



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:48 PM GMT

PDB ID : 4JIM
Title : Native Crystal Strucutre of N10-Formyltetrahydrofolate Synthetase
Authors : Celeste, L.R.; Lovelace, L.L.; Lebioda, L.
Deposited on : 2013-03-06
Resolution : 2.10 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

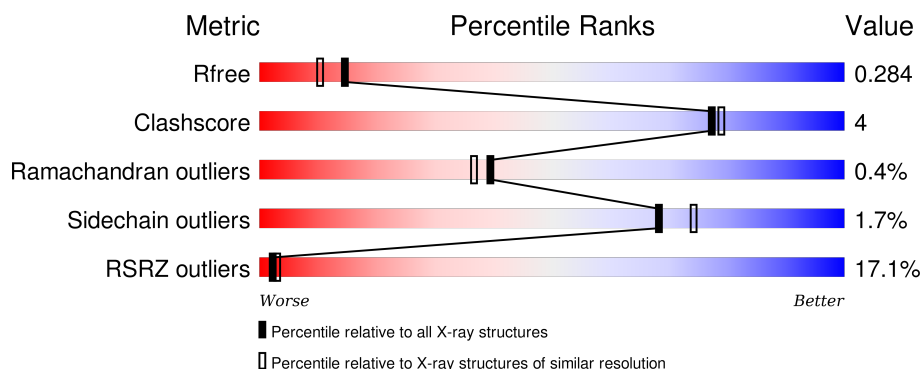
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	559	
1	B	559	

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
3	SO4	A	705	-	-	-	X
3	SO4	A	707	-	-	-	X



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

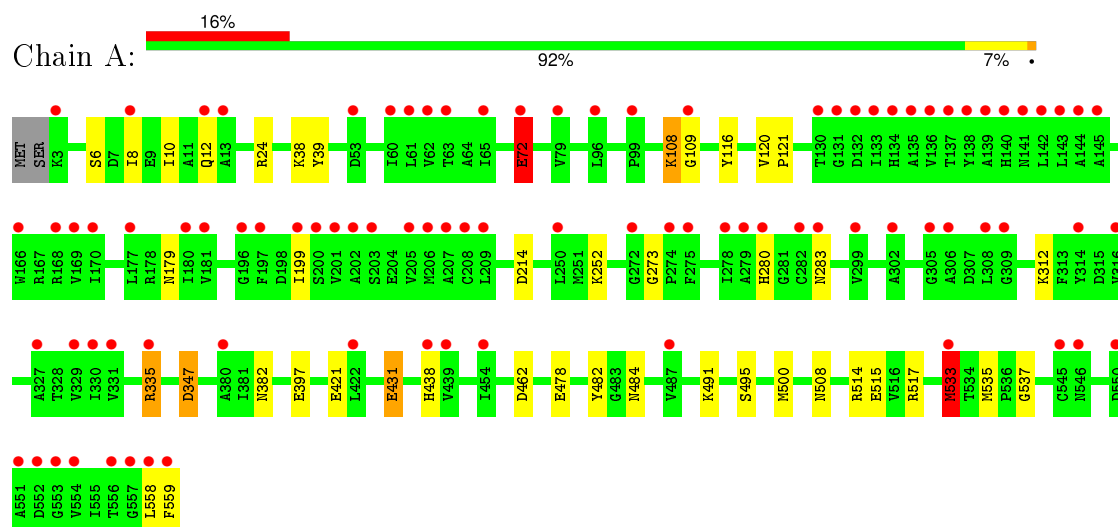
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	153	Total	O	0	0
			153	153		
4	A	399	Total	O	0	0
			399	399		

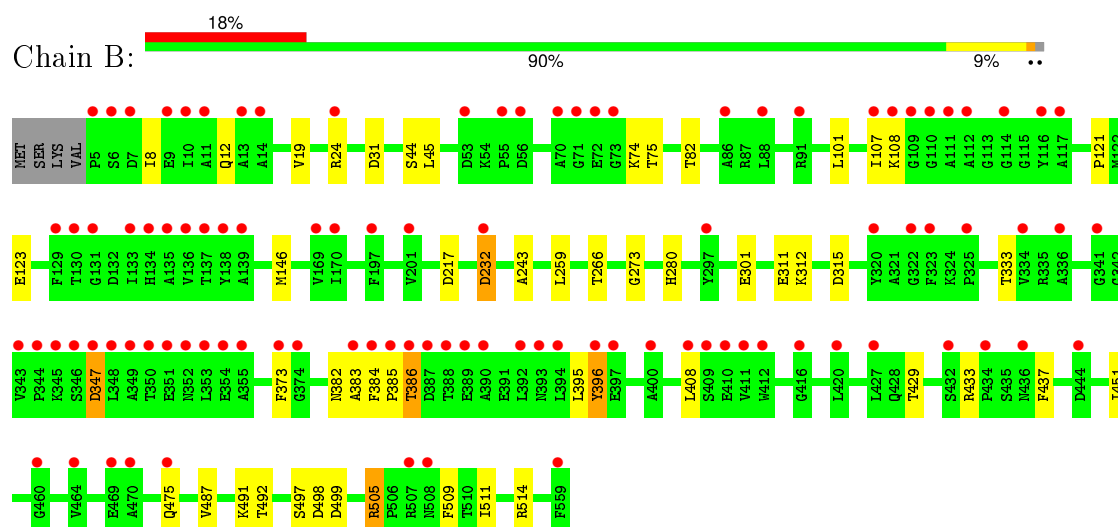
3 Residue-property plots

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: Formate--tetrahydrofolate ligase



- Molecule 1: Formate--tetrahydrofolate ligase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	161.19Å 161.19Å 256.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.92 – 2.10 40.89 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.92-2.10) 99.9 (40.89-2.10)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.59 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.234 , 0.274 0.243 , 0.284	Depositor DCC
R_{free} test set	3753 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	29.1	Xtriage
Anisotropy	0.152	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 52.3	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 74517 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8890	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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: SO4, TOE

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	0.95	12/4255 (0.3%)	0.92	10/5771 (0.2%)
1	B	0.80	1/4168 (0.0%)	0.85	2/5669 (0.0%)
All	All	0.88	13/8423 (0.2%)	0.89	12/11440 (0.1%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	438[A]	HIS	N-CA	-9.07	1.28	1.46
1	A	438[B]	HIS	N-CA	-9.07	1.28	1.46
1	B	509	PHE	C-O	7.76	1.38	1.23
1	A	478	GLU	CD-OE2	6.51	1.32	1.25
1	A	431	GLU	CD-OE1	6.37	1.32	1.25
1	A	421	GLU	CD-OE1	5.71	1.31	1.25
1	A	431	GLU	CD-OE2	5.51	1.31	1.25
1	A	335[A]	ARG	N-CA	-5.33	1.35	1.46
1	A	335[B]	ARG	N-CA	-5.33	1.35	1.46
1	A	72	GLU	CD-OE2	5.31	1.31	1.25
1	A	72	GLU	CD-OE1	5.18	1.31	1.25
1	A	537	GLY	C-O	-5.05	1.15	1.23
1	A	397	GLU	CG-CD	5.04	1.59	1.51

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	347	ASP	CB-CG-OD1	6.67	124.31	118.30
1	A	438[A]	HIS	N-CA-CB	6.43	122.18	110.60
1	A	438[B]	HIS	N-CA-CB	6.43	122.18	110.60
1	B	217	ASP	CB-CG-OD2	-6.43	112.51	118.30
1	A	214	ASP	CB-CG-OD1	6.09	123.78	118.30
1	A	514	ARG	NE-CZ-NH2	-5.41	117.59	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	514	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	A	335[A]	ARG	NE-CZ-NH1	5.28	122.94	120.30
1	A	335[B]	ARG	NE-CZ-NH1	5.28	122.94	120.30
1	A	533[A]	MET	CB-CG-SD	5.08	127.63	112.40
1	A	533[B]	MET	CB-CG-SD	5.08	127.63	112.40
1	B	232	ASP	CB-CG-OD1	-5.01	113.80	118.30

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	4174	0	4231	27	1
1	B	4091	0	4067	35	0
2	A	22	0	32	1	0
2	B	11	0	16	1	0
3	A	25	0	0	0	0
3	B	15	0	0	2	0
4	A	399	0	0	9	0
4	B	153	0	0	7	0
All	All	8890	0	8346	62	1

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 (62) 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:505[A]:ARG:HG2	1:B:505[A]:ARG:HH21	1.17	1.09
1:B:505[A]:ARG:CG	1:B:505[A]:ARG:HH21	1.73	0.99
1:A:108:LYS:CB	1:A:109:GLY:HA2	2.03	0.89
1:A:533[B]:MET:SD	4:A:1174:HOH:O	2.32	0.86
1:A:72:GLU:HG3	4:A:875:HOH:O	1.79	0.82
1:B:429:THR:HG23	1:B:433:ARG:HD3	1.66	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:491:LYS:NZ	1:B:497:SER:O	2.20	0.74
1:A:500:MET:CE	4:A:1180:HOH:O	2.35	0.73
1:A:8:ILE:O	1:A:12:GLN:HG2	1.89	0.73
1:B:505[A]:ARG:HG2	1:B:505[A]:ARG:NH2	1.96	0.71
1:A:491:LYS:NZ	1:A:495:SER:O	2.30	0.65
1:A:108:LYS:CB	1:A:109:GLY:CA	2.76	0.65
1:B:107:ILE:HD12	1:B:108:LYS:CB	2.28	0.63
1:A:500:MET:HE3	4:A:1180:HOH:O	1.96	0.61
1:A:515:GLU:OE1	1:A:517:ARG:CZ	2.48	0.61
1:B:383:ALA:HB2	1:B:408:LEU:HD11	1.84	0.60
1:B:505[A]:ARG:NH1	4:B:786:HOH:O	2.35	0.60
1:A:515:GLU:OE1	1:A:517:ARG:NH1	2.35	0.59
1:A:484:ASN:OD1	4:A:1084:HOH:O	2.17	0.57
1:A:252:LYS:HE2	1:B:123:GLU:OE1	2.04	0.57
1:B:347:ASP:N	1:B:347:ASP:OD1	2.38	0.57
1:B:8:ILE:O	1:B:12:GLN:HG2	2.07	0.55
1:B:333:THR:HG22	1:B:382:ASN:HB3	1.89	0.55
1:B:384:PHE:O	1:B:386:THR:N	2.41	0.53
1:A:6:SER:O	1:A:10:ILE:HG12	2.08	0.53
1:A:179:ASN:ND2	4:A:1090:HOH:O	2.43	0.52
1:A:558:LEU:HD23	1:A:559:PHE:CE2	2.45	0.52
1:B:44:SER:HB2	3:B:604:SO4:O3	2.11	0.51
1:B:45:LEU:HA	4:B:768:HOH:O	2.09	0.51
1:B:505[A]:ARG:CG	1:B:505[A]:ARG:NH2	2.45	0.51
1:B:514:ARG:NH2	4:B:734:HOH:O	2.43	0.51
1:B:498:ASP:OD2	1:B:511:ILE:HA	2.11	0.50
1:B:280:HIS:HA	1:B:312:LYS:HD3	1.94	0.50
1:B:499:ASP:C	1:B:499:ASP:OD1	2.50	0.50
1:B:101:LEU:HD23	1:B:101:LEU:C	2.33	0.49
1:A:10:ILE:HD12	1:A:116:TYR:CE2	2.49	0.48
1:B:74:LYS:NZ	3:B:603:SO4:O4	2.47	0.47
1:A:10:ILE:HD12	1:A:116:TYR:CZ	2.49	0.47
1:A:120:VAL:HB	1:A:121:PRO:HA	1.98	0.46
1:A:38:LYS:HG3	1:A:39:TYR:CZ	2.52	0.45
1:A:462:ASP:C	4:A:901:HOH:O	2.55	0.45
1:B:82:THR:HG22	1:B:266:THR:HG21	1.98	0.45
1:B:121:PRO:HA	4:B:805:HOH:O	2.15	0.45
1:B:74:LYS:HB3	4:B:803:HOH:O	2.16	0.45
2:B:601:TOE:H12	4:B:836:HOH:O	2.16	0.44
1:B:373:PHE:O	1:B:437:PHE:HA	2.17	0.44
1:B:75:THR:HA	1:B:301:GLU:OE1	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:482:TYR:CE1	2:A:702:TOE:H7	2.53	0.44
1:B:505[A]:ARG:HH21	1:B:505[A]:ARG:HG3	1.73	0.44
1:B:19:VAL:HG13	1:B:259:LEU:HD23	2.00	0.44
1:A:558:LEU:HG	1:A:559:PHE:CZ	2.54	0.43
1:A:347:ASP:HB3	4:A:913:HOH:O	2.18	0.43
1:B:395:LEU:O	1:B:396:TYR:C	2.57	0.42
1:B:492:THR:HG22	4:B:779:HOH:O	2.18	0.42
4:A:1186:HOH:O	1:B:31:ASP:HA	2.19	0.42
1:A:535:MET:O	1:A:535:MET:HG3	2.19	0.41
1:B:146:MET:HE3	1:B:243:ALA:CB	2.49	0.41
1:A:280:HIS:HA	1:A:312:LYS:HD3	2.01	0.41
1:B:451:ILE:HG12	1:B:487:VAL:HG11	2.03	0.40
1:A:335[B]:ARG:HD2	1:A:335[B]:ARG:HH11	1.77	0.40
1:A:462:ASP:OD2	1:A:508:ASN:HA	2.22	0.40
1:B:311:GLU:OE2	1:B:315:ASP:OD2	2.38	0.40

All (1) 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:A:462:ASP:O	1:A:462:ASP:O[17_554]	2.02	0.18

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	558/559 (100%)	537 (96%)	19 (3%)	2 (0%)	39	37
1	B	555/559 (99%)	521 (94%)	32 (6%)	2 (0%)	39	37
All	All	1113/1118 (100%)	1058 (95%)	51 (5%)	4 (0%)	39	37

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	385	PRO
1	A	108	LYS
1	B	273	GLY
1	A	273	GLY

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	432/441 (98%)	424 (98%)	8 (2%)	65	70
1	B	414/441 (94%)	405 (98%)	9 (2%)	60	64
All	All	846/882 (96%)	829 (98%)	17 (2%)	68	68

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	ARG
1	A	72	GLU
1	A	199	ILE
1	A	283	ASN
1	A	382	ASN
1	A	431	GLU
1	A	533[A]	MET
1	A	533[B]	MET
1	B	24[A]	ARG
1	B	24[B]	ARG
1	B	232	ASP
1	B	347	ASP
1	B	386	THR
1	B	396	TYR
1	B	475	GLN
1	B	505[A]	ARG
1	B	505[B]	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	A	179	ASN
1	B	382	ASN
1	B	442	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

11 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	TOE	A	701	-	10,10,10	1.25	0	9,9,9	1.69	3 (33%)
2	TOE	A	702	-	10,10,10	0.77	0	9,9,9	0.82	0
3	SO4	A	703	-	4,4,4	1.09	0	6,6,6	0.39	0
3	SO4	A	704	-	4,4,4	1.37	0	6,6,6	0.98	1 (16%)
3	SO4	A	705	-	4,4,4	1.42	0	6,6,6	0.50	0
3	SO4	A	706	-	4,4,4	0.74	0	6,6,6	0.21	0
3	SO4	A	707	-	4,4,4	0.77	0	6,6,6	0.36	0
2	TOE	B	601	-	10,10,10	0.68	0	9,9,9	0.54	0
3	SO4	B	602	-	4,4,4	0.97	0	6,6,6	1.05	0
3	SO4	B	603	-	4,4,4	0.89	0	6,6,6	1.69	1 (16%)
3	SO4	B	604	-	4,4,4	0.61	0	6,6,6	0.22	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	TOE	A	701	-	-	0/8/8/8	0/0/0/0
2	TOE	A	702	-	-	0/8/8/8	0/0/0/0
3	SO4	A	703	-	-	0/0/0/0	0/0/0/0
3	SO4	A	704	-	-	0/0/0/0	0/0/0/0
3	SO4	A	705	-	-	0/0/0/0	0/0/0/0
3	SO4	A	706	-	-	0/0/0/0	0/0/0/0
3	SO4	A	707	-	-	0/0/0/0	0/0/0/0
2	TOE	B	601	-	-	0/8/8/8	0/0/0/0
3	SO4	B	602	-	-	0/0/0/0	0/0/0/0
3	SO4	B	603	-	-	0/0/0/0	0/0/0/0
3	SO4	B	604	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	704	SO4	O4-S-O3	-2.02	100.75	108.98
2	A	701	TOE	OC'-CD'-CE'	2.06	119.54	110.36
2	A	701	TOE	OF'-CG'-CH'	2.85	123.04	110.36
2	A	701	TOE	CE'-OF'-CG'	3.23	127.21	113.31
3	B	603	SO4	O2-S-O1	3.84	121.66	109.50

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	702	TOE	1	0
2	B	601	TOE	1	0
3	B	603	SO4	1	0
3	B	604	SO4	1	0

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	557/559 (99%)	0.88	90 (16%) 3 3	41, 43, 47, 64	3 (0%)
1	B	555/559 (99%)	1.01	100 (18%) 2 2	39, 54, 107, 143	1 (0%)
All	All	1112/1118 (99%)	0.95	190 (17%) 2 3	39, 44, 97, 143	4 (0%)

All (190) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	353	LEU	7.0
1	B	384	PHE	6.7
1	A	438[A]	HIS	6.2
1	B	396	TYR	6.2
1	A	533[A]	MET	5.7
1	A	559	PHE	5.5
1	B	432	SER	5.5
1	A	169	VAL	5.2
1	B	559	PHE	5.2
1	B	355	ALA	5.1
1	A	136	VAL	5.0
1	B	343	VAL	5.0
1	A	335[A]	ARG	5.0
1	B	110	GLY	4.9
1	B	5	PRO	4.9
1	B	6	SER	4.9
1	B	389	GLU	4.8
1	A	133	ILE	4.8
1	B	409	SER	4.7
1	B	385	PRO	4.7
1	B	387	ASP	4.6
1	B	388	THR	4.6
1	B	350	THR	4.6
1	B	348	LEU	4.5

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Mol	Chain	Res	Type	RSRZ
1	B	386	THR	4.4
1	A	137	THR	4.3
1	A	201	VAL	4.3
1	B	116	TYR	4.3
1	B	392	LEU	4.3
1	B	346	SER	4.2
1	A	205	VAL	4.2
1	B	394	LEU	4.1
1	A	170	ILE	4.0
1	B	411	VAL	4.0
1	B	475	GLN	4.0
1	B	349	ALA	3.9
1	B	322	GLY	3.9
1	A	135	ALA	3.8
1	B	73	GLY	3.7
1	B	344	PRO	3.7
1	B	136	VAL	3.6
1	B	170	ILE	3.6
1	B	133	ILE	3.6
1	B	334	VAL	3.5
1	B	169	VAL	3.5
1	B	55	PRO	3.4
1	A	329	VAL	3.4
1	A	551	ALA	3.4
1	B	336	ALA	3.4
1	A	142	LEU	3.3
1	A	274	PRO	3.3
1	A	557	GLY	3.3
1	B	345	LYS	3.3
1	A	197	PHE	3.3
1	B	412	TRP	3.3
1	B	137	THR	3.3
1	A	556	THR	3.2
1	A	250	LEU	3.2
1	B	393	ASN	3.2
1	A	554	VAL	3.2
1	A	139	ALA	3.2
1	B	114	GLY	3.2
1	A	207	ALA	3.1
1	A	206	MET	3.1
1	B	323	PHE	3.1
1	B	351	GLU	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	208	CYS	3.1
1	B	13	ALA	3.1
1	B	427	LEU	3.0
1	A	308	LEU	3.0
1	A	200	SER	3.0
1	B	341	GLY	3.0
1	B	347	ASP	3.0
1	B	10	ILE	2.9
1	B	53	ASP	2.9
1	A	282	CYS	2.9
1	B	297	TYR	2.9
1	B	434	PRO	2.9
1	A	553	GLY	2.9
1	B	420	LEU	2.9
1	A	141	ASN	2.9
1	B	107	ILE	2.9
1	A	302	ALA	2.9
1	B	91	ARG	2.8
1	B	88	LEU	2.8
1	B	14	ALA	2.8
1	A	272	GLY	2.8
1	B	436	ASN	2.8
1	B	112	ALA	2.7
1	B	72	GLU	2.7
1	A	552	ASP	2.7
1	A	199	ILE	2.7
1	A	331	VAL	2.7
1	A	12	GLN	2.7
1	B	130	THR	2.7
1	B	9	GLU	2.7
1	B	469	GLU	2.7
1	B	7	ASP	2.7
1	A	13	ALA	2.7
1	B	134	HIS	2.7
1	A	138	TYR	2.7
1	A	202	ALA	2.6
1	A	279	ALA	2.6
1	A	203	SER	2.6
1	B	320	TYR	2.6
1	B	373	PHE	2.6
1	B	374	GLY	2.6
1	A	140	HIS	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	383	ALA	2.6
1	A	61	LEU	2.6
1	A	96	LEU	2.5
1	A	177	LEU	2.5
1	B	109	GLY	2.5
1	A	62	VAL	2.5
1	A	166	TRP	2.5
1	B	397	GLU	2.5
1	B	410	GLU	2.5
1	B	71	GLY	2.5
1	A	65	ILE	2.5
1	A	144	ALA	2.4
1	B	117	ALA	2.4
1	A	131	GLY	2.4
1	B	111	ALA	2.4
1	A	134	HIS	2.4
1	B	408	LEU	2.4
1	B	464	VAL	2.4
1	A	275	PHE	2.4
1	A	558	LEU	2.4
1	A	299	VAL	2.4
1	B	56	ASP	2.4
1	A	309	GLY	2.4
1	A	280	HIS	2.3
1	B	135	ALA	2.3
1	A	454	ILE	2.3
1	B	108	LYS	2.3
1	B	352	ASN	2.3
1	B	86	ALA	2.3
1	A	8	ILE	2.3
1	A	546	ASN	2.3
1	A	63	THR	2.3
1	B	325	PRO	2.3
1	A	306	ALA	2.3
1	A	487	VAL	2.3
1	B	70	ALA	2.3
1	A	545	CYS	2.3
1	A	305	GLY	2.2
1	B	201	VAL	2.2
1	A	130	THR	2.2
1	B	131	GLY	2.2
1	A	422	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	99	PRO	2.2
1	A	278	ILE	2.2
1	B	139	ALA	2.2
1	A	550	ASP	2.2
1	B	232	ASP	2.2
1	B	197	PHE	2.2
1	A	168	ARG	2.2
1	B	11	ALA	2.2
1	B	470	ALA	2.2
1	A	3	LYS	2.2
1	A	314	TYR	2.2
1	B	138	TYR	2.2
1	A	143	LEU	2.2
1	A	53	ASP	2.2
1	A	439	VAL	2.2
1	A	180	ILE	2.2
1	B	507	ARG	2.1
1	A	209	LEU	2.1
1	A	132	ASP	2.1
1	B	129	PHE	2.1
1	B	390	ALA	2.1
1	B	444	ASP	2.1
1	A	380	ALA	2.1
1	B	354	GLU	2.1
1	B	24[A]	ARG	2.1
1	B	400	ALA	2.1
1	A	109	GLY	2.1
1	A	181	VAL	2.1
1	B	460	GLY	2.1
1	A	60	ILE	2.1
1	A	316	VAL	2.1
1	A	145	ALA	2.1
1	B	508	ASN	2.1
1	B	416	GLY	2.0
1	A	330	ILE	2.0
1	A	72	GLU	2.0
1	A	327	ALA	2.0
1	A	283	ASN	2.0
1	A	79	VAL	2.0
1	A	196	GLY	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	SO4	A	707	5/5	0.92	0.33	9.81	59,77,92,98	0
3	SO4	A	705	5/5	0.97	0.24	4.22	34,38,51,60	0
2	TOE	B	601	11/11	0.93	0.15	1.93	33,38,46,48	0
3	SO4	A	704	5/5	0.87	0.17	-0.16	42,43,50,61	0
3	SO4	B	602	5/5	0.95	0.11	-0.45	37,38,47,49	0
2	TOE	A	701	11/11	0.90	0.14	-0.57	23,34,47,47	0
3	SO4	B	603	5/5	0.94	0.11	-0.94	41,55,63,64	0
3	SO4	A	703	5/5	0.98	0.09	-4.58	21,25,26,40	0
3	SO4	B	604	5/5	0.82	0.31	-	80,86,102,113	0
2	TOE	A	702	11/11	0.92	0.22	-	33,39,59,60	0
3	SO4	A	706	5/5	0.88	0.34	-	67,72,83,101	0

6.5 Other polymers [i](#)

There are no such residues in this entry.