0.52
1:C:708:ILE:CG2	1:C:872:PHE:CE1	2.93	0.52
1:A:648:PHE:CD1	1:A:658:ILE:HD12	2.44	0.52
1:C:648:PHE:CD1	1:C:658:ILE:HD12	2.44	0.52
1:B:708:ILE:CG2	1:B:872:PHE:HE1	2.23	0.51
1:A:483:ARG:HB3	1:D:548:ASN:HD21	1.75	0.51
1:A:669:LYS:NZ	2:A:3007:SO4:O1	2.36	0.51
1:C:696:MET:CE	1:C:700:PHE:CE2	2.94	0.50
1:B:411:LYS:HG2	1:B:413:VAL:HG22	1.94	0.50
1:D:692:VAL:O	1:D:696:MET:HG2	2.10	0.50
1:D:517:LEU:HD11	1:D:701:PHE:CZ	2.47	0.49
1:C:708:ILE:HG21	1:C:872:PHE:HE1	1.77	0.49
1:D:799:TYR:CE2	1:D:803[A]:GLU:HG3	2.47	0.49
1:C:643:VAL:O	1:C:668:LYS:HE3	2.12	0.48
5:B:1109:HOH:O	1:D:483:ARG:HD3	2.13	0.48
1:A:676:GLY:HA2	1:A:832[A]:PHE:CG	2.47	0.48
1:D:632:LEU:HD23	1:D:632:LEU:C	2.34	0.48
1:B:756[B]:ARG:NH1	1:B:756[B]:ARG:HG3	2.26	0.48
1:A:620[A]:LYS:HG3	1:A:620[A]:LYS:O	2.14	0.48
1:C:675:GLY:HA2	4:C:1009[B]:TAP:H72N	1.79	0.48
1:B:643:VAL:O	1:B:668:LYS:HE3	2.14	0.47
1:A:720:HIS:CD2	1:A:814:ARG:HG2	2.50	0.47
1:B:799:TYR:CZ	1:B:803[B]:GLU:HG3	2.50	0.47
1:A:708:ILE:CG2	1:A:872:PHE:HE1	2.28	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:819:ASP:C	1:C:819:ASP:OD2	2.52	0.47
1:C:720:HIS:CD2	1:C:814:ARG:HG3	2.50	0.47
1:C:793:ASP:OD2	1:C:814:ARG:NH2	2.48	0.46
1:D:597:LYS:HG3	1:D:597:LYS:O	2.14	0.46
1:A:676:GLY:HA2	1:A:832[A]:PHE:CD1	2.51	0.46
1:A:806:PHE:CZ	4:A:3009[A]:TAP:H2D	2.51	0.45
1:B:806:PHE:CZ	4:B:1010[A]:TAP:H2D	2.51	0.45
1:C:424:LEU:O	1:C:430:PHE:HA	2.17	0.45
1:B:794:VAL:HG21	1:B:811:ILE:HG23	1.98	0.44
1:A:696:MET:SD	1:A:730:GLU:HB2	2.56	0.44
1:D:794:VAL:HG21	1:D:811:ILE:HG23	2.00	0.44
1:B:676:GLY:HA2	1:B:832[B]:PHE:CG	2.52	0.44
1:B:738:ASN:HB3	1:B:741[A]:GLU:HG3	1.99	0.44
1:B:536:TRP:CE2	1:D:539:LYS:HD3	2.53	0.44
1:A:632:LEU:C	1:A:632:LEU:HD23	2.37	0.44
1:D:491[B]:ARG:HH12	1:D:498:GLN:HE22	1.65	0.43
1:C:660:LYS:HG3	1:D:660:LYS:HG3	2.00	0.43
1:A:794:VAL:HG21	1:A:811:ILE:HG23	1.99	0.43
1:C:866:THR:OG1	3:C:1008:GOL:H11	2.18	0.43
1:C:708:ILE:HG21	1:C:872:PHE:CE1	2.53	0.43
1:C:781[B]:ARG:HD2	1:C:785:PHE:CD2	2.54	0.43
1:C:883:ASP:OD1	5:C:1357:HOH:O	2.21	0.43
1:A:901:GLU:OE2	5:A:3141:HOH:O	2.22	0.43
1:A:824:LEU:HD21	1:A:849:VAL:HG13	2.00	0.42
1:D:646:ILE:HG12	1:D:668:LYS:HD3	2.01	0.42
1:B:529:THR:HG21	1:B:582:TRP:HA	2.01	0.42
1:B:539:LYS:HD3	1:D:536:TRP:CE2	2.54	0.42
1:C:597:LYS:O	1:C:597:LYS:HG3	2.18	0.42
1:C:664:LEU:HD21	1:D:660:LYS:HD3	2.00	0.42
1:B:708:ILE:CG2	1:B:872:PHE:CE1	3.01	0.42
1:D:600:GLN:NE2	5:D:1182:HOH:O	2.53	0.42
1:A:569:ILE:CD1	1:A:594:VAL:HG21	2.50	0.41
1:A:638[A]:SER:HB3	1:A:661:SER:HB2	2.01	0.41
1:B:781[A]:ARG:HD2	1:B:785:PHE:CD2	2.56	0.41
1:D:529:THR:HG21	1:D:582:TRP:HA	2.02	0.41
1:C:529:THR:HG21	1:C:582:TRP:HA	2.03	0.41
1:C:901:GLU:OE2	5:C:1150:HOH:O	2.22	0.41
1:A:539:LYS:HD3	1:C:536:TRP:CE2	2.56	0.41
1:D:803[B]:GLU:HG3	1:D:804:GLU:N	2.37	0.40
1:A:582:TRP:CD1	1:A:884:LEU:HD21	2.57	0.40
1:C:806:PHE:CZ	4:C:1009[A]:TAP:H2D	2.57	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	518/517 (100%)	503 (97%)	15 (3%)	0	100	100
1	B	519/517 (100%)	501 (96%)	18 (4%)	0	100	100
1	C	516/517 (100%)	496 (96%)	20 (4%)	0	100	100
1	D	517/517 (100%)	500 (97%)	17 (3%)	0	100	100
All	All	2070/2068 (100%)	2000 (97%)	70 (3%)	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	430/426 (101%)	418 (97%)	12 (3%)	51	66
1	B	431/426 (101%)	422 (98%)	9 (2%)	61	76
1	C	428/426 (100%)	417 (97%)	11 (3%)	54	69
1	D	429/426 (101%)	421 (98%)	8 (2%)	65	79
All	All	1718/1704 (101%)	1678 (98%)	40 (2%)	58	73

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	507	GLU
1	A	547	ILE
1	A	569	ILE
1	A	673	GLU
1	A	703	LYS
1	A	735	LYS
1	A	745	ASN
1	A	756[A]	ARG
1	A	756[B]	ARG
1	A	781[A]	ARG
1	A	781[B]	ARG
1	A	806	PHE
1	B	413	VAL
1	B	517	LEU
1	B	547	ILE
1	B	597	LYS
1	B	673	GLU
1	B	703	LYS
1	B	806	PHE
1	B	819	ASP
1	B	845	LYS
1	C	517	LEU
1	C	547	ILE
1	C	569	ILE
1	C	667	VAL
1	C	673	GLU
1	C	703	LYS
1	C	725[A]	GLN
1	C	725[B]	GLN
1	C	777	ASN
1	C	806	PHE
1	C	819	ASP
1	D	418	LEU
1	D	547	ILE
1	D	569	ILE
1	D	673	GLU
1	D	703	LYS
1	D	742	ARG
1	D	745	ASN
1	D	806	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	457	GLN
1	A	706	ASN
1	A	745	ASN
1	B	457	GLN
1	B	600	GLN
1	B	706	ASN
1	C	600	GLN
1	C	706	ASN
1	C	797	HIS
1	D	407	ASN
1	D	498	GLN
1	D	548	ASN
1	D	600	GLN
1	D	706	ASN
1	D	745	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](#)

40 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	SO4	A	3001	-	4,4,4	0.23	0	6,6,6	0.20	0
2	SO4	A	3002	-	4,4,4	0.29	0	6,6,6	0.68	0
2	SO4	A	3003	-	4,4,4	0.20	0	6,6,6	0.13	0
2	SO4	A	3004	-	4,4,4	0.18	0	6,6,6	0.21	0
2	SO4	A	3005	-	4,4,4	0.19	0	6,6,6	0.36	0
2	SO4	A	3006	-	4,4,4	0.18	0	6,6,6	0.18	0
2	SO4	A	3007	-	4,4,4	0.24	0	6,6,6	0.60	0
3	GOL	A	3008	-	5,5,5	0.32	0	5,5,5	0.33	0
4	TAP	A	3009[A]	-	35,42,52	0.97	2 (5%)	46,65,80	1.57	6 (13%)
4	TAP	A	3009[B]	-	42,52,52	1.57	5 (11%)	53,80,80	1.64	8 (15%)
2	SO4	B	1001	-	4,4,4	0.37	0	6,6,6	0.26	0
2	SO4	B	1002	-	4,4,4	0.16	0	6,6,6	0.13	0
2	SO4	B	1003	-	4,4,4	0.08	0	6,6,6	0.22	0
2	SO4	B	1004	-	4,4,4	0.18	0	6,6,6	0.13	0
2	SO4	B	1005	-	4,4,4	0.21	0	6,6,6	0.19	0
2	SO4	B	1006	-	4,4,4	0.19	0	6,6,6	0.24	0
2	SO4	B	1007	-	4,4,4	0.21	0	6,6,6	0.22	0
2	SO4	B	1008	-	4,4,4	0.15	0	6,6,6	0.39	0
3	GOL	B	1009	-	5,5,5	0.31	0	5,5,5	0.28	0
4	TAP	B	1010[A]	-	35,42,52	0.88	2 (5%)	46,65,80	1.61	5 (10%)
4	TAP	B	1010[B]	-	42,52,52	1.49	3 (7%)	53,80,80	1.59	7 (13%)
2	SO4	C	1001	-	4,4,4	0.23	0	6,6,6	0.17	0
2	SO4	C	1002	-	4,4,4	0.24	0	6,6,6	0.56	0
2	SO4	C	1003	-	4,4,4	0.19	0	6,6,6	0.17	0
2	SO4	C	1004	-	4,4,4	0.19	0	6,6,6	0.27	0
2	SO4	C	1005	-	4,4,4	0.21	0	6,6,6	0.22	0
2	SO4	C	1006	-	4,4,4	0.24	0	6,6,6	0.81	0
2	SO4	C	1007	-	4,4,4	0.23	0	6,6,6	0.66	0
3	GOL	C	1008	-	5,5,5	0.36	0	5,5,5	0.43	0
4	TAP	C	1009[A]	-	35,42,52	0.93	2 (5%)	46,65,80	1.54	5 (10%)
4	TAP	C	1009[B]	-	42,52,52	1.54	3 (7%)	53,80,80	1.59	8 (15%)
2	SO4	D	1001	-	4,4,4	0.17	0	6,6,6	0.59	0
2	SO4	D	1002	-	4,4,4	0.18	0	6,6,6	0.14	0
2	SO4	D	1003	-	4,4,4	0.16	0	6,6,6	0.25	0
2	SO4	D	1004	-	4,4,4	0.15	0	6,6,6	0.20	0
2	SO4	D	1005	-	4,4,4	0.23	0	6,6,6	0.26	0
2	SO4	D	1006	-	4,4,4	0.19	0	6,6,6	0.56	0
3	GOL	D	1007	-	5,5,5	0.30	0	5,5,5	0.30	0
4	TAP	D	1008[A]	-	35,42,52	0.90	2 (5%)	46,65,80	1.79	7 (15%)
4	TAP	D	1008[B]	-	42,52,52	1.54	4 (9%)	53,80,80	1.74	10 (18%)

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	SO4	A	3001	-	-	0/0/0/0	0/0/0/0
2	SO4	A	3002	-	-	0/0/0/0	0/0/0/0
2	SO4	A	3003	-	-	0/0/0/0	0/0/0/0
2	SO4	A	3004	-	-	0/0/0/0	0/0/0/0
2	SO4	A	3005	-	-	0/0/0/0	0/0/0/0
2	SO4	A	3006	-	-	0/0/0/0	0/0/0/0
2	SO4	A	3007	-	-	0/0/0/0	0/0/0/0
3	GOL	A	3008	-	-	0/4/4/4	0/0/0/0
4	TAP	A	3009[A]	-	-	0/23/56/67	0/4/4/5
4	TAP	A	3009[B]	-	-	0/27/67/67	0/5/5/5
2	SO4	B	1001	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1002	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1003	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1004	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1005	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1006	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1007	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1008	-	-	0/0/0/0	0/0/0/0
3	GOL	B	1009	-	-	0/4/4/4	0/0/0/0
4	TAP	B	1010[A]	-	-	0/23/56/67	0/4/4/5
4	TAP	B	1010[B]	-	-	0/27/67/67	0/5/5/5
2	SO4	C	1001	-	-	0/0/0/0	0/0/0/0
2	SO4	C	1002	-	-	0/0/0/0	0/0/0/0
2	SO4	C	1003	-	-	0/0/0/0	0/0/0/0
2	SO4	C	1004	-	-	0/0/0/0	0/0/0/0
2	SO4	C	1005	-	-	0/0/0/0	0/0/0/0
2	SO4	C	1006	-	-	0/0/0/0	0/0/0/0
2	SO4	C	1007	-	-	0/0/0/0	0/0/0/0
3	GOL	C	1008	-	-	0/4/4/4	0/0/0/0
4	TAP	C	1009[A]	-	-	0/23/56/67	0/4/4/5
4	TAP	C	1009[B]	-	-	0/27/67/67	0/5/5/5
2	SO4	D	1001	-	-	0/0/0/0	0/0/0/0
2	SO4	D	1002	-	-	0/0/0/0	0/0/0/0
2	SO4	D	1003	-	-	0/0/0/0	0/0/0/0
2	SO4	D	1004	-	-	0/0/0/0	0/0/0/0
2	SO4	D	1005	-	-	0/0/0/0	0/0/0/0
2	SO4	D	1006	-	-	0/0/0/0	0/0/0/0
3	GOL	D	1007	-	-	0/4/4/4	0/0/0/0
4	TAP	D	1008[A]	-	-	0/23/56/67	0/4/4/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	TAP	D	1008[B]	-	-	0/27/67/67	0/5/5/5

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	1009[B]	TAP	C7N-S7N	-7.71	1.56	1.67
4	D	1008[B]	TAP	C7N-S7N	-7.71	1.56	1.67
4	A	3009[B]	TAP	C7N-S7N	-7.51	1.56	1.67
4	B	1010[B]	TAP	C7N-S7N	-7.44	1.56	1.67
4	A	3009[B]	TAP	C7N-N7N	2.06	1.33	1.32
4	D	1008[B]	TAP	C7N-N7N	2.15	1.34	1.32
4	C	1009[B]	TAP	O4B-C1B	2.18	1.44	1.41
4	C	1009[A]	TAP	O4B-C1B	2.18	1.44	1.41
4	A	3009[B]	TAP	O4D-C1D	2.19	1.44	1.41
4	A	3009[A]	TAP	O4B-C1B	2.29	1.44	1.41
4	A	3009[B]	TAP	O4B-C1B	2.29	1.44	1.41
4	B	1010[A]	TAP	O4B-C1B	2.32	1.44	1.41
4	B	1010[B]	TAP	O4B-C1B	2.32	1.44	1.41
4	D	1008[A]	TAP	O4B-C1B	2.37	1.44	1.41
4	D	1008[B]	TAP	O4B-C1B	2.37	1.44	1.41
4	B	1010[A]	TAP	C5A-C4A	3.28	1.47	1.40
4	B	1010[B]	TAP	C5A-C4A	3.28	1.47	1.40
4	D	1008[A]	TAP	C5A-C4A	3.41	1.48	1.40
4	D	1008[B]	TAP	C5A-C4A	3.41	1.48	1.40
4	C	1009[B]	TAP	C5A-C4A	3.59	1.48	1.40
4	C	1009[A]	TAP	C5A-C4A	3.59	1.48	1.40
4	A	3009[A]	TAP	C5A-C4A	3.66	1.48	1.40
4	A	3009[B]	TAP	C5A-C4A	3.66	1.48	1.40

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1008[A]	TAP	N3A-C2A-N1A	-8.15	122.65	128.89
4	D	1008[B]	TAP	N3A-C2A-N1A	-8.15	122.65	128.89
4	B	1010[A]	TAP	N3A-C2A-N1A	-7.19	123.39	128.89
4	B	1010[B]	TAP	N3A-C2A-N1A	-7.19	123.39	128.89
4	A	3009[A]	TAP	N3A-C2A-N1A	-6.63	123.82	128.89
4	A	3009[B]	TAP	N3A-C2A-N1A	-6.63	123.82	128.89
4	C	1009[B]	TAP	N3A-C2A-N1A	-6.36	124.03	128.89
4	C	1009[A]	TAP	N3A-C2A-N1A	-6.36	124.03	128.89
4	D	1008[A]	TAP	PN-O3-PA	-3.77	122.15	132.73
4	A	3009[A]	TAP	C4A-C5A-N7A	-3.37	106.38	109.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	3009[B]	TAP	C4A-C5A-N7A	-3.37	106.38	109.48
4	B	1010[A]	TAP	PN-O3-PA	-3.34	123.34	132.73
4	D	1008[A]	TAP	C4A-C5A-N7A	-3.34	106.40	109.48
4	D	1008[B]	TAP	C4A-C5A-N7A	-3.34	106.40	109.48
4	C	1009[A]	TAP	PN-O3-PA	-3.31	123.45	132.73
4	A	3009[B]	TAP	PN-O3-PA	-3.26	123.59	132.73
4	C	1009[B]	TAP	PN-O3-PA	-3.25	123.59	132.73
4	D	1008[A]	TAP	C1B-N9A-C4A	-3.16	122.18	126.94
4	D	1008[B]	TAP	C1B-N9A-C4A	-3.16	122.18	126.94
4	B	1010[A]	TAP	C1B-N9A-C4A	-3.08	122.30	126.94
4	B	1010[B]	TAP	C1B-N9A-C4A	-3.08	122.30	126.94
4	A	3009[B]	TAP	S7N-C7N-N7N	-2.92	118.00	122.23
4	D	1008[B]	TAP	PN-O3-PA	-2.90	124.58	132.73
4	C	1009[B]	TAP	C4N-C3N-C7N	-2.84	118.49	120.95
4	A	3009[A]	TAP	PN-O3-PA	-2.81	124.84	132.73
4	C	1009[B]	TAP	S7N-C7N-N7N	-2.69	118.34	122.23
4	B	1010[B]	TAP	PN-O3-PA	-2.56	125.54	132.73
4	B	1010[A]	TAP	C4A-C5A-N7A	-2.50	107.18	109.48
4	B	1010[B]	TAP	C4A-C5A-N7A	-2.50	107.18	109.48
4	B	1010[B]	TAP	S7N-C7N-N7N	-2.48	118.63	122.23
4	D	1008[B]	TAP	S7N-C7N-N7N	-2.33	118.86	122.23
4	A	3009[A]	TAP	O4B-C1B-C2B	-2.33	102.39	106.60
4	A	3009[B]	TAP	O4B-C1B-C2B	-2.33	102.39	106.60
4	C	1009[B]	TAP	C4A-C5A-N7A	-2.12	107.53	109.48
4	C	1009[A]	TAP	C4A-C5A-N7A	-2.12	107.53	109.48
4	C	1009[B]	TAP	C1B-N9A-C4A	-2.07	123.81	126.94
4	C	1009[A]	TAP	C1B-N9A-C4A	-2.07	123.81	126.94
4	D	1008[B]	TAP	O4D-C1D-N1N	2.02	110.36	108.13
4	A	3009[A]	TAP	C1D-C2D-C3D	2.08	104.99	101.64
4	D	1008[A]	TAP	O4B-C1B-N9A	2.42	113.17	108.10
4	D	1008[B]	TAP	O4B-C1B-N9A	2.42	113.17	108.10
4	D	1008[A]	TAP	O3X-P2B-O2X	2.42	116.60	107.38
4	D	1008[B]	TAP	O3X-P2B-O2X	2.42	116.60	107.38
4	C	1009[B]	TAP	O4B-C1B-N9A	2.59	113.52	108.10
4	C	1009[A]	TAP	O4B-C1B-N9A	2.59	113.52	108.10
4	B	1010[A]	TAP	O4B-C1B-N9A	2.65	113.65	108.10
4	B	1010[B]	TAP	O4B-C1B-N9A	2.65	113.65	108.10
4	A	3009[A]	TAP	O4B-C1B-N9A	2.68	113.70	108.10
4	A	3009[B]	TAP	O4B-C1B-N9A	2.68	113.70	108.10
4	D	1008[A]	TAP	C2A-N1A-C6A	2.70	123.59	118.77
4	D	1008[B]	TAP	C2A-N1A-C6A	2.70	123.59	118.77
4	B	1010[B]	TAP	C3N-C7N-N7N	2.72	120.76	116.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1008[B]	TAP	C3N-C7N-N7N	2.87	120.98	116.78
4	C	1009[B]	TAP	C3N-C7N-N7N	3.29	121.60	116.78
4	A	3009[B]	TAP	C3N-C7N-N7N	3.31	121.63	116.78
4	A	3009[B]	TAP	O4D-C1D-N1N	3.91	112.42	108.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

10 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	3007	SO4	1	0
4	A	3009[A]	TAP	2	0
4	B	1010[A]	TAP	2	0
4	B	1010[B]	TAP	2	0
2	C	1001	SO4	1	0
3	C	1008	GOL	1	0
4	C	1009[A]	TAP	2	0
4	C	1009[B]	TAP	4	0
4	D	1008[A]	TAP	1	0
4	D	1008[B]	TAP	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	498/517 (96%)	-0.21	4 (0%) 87 90	15, 27, 39, 46	20 (4%)
1	B	498/517 (96%)	-0.14	9 (1%) 71 77	16, 29, 39, 57	17 (3%)
1	C	498/517 (96%)	-0.36	1 (0%) 95 97	15, 24, 33, 48	18 (3%)
1	D	498/517 (96%)	-0.34	6 (1%) 81 85	15, 27, 37, 55	19 (3%)
All	All	1992/2068 (96%)	-0.27	20 (1%) 84 87	15, 27, 38, 57	74 (3%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	405	VAL	4.5
1	A	405	VAL	4.4
1	B	405	VAL	4.0
1	D	415	LYS	3.4
1	B	415	LYS	3.1
1	C	405	VAL	3.0
1	B	743[A]	ASP	2.6
1	B	742	ARG	2.6
1	B	413	VAL	2.4
1	B	436	SER	2.3
1	A	887	ALA	2.3
1	D	742	ARG	2.3
1	D	552	PRO	2.2
1	A	797	HIS	2.2
1	A	435	GLY	2.1
1	D	551	ARG	2.1
1	D	490[A]	TYR	2.0
1	B	452[A]	GLN	2.0
1	B	893	LEU	2.0
1	B	435	GLY	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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
2	SO4	B	1006	5/5	0.85	0.23	34.89	45,47,48,48	5
2	SO4	A	3005	5/5	0.81	0.30	13.88	45,47,47,47	5
2	SO4	D	1004	5/5	0.88	0.23	7.25	41,43,43,44	5
2	SO4	A	3006	5/5	0.86	0.26	7.18	53,54,54,54	5
3	GOL	D	1007	6/6	0.84	0.20	3.80	68,69,69,69	0
3	GOL	C	1008	6/6	0.87	0.23	3.04	61,62,63,64	0
3	GOL	A	3008	6/6	0.89	0.25	2.90	60,61,62,62	0
4	TAP	C	1009[B]	48/48	0.93	0.17	1.79	33,40,53,57	25
4	TAP	C	1009[A]	39/48	0.93	0.17	1.79	33,37,40,41	16
4	TAP	B	1010[A]	39/48	0.92	0.18	1.71	41,45,51,51	16
4	TAP	B	1010[B]	48/48	0.92	0.18	1.71	41,47,63,65	25
3	GOL	B	1009	6/6	0.92	0.20	1.63	58,60,61,61	0
4	TAP	A	3009[A]	39/48	0.92	0.18	1.58	35,41,51,51	16
2	SO4	A	3007	5/5	0.93	0.22	1.53	36,43,43,44	5
4	TAP	A	3009[B]	48/48	0.92	0.18	1.48	35,45,61,64	25
4	TAP	D	1008[B]	48/48	0.92	0.17	1.45	32,40,53,58	25
4	TAP	D	1008[A]	39/48	0.92	0.17	1.45	32,39,47,48	16
2	SO4	C	1006	5/5	0.97	0.17	1.41	39,41,42,42	5
2	SO4	C	1005	5/5	0.94	0.17	1.19	43,44,44,44	5
2	SO4	D	1006	5/5	0.96	0.18	1.15	34,39,40,41	5
2	SO4	B	1008	5/5	0.95	0.24	1.06	35,40,42,42	5
2	SO4	D	1005	5/5	0.88	0.23	0.65	45,45,45,46	5
2	SO4	B	1001	5/5	0.98	0.14	0.44	26,26,27,27	5
2	SO4	C	1007	5/5	0.95	0.11	0.36	30,34,36,36	5
2	SO4	B	1007	5/5	0.92	0.18	0.13	44,44,46,46	5

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	SO4	C	1001	5/5	0.95	0.13	0.04	38,39,40,41	5
2	SO4	B	1003	5/5	0.94	0.17	-	41,41,44,44	5
2	SO4	D	1003	5/5	0.73	0.25	-	48,48,50,50	5
2	SO4	A	3003	5/5	0.80	0.28	-	85,85,86,86	5
2	SO4	A	3002	5/5	0.94	0.13	-	34,36,38,39	5
2	SO4	B	1002	5/5	0.97	0.09	-	34,35,37,37	5
2	SO4	A	3004	5/5	0.90	0.14	-	52,53,53,54	5
2	SO4	D	1002	5/5	0.86	0.27	-	86,86,87,87	5
2	SO4	C	1002	5/5	0.95	0.13	-	38,38,40,40	5
2	SO4	A	3001	5/5	0.98	0.14	-	37,38,39,39	5
2	SO4	B	1005	5/5	0.87	0.17	-	49,50,51,51	5
2	SO4	D	1001	5/5	0.95	0.13	-	33,34,36,36	5
2	SO4	C	1004	5/5	0.87	0.23	-	52,52,52,52	5
2	SO4	C	1003	5/5	0.92	0.20	-	76,77,77,77	5
2	SO4	B	1004	5/5	0.93	0.22	-	96,96,97,97	0

6.5 Other polymers [i](#)

There are no such residues in this entry.