



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 7, 2020 – 02:26 PM BST

PDB ID : 4JIM
Title : Native Crystal Structure 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.14.2
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.14.2

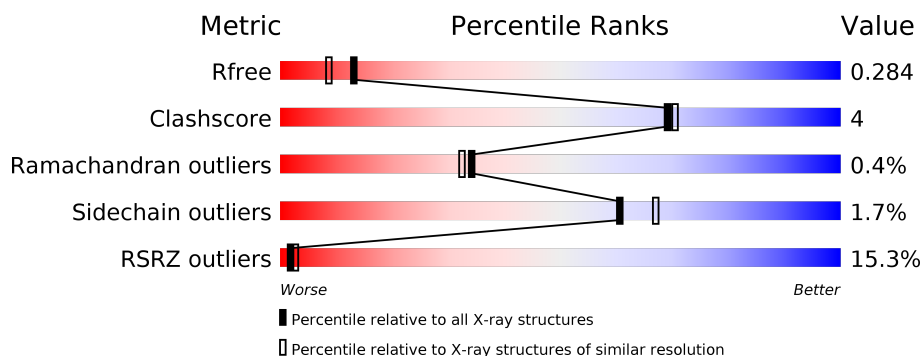
1 Overall quality at a glance

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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (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 respectively. 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	 14% 92% 7% •
1	B	559	 17% 90% 9% ••

2 Entry composition [i](#)

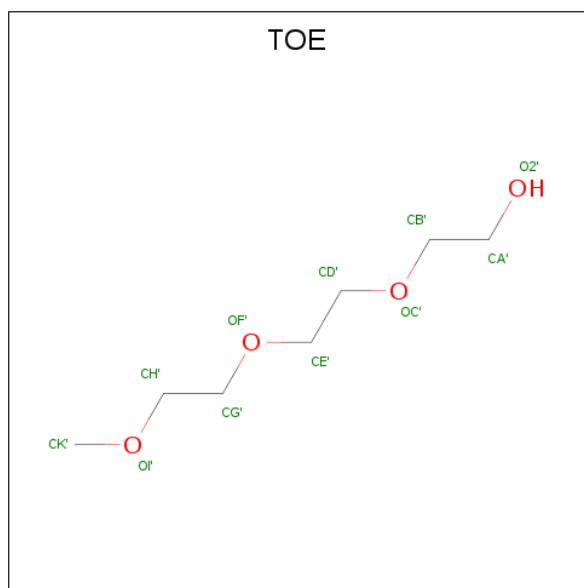
There are 4 unique types of molecules in this entry. The entry contains 8890 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Formate--tetrahydrofolate ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	557	Total	C	N	O	S	20	3	0
			4174	2652	716	784	22			
1	B	555	Total	C	N	O	S	8	2	0
			4091	2595	700	775	21			

- Molecule 2 is 2-[2-(2-METHOXY-ETHOXY)-ETHOXY]-ETHOXYL (three-letter code: TOE) (formula: C₇H₁₆O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			11	7	4		
2	A	1	Total	C	O	0	0
			11	7	4		
2	B	1	Total	C	O	0	0
			11	7	4		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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		

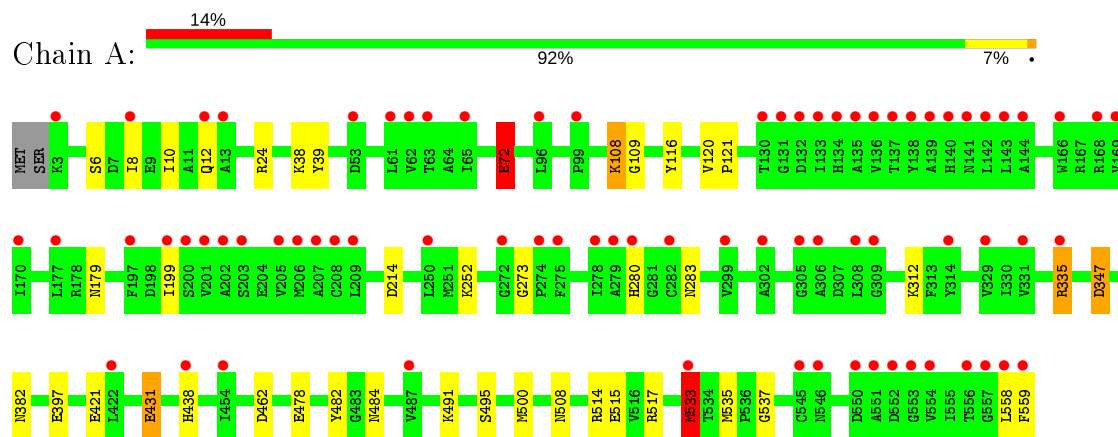
- Molecule 4 is water.

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

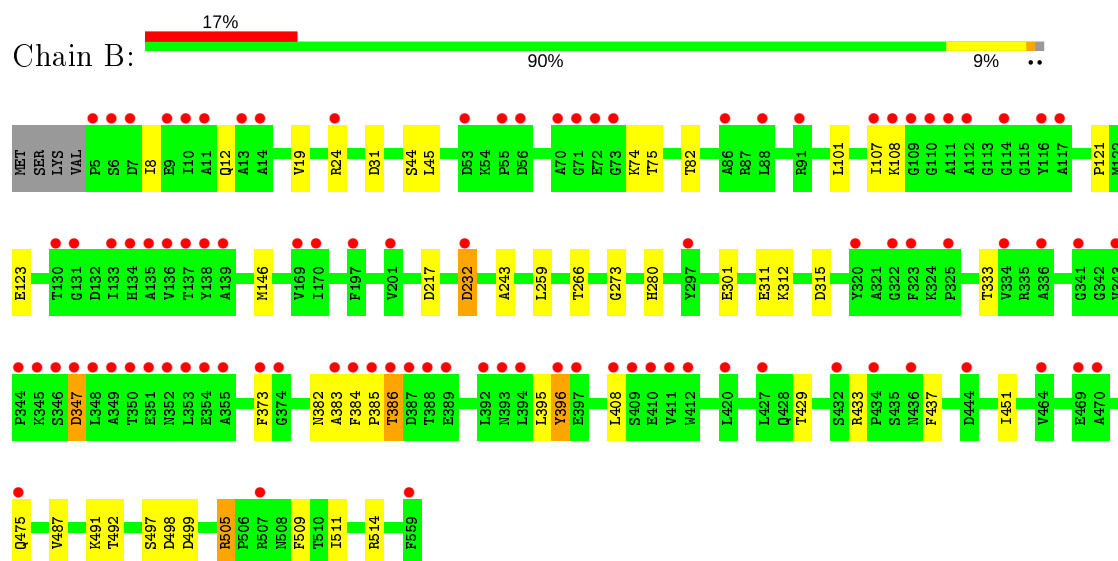
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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	Depositor
R, R_{free}	0.234 , 0.274 0.243 , 0.284	Depositor DCC
R_{free} test set	3753 reflections (5.04%)	wwPDB-VP
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
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	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, 0.333 respectively for untwinned datasets, and 0.375, 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:A:482:TYR:CE1	2:A:702:TOE:H7	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:THR:HA	1:B:301:GLU:OE1	2.17	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:347:ASP:HB3	4:A:913:HOH:O	2.18	0.43
1:A:558:LEU:HG	1:A:559:PHE:CZ	2.54	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%)	34	32
1	B	555/559 (99%)	521 (94%)	32 (6%)	2 (0%)	34	32
All	All	1113/1118 (100%)	1058 (95%)	51 (5%)	4 (0%)	34	32

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%)	57	63
1	B	414/441 (94%)	405 (98%)	9 (2%)	52	57
All	All	846/882 (96%)	829 (98%)	17 (2%)	60	60

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 [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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	$\# Z > 2$
3	SO4	B	602	-	4,4,4	0.70	0	6,6,6	0.98	0
2	TOE	A	702	-	10,10,10	0.77	0	9,9,9	0.82	0
3	SO4	A	705	-	4,4,4	0.95	0	6,6,6	0.67	0
2	TOE	B	601	-	10,10,10	0.67	0	9,9,9	0.53	0
3	SO4	B	604	-	4,4,4	0.44	0	6,6,6	0.21	0
2	TOE	A	701	-	10,10,10	1.24	0	9,9,9	1.67	3 (33%)
3	SO4	A	704	-	4,4,4	0.92	0	6,6,6	1.26	0
3	SO4	B	603	-	4,4,4	0.69	0	6,6,6	1.04	0
3	SO4	A	703	-	4,4,4	0.73	0	6,6,6	0.42	0
3	SO4	A	707	-	4,4,4	0.51	0	6,6,6	0.37	0
3	SO4	A	706	-	4,4,4	0.53	0	6,6,6	0.34	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	-	-	5/8/8/8	-
2	TOE	A	702	-	-	7/8/8/8	-
2	TOE	B	601	-	-	3/8/8/8	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	TOE	CG'-OF'-CE'	3.21	127.21	113.29
2	A	701	TOE	OF'-CG'-CH'	2.81	123.04	110.39
2	A	701	TOE	OC'-CD'-CE'	2.03	119.54	110.39

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	702	TOE	OC'-CD'-CE'-OF'
2	B	601	TOE	OF'-CG'-CH'-OI'
2	A	701	TOE	CH'-CG'-OF'-CE'
2	B	601	TOE	O2'-CA'-CB'-OC'
2	A	701	TOE	O2'-CA'-CB'-OC'
2	A	702	TOE	CE'-CD'-OC'-CB'
2	A	701	TOE	OC'-CD'-CE'-OF'
2	A	702	TOE	OF'-CG'-CH'-OI'
2	B	601	TOE	CD'-CE'-OF'-CG'
2	A	702	TOE	CG'-CH'-OI'-CK'
2	A	702	TOE	O2'-CA'-CB'-OC'
2	A	702	TOE	CA'-CB'-OC'-CD'
2	A	701	TOE	OF'-CG'-CH'-OI'
2	A	701	TOE	CE'-CD'-OC'-CB'
2	A	702	TOE	CD'-CE'-OF'-CG'

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	604	SO4	1	0
3	B	603	SO4	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	557/559 (99%)	0.80	76 (13%)  	41, 43, 47, 64	3 (0%)
1	B	555/559 (99%)	0.92	94 (16%)  	39, 54, 107, 143	1 (0%)
All	All	1112/1118 (99%)	0.86	170 (15%)  	39, 44, 97, 143	4 (0%)

All (170) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	353	LEU	6.6
1	B	384	PHE	6.3
1	A	438[A]	HIS	6.0
1	B	396	TYR	6.0
1	A	533[A]	MET	5.6
1	B	432	SER	5.3
1	A	559	PHE	5.2
1	A	169	VAL	4.9
1	B	559	PHE	4.9
1	B	355	ALA	4.8
1	A	335[A]	ARG	4.8
1	B	343	VAL	4.7
1	B	5	PRO	4.7
1	A	136	VAL	4.7
1	B	110	GLY	4.7
1	B	6	SER	4.7
1	B	389	GLU	4.7
1	A	133	ILE	4.5
1	B	385	PRO	4.5
1	B	409	SER	4.5
1	B	387	ASP	4.4
1	B	388	THR	4.4
1	B	350	THR	4.4
1	B	348	LEU	4.3

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Mol	Chain	Res	Type	RSRZ
1	B	386	THR	4.2
1	A	137	THR	4.1
1	B	116	TYR	4.2
1	B	346	SER	4.0
1	B	392	LEU	4.0
1	A	201	VAL	4.0
1	A	205	VAL	3.9
1	B	394	LEU	3.9
1	B	475	GLN	3.8
1	A	170	ILE	3.8
1	B	411	VAL	3.7
1	B	322	GLY	3.7
1	B	349	ALA	3.7
1	A	135	ALA	3.6
1	B	344	PRO	3.6
1	B	73	GLY	3.5
1	B	136	VAL	3.4
1	B	170	ILE	3.4
1	B	133	ILE	3.3
1	B	55	PRO	3.3
1	B	334	VAL	3.3
1	B	169	VAL	3.2
1	A	329	VAL	3.2
1	A	551	ALA	3.2
1	A	274	PRO	3.2
1	B	336	ALA	3.2
1	B	345	LYS	3.2
1	A	142	LEU	3.1
1	A	557	GLY	3.1
1	B	137	THR	3.1
1	A	197	PHE	3.1
1	B	393	ASN	3.1
1	B	412	TRP	3.1
1	A	556	THR	3.1
1	A	206	MET	3.0
1	A	250	LEU	3.0
1	B	351	GLU	3.0
1	B	114	GLY	3.0
1	A	139	ALA	3.0
1	A	207	ALA	3.0
1	A	554	VAL	2.9
1	B	323	PHE	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	208	CYS	2.9
1	B	13	ALA	2.9
1	B	427	LEU	2.9
1	A	200	SER	2.9
1	A	308	LEU	2.9
1	B	347	ASP	2.8
1	B	341	GLY	2.8
1	B	53	ASP	2.8
1	A	282	CYS	2.8
1	B	434	PRO	2.8
1	A	141	ASN	2.8
1	B	297	TYR	2.8
1	B	10	ILE	2.7
1	A	553	GLY	2.7
1	B	420	LEU	2.7
1	B	91	ARG	2.7
1	A	302	ALA	2.7
1	B	436	ASN	2.7
1	B	107	ILE	2.7
1	B	88	LEU	2.6
1	A	272	GLY	2.6
1	B	72	GLU	2.6
1	B	14	ALA	2.6
1	A	12	GLN	2.6
1	A	552	ASP	2.6
1	B	9	GLU	2.6
1	B	469	GLU	2.6
1	B	134	HIS	2.6
1	B	112	ALA	2.5
1	A	138	TYR	2.5
1	B	130	THR	2.5
1	B	7	ASP	2.5
1	B	320	TYR	2.5
1	A	203	SER	2.5
1	A	199	ILE	2.5
1	A	140	HIS	2.5
1	B	374	GLY	2.5
1	A	13	ALA	2.5
1	A	331	VAL	2.5
1	B	373	PHE	2.5
1	A	202	ALA	2.5
1	A	279	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	383	ALA	2.4
1	A	61	LEU	2.4
1	B	410	GLU	2.4
1	B	397	GLU	2.4
1	B	109	GLY	2.3
1	A	96	LEU	2.3
1	A	166	TRP	2.3
1	A	177	LEU	2.3
1	B	71	GLY	2.3
1	A	134	HIS	2.3
1	A	131	GLY	2.3
1	A	62	VAL	2.3
1	A	144	ALA	2.3
1	A	65	ILE	2.3
1	B	111	ALA	2.3
1	B	117	ALA	2.3
1	B	56	ASP	2.3
1	B	408	LEU	2.2
1	A	280	HIS	2.2
1	B	352	ASN	2.2
1	B	108	LYS	2.2
1	A	558	LEU	2.2
1	A	275	PHE	2.2
1	A	309	GLY	2.2
1	A	546	ASN	2.2
1	B	464	VAL	2.2
1	B	325	PRO	2.2
1	A	299	VAL	2.2
1	A	63	THR	2.2
1	B	135	ALA	2.2
1	A	306	ALA	2.2
1	B	86	ALA	2.2
1	A	454	ILE	2.1
1	A	545	CYS	2.1
1	A	8	ILE	2.1
1	A	305	GLY	2.1
1	A	99	PRO	2.1
1	A	130	THR	2.1
1	B	70	ALA	2.1
1	A	168	ARG	2.1
1	A	487	VAL	2.1
1	B	131	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	3	LYS	2.1
1	A	550	ASP	2.1
1	B	232	ASP	2.1
1	A	422	LEU	2.1
1	A	314	TYR	2.1
1	B	138	TYR	2.1
1	B	201	VAL	2.1
1	B	139	ALA	2.1
1	B	507	ARG	2.1
1	B	197	PHE	2.1
1	A	53	ASP	2.0
1	A	278	ILE	2.0
1	B	11	ALA	2.0
1	B	470	ALA	2.0
1	A	132	ASP	2.0
1	A	143	LEU	2.0
1	B	354	GLU	2.0
1	B	24[A]	ARG	2.0
1	B	444	ASP	2.0
1	A	209	LEU	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 monosaccharides 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. 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	B-factors(\AA^2)	Q<0.9
3	SO4	B	604	5/5	0.82	0.31	80,86,102,113	0
3	SO4	A	704	5/5	0.87	0.17	42,43,50,61	0
3	SO4	A	706	5/5	0.89	0.34	67,72,83,101	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	TOE	A	701	11/11	0.90	0.14	23,34,47,47	0
3	SO4	A	707	5/5	0.92	0.33	59,77,92,98	0
2	TOE	A	702	11/11	0.92	0.22	33,39,59,60	0
2	TOE	B	601	11/11	0.93	0.15	33,38,46,48	0
3	SO4	B	603	5/5	0.94	0.11	41,55,63,64	0
3	SO4	B	602	5/5	0.95	0.11	37,38,47,49	0
3	SO4	A	705	5/5	0.97	0.24	34,38,51,60	0
3	SO4	A	703	5/5	0.98	0.09	21,25,26,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.