



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 7, 2017 – 01:24 AM EDT

PDB ID : 5TCG
Title : Crystal structure of tryptophan synthase from *M. tuberculosis* - aminoacylate-bound form
Authors : Michalska, K.; Maltseva, N.; Jedrzejczak, R.; Joachimiak, A.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : unknown
Resolution : 2.40 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20029824

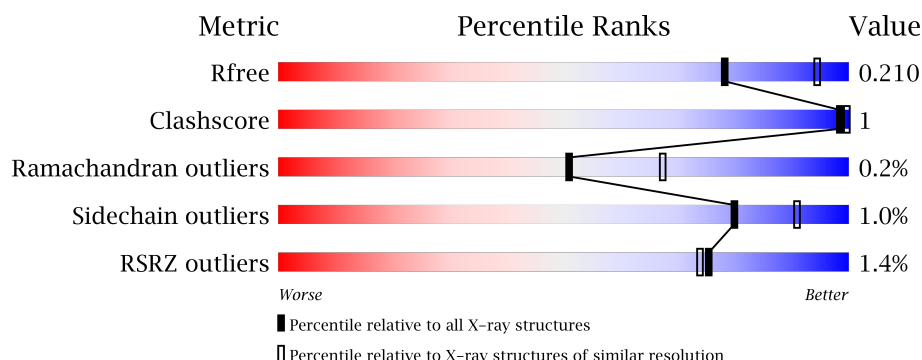
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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.



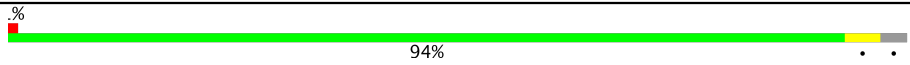
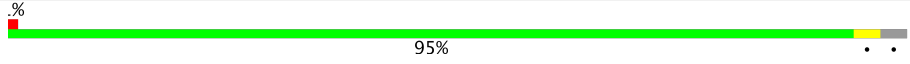
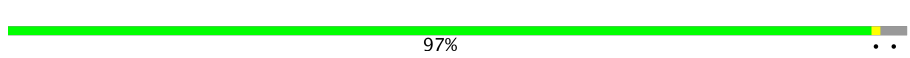
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	3166 (2.40-2.40)
Clashscore	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (2.40-2.40)
RSRZ outliers	101464	3195 (2.40-2.40)

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	276	<div> <div>0.2%</div> <div>89%</div> <div>10%</div> </div>
1	C	276	<div> <div>3%</div> <div>87%</div> <div>10%</div> </div>
1	E	276	<div> <div>4%</div> <div>87%</div> <div>11%</div> </div>
1	G	276	<div> <div>0.2%</div> <div>87%</div> <div>10%</div> </div>
2	B	410	<div> <div>94%</div> <div>10%</div> </div>

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Mol	Chain	Length	Quality of chain
2	D	410	
2	F	410	
2	H	410	

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	FMT	B	506	-	-	-	X
4	FMT	B	508	-	-	-	X
4	FMT	C	302	-	-	-	X
4	FMT	F	505	-	-	-	X
4	FMT	G	302	-	-	-	X
4	FMT	G	304	-	-	-	X
4	FMT	H	508	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 20280 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 Tryptophan synthase alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	249	Total	C	N	O	S	0	1	0
			1812	1136	325	346	5			
1	G	248	Total	C	N	O	S	0	1	0
			1802	1131	323	342	6			
1	E	245	Total	C	N	O	S	0	1	0
			1784	1120	320	338	6			
1	C	249	Total	C	N	O	S	0	1	0
			1811	1136	324	345	6			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	271	HIS	-	expression tag	UNP P9WFY1
A	272	HIS	-	expression tag	UNP P9WFY1
A	273	HIS	-	expression tag	UNP P9WFY1
A	274	HIS	-	expression tag	UNP P9WFY1
A	275	HIS	-	expression tag	UNP P9WFY1
A	276	HIS	-	expression tag	UNP P9WFY1
G	271	HIS	-	expression tag	UNP P9WFY1
G	272	HIS	-	expression tag	UNP P9WFY1
G	273	HIS	-	expression tag	UNP P9WFY1
G	274	HIS	-	expression tag	UNP P9WFY1
G	275	HIS	-	expression tag	UNP P9WFY1
G	276	HIS	-	expression tag	UNP P9WFY1
E	271	HIS	-	expression tag	UNP P9WFY1
E	272	HIS	-	expression tag	UNP P9WFY1
E	273	HIS	-	expression tag	UNP P9WFY1
E	274	HIS	-	expression tag	UNP P9WFY1
E	275	HIS	-	expression tag	UNP P9WFY1
E	276	HIS	-	expression tag	UNP P9WFY1
C	271	HIS	-	expression tag	UNP P9WFY1
C	272	HIS	-	expression tag	UNP P9WFY1
C	273	HIS	-	expression tag	UNP P9WFY1

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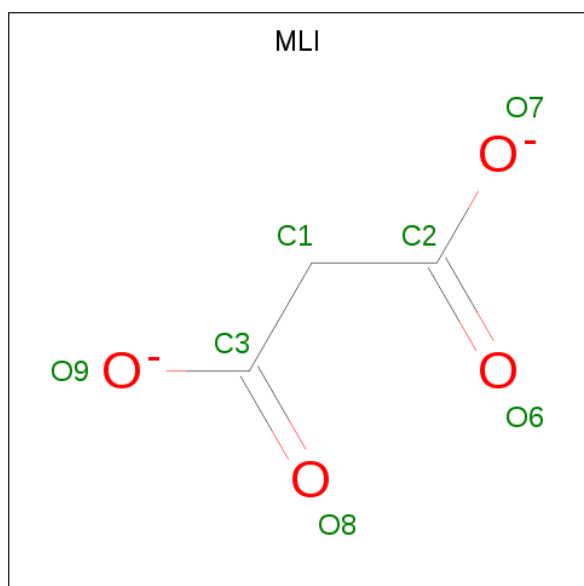
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Chain	Residue	Modelled	Actual	Comment	Reference
C	274	HIS	-	expression tag	UNP P9WFY1
C	275	HIS	-	expression tag	UNP P9WFY1
C	276	HIS	-	expression tag	UNP P9WFY1

- Molecule 2 is a protein called Tryptophan synthase beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	399	Total	C	N	O	S	0	3	0
			3000	1873	544	569	14			
2	H	399	Total	C	N	O	S	0	2	0
			2989	1867	541	567	14			
2	F	399	Total	C	N	O	S	0	3	0
			3000	1872	545	570	13			
2	D	399	Total	C	N	O	S	0	2	0
			2994	1869	546	566	13			

- Molecule 3 is MALONATE ION (three-letter code: MLI) (formula: $C_3H_2O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	3	4		
3	G	1	Total	C	O	0	0
			7	3	4		
3	E	1	Total	C	O	0	0
			7	3	4		
3	C	1	Total	C	O	0	0
			7	3	4		

- Molecule 4 is FORMIC ACID (three-letter code: FMT) (formula: CH₂O₂).



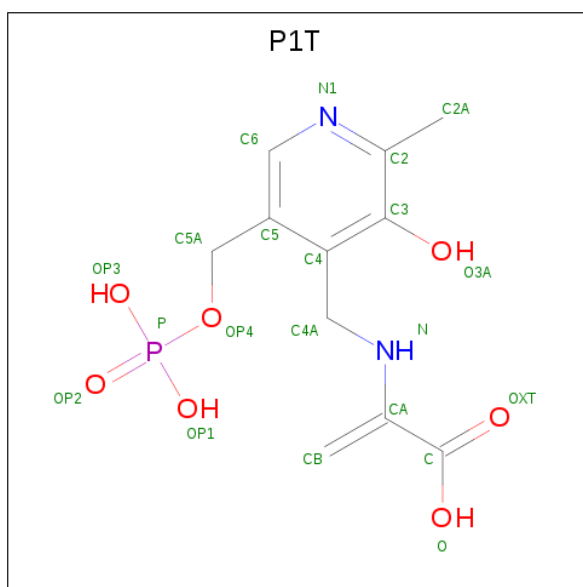
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			3	1	2		
4	B	1	Total	C	O	0	0
			3	1	2		
4	B	1	Total	C	O	0	0
			3	1	2		
4	B	1	Total	C	O	0	0
			3	1	2		
4	B	1	Total	C	O	0	0
			3	1	2		
4	B	1	Total	C	O	0	0
			3	1	2		
4	G	1	Total	C	O	0	0
			3	1	2		
4	G	1	Total	C	O	0	0
			3	1	2		
4	G	1	Total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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	H	1	Total	C	O	0	0
			3	1	2		
4	H	1	Total	C	O	0	0
			3	1	2		
4	H	1	Total	C	O	0	0
			3	1	2		
4	F	1	Total	C	O	0	0
			3	1	2		
4	F	1	Total	C	O	0	0
			3	1	2		
4	F	1	Total	C	O	0	0
			3	1	2		
4	F	1	Total	C	O	0	0
			3	1	2		
4	C	1	Total	C	O	0	0
			3	1	2		
4	D	1	Total	C	O	0	0
			3	1	2		
4	D	1	Total	C	O	0	0
			3	1	2		
4	D	1	Total	C	O	0	0
			3	1	2		
4	D	1	Total	C	O	0	0
			3	1	2		

- Molecule 5 is 2-[(3-HYDROXY-2-METHYL-5-[(PHOSPHONOOXY)METHYL]PYRIDIN-4-YL)METHYL)AMINO]ACRYLIC ACID (three-letter code: P1T) (formula: C₁₁H₁₅N₂O₇P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	P	0	0
			21	11	2	7	1		
5	H	1	Total	C	N	O	P	0	0
			21	11	2	7	1		
5	F	1	Total	C	N	O	P	0	0
			21	11	2	7	1		
5	D	1	Total	C	N	O	P	0	0
			21	11	2	7	1		

- Molecule 6 is CESIUM ION (three-letter code: CS) (formula: Cs).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	H	1	Total	Cs	0	0
			1	1		
6	B	2	Total	Cs	0	0
			2	2		
6	D	3	Total	Cs	0	0
			3	3		
6	F	3	Total	Cs	0	0
			3	3		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	57	Total	O	0	0
			57	57		

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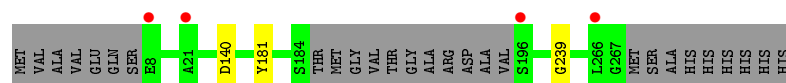
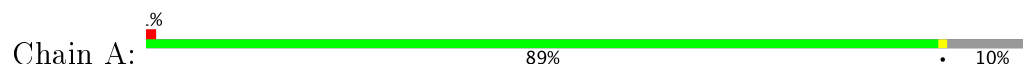
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	170	Total 170	O 170	0	0
7	G	82	Total 82	O 82	0	2
7	H	186	Total 187	O 187	0	2
7	E	39	Total 39	O 39	0	0
7	F	140	Total 140	O 140	0	2
7	C	63	Total 63	O 63	0	0
7	D	153	Total 154	O 154	0	1

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 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.

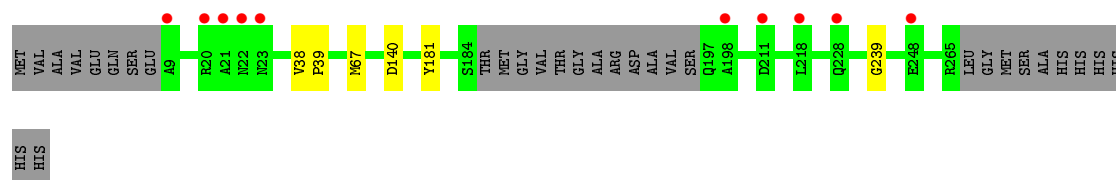
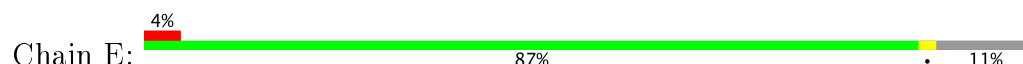
- Molecule 1: Tryptophan synthase alpha chain



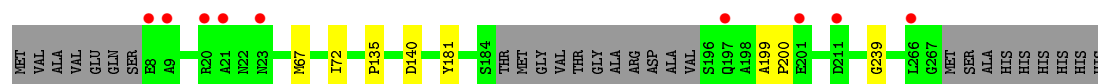
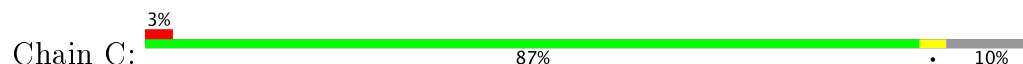
- Molecule 1: Tryptophan synthase alpha chain



- Molecule 1: Tryptophan synthase alpha chain



- Molecule 1: Tryptophan synthase alpha chain



- Molecule 2: Tryptophan synthase beta chain



● Molecule 2: Tryptophan synthase beta chain

Chain H:

97%



● Molecule 2: Tryptophan synthase beta chain

Chain F:

.%

95%



● Molecule 2: Tryptophan synthase beta chain

Chain D:

.%

94%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	135.18Å 159.43Å 165.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.40 29.91 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.5 (30.00-2.40) 99.6 (29.91-2.40)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.77 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.172 , 0.206 0.178 , 0.210	Depositor DCC
R_{free} test set	2704 reflections (1.98%)	DCC
Wilson B-factor (Å ²)	27.8	Xtriage
Anisotropy	0.248	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.010 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	20280	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.62% 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: CS, FMT, MLI, P1T

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	0.55	0/1841	0.71	0/2512
1	C	0.53	0/1840	0.71	0/2510
1	E	0.51	0/1813	0.68	0/2474
1	G	0.56	0/1831	0.73	0/2498
2	B	0.59	0/3060	0.75	0/4145
2	D	0.59	0/3054	0.74	0/4137
2	F	0.58	0/3060	0.73	1/4146 (0.0%)
2	H	0.61	0/3049	0.74	0/4131
All	All	0.57	0/19548	0.73	1/26553 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	280	ARG	NE-CZ-NH2	5.07	122.83	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	1812	0	1829	0	0
1	C	1811	0	1830	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1784	0	1805	2	0
1	G	1802	0	1824	3	0
2	B	3000	0	2920	7	0
2	D	2994	0	2919	6	0
2	F	3000	0	2917	3	0
2	H	2989	0	2910	1	0
3	A	7	0	2	0	0
3	C	7	0	2	0	0
3	E	7	0	2	0	0
3	G	7	0	2	0	0
4	A	3	0	1	0	0
4	B	15	0	5	0	0
4	C	3	0	1	0	0
4	D	12	0	4	0	0
4	F	12	0	4	0	0
4	G	9	0	3	0	0
4	H	21	0	7	0	0
5	B	21	0	11	0	0
5	D	21	0	12	0	0
5	F	21	0	12	0	0
5	H	21	0	11	0	0
6	B	2	0	0	0	0
6	D	3	0	0	0	0
6	F	3	0	0	0	0
6	H	1	0	0	0	0
7	A	57	0	0	0	0
7	B	170	0	0	0	0
7	C	63	0	0	1	0
7	D	154	0	0	1	0
7	E	39	0	0	0	0
7	F	140	0	0	0	0
7	G	82	0	0	0	0
7	H	187	0	0	0	0
All	All	20280	0	19033	25	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 1.

All (25) 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)
2:B:406:LEU:O	2:B:407:LEU:HB2	2.04	0.56
2:B:265[B]:ARG:NH2	2:B:378[B]:GLU:OE1	2.35	0.56
2:D:207:GLY:HA2	2:D:294:HIS:O	2.14	0.47
2:D:129:HIS:CE1	2:D:203:GLY:HA2	2.50	0.46
1:E:67[B]:MET:HE2	1:E:67[B]:MET:HB3	1.86	0.46
2:H:207:GLY:HA2	2:H:294:HIS:O	2.16	0.46
1:C:199:ALA:HB3	1:C:200:PRO:HD3	1.98	0.45
2:B:129:HIS:CE1	2:B:203:GLY:HA2	2.52	0.44
2:F:34:LEU:HG	2:F:192:VAL:HG22	1.99	0.44
1:C:72:ILE:HD12	7:C:411:HOH:O	2.16	0.44
2:B:123:GLU:HG3	2:B:184:ILE:CG1	2.48	0.43
1:E:38:VAL:HB	1:E:39:PRO:HD3	2.00	0.43
1:G:72:ILE:O	1:G:76:THR:HG23	2.18	0.43
2:D:123:GLU:HG3	2:D:184:ILE:CG1	2.49	0.42
2:F:284:THR:HG22	2:F:299:TYR:HB3	2.01	0.42
1:G:134:THR:HB	1:G:137:LEU:HB3	2.02	0.41
2:B:128:GLN:OE1	2:B:395:LYS:HD3	2.21	0.41
2:B:123:GLU:HG3	2:B:184:ILE:HG12	2.03	0.41
1:C:67[B]:MET:SD	1:C:135:PRO:HB2	2.61	0.41
2:D:93:ASP:HB2	2:D:392:ARG:HB3	2.02	0.41
2:B:146:ILE:HG21	2:B:163:MET:SD	2.61	0.41
2:D:62:GLN:NE2	7:D:602:HOH:O	2.52	0.41
2:D:159:ASN:OD1	2:D:162:ARG:NH1	2.54	0.40
2:F:285:PHE:CZ	2:F:323:VAL:HG11	2.55	0.40
1:G:61:PRO:HG2	1:G:84:VAL:HG11	2.02	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	246/276 (89%)	243 (99%)	2 (1%)	1 (0%)	38 54

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	246/276 (89%)	241 (98%)	4 (2%)	1 (0%)	38	54
1	E	242/276 (88%)	238 (98%)	3 (1%)	1 (0%)	38	54
1	G	245/276 (89%)	241 (98%)	3 (1%)	1 (0%)	38	54
2	B	400/410 (98%)	390 (98%)	10 (2%)	0	100	100
2	D	399/410 (97%)	389 (98%)	10 (2%)	0	100	100
2	F	400/410 (98%)	389 (97%)	11 (3%)	0	100	100
2	H	399/410 (97%)	392 (98%)	7 (2%)	0	100	100
All	All	2577/2744 (94%)	2523 (98%)	50 (2%)	4 (0%)	51	67

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	239	GLY
1	A	239	GLY
1	G	239	GLY
1	E	239	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	180/200 (90%)	178 (99%)	2 (1%)	78	90
1	C	180/200 (90%)	178 (99%)	2 (1%)	78	90
1	E	177/200 (88%)	175 (99%)	2 (1%)	78	90
1	G	179/200 (90%)	177 (99%)	2 (1%)	78	90
2	B	298/302 (99%)	296 (99%)	2 (1%)	87	94
2	D	297/302 (98%)	292 (98%)	5 (2%)	66	82
2	F	298/302 (99%)	294 (99%)	4 (1%)	73	87
2	H	297/302 (98%)	296 (100%)	1 (0%)	94	98
All	All	1906/2008 (95%)	1886 (99%)	20 (1%)	80	91

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	140	ASP
1	A	181	TYR
2	B	80	GLN
2	B	200	TYR
1	G	140	ASP
1	G	181	TYR
2	H	300	LEU
1	E	140	ASP
1	E	181	TYR
2	F	23	SER
2	F	200	TYR
2	F	300	LEU
2	F	348	ASP
1	C	140	ASP
1	C	181	TYR
2	D	23	SER
2	D	164[A]	ARG
2	D	164[B]	ARG
2	D	200	TYR
2	D	256	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 42 ligands modelled in this entry, 9 are monoatomic - leaving 33 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	MLI	A	301	-	0,6,6	0.00	-	0,7,7	0.00	-
4	FMT	A	302	-	0,2,2	0.00	-	0,1,1	0.00	-
5	P1T	B	501	-	19,21,21	3.20	4 (21%)	22,30,30	1.43	4 (18%)
4	FMT	B	504	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	B	505	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	B	506	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	B	507	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	B	508	-	0,2,2	0.00	-	0,1,1	0.00	-
3	MLI	C	301	-	0,6,6	0.00	-	0,7,7	0.00	-
4	FMT	C	302	-	0,2,2	0.00	-	0,1,1	0.00	-
5	P1T	D	501	-	19,21,21	3.52	4 (21%)	22,30,30	1.57	3 (13%)
4	FMT	D	505	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	D	506	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	D	507	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	D	508	-	0,2,2	0.00	-	0,1,1	0.00	-
3	MLI	E	301	-	0,6,6	0.00	-	0,7,7	0.00	-
5	P1T	F	501	-	19,21,21	3.15	4 (21%)	22,30,30	1.51	3 (13%)
4	FMT	F	505	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	F	506	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	F	507	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	F	508	-	0,2,2	0.00	-	0,1,1	0.00	-
3	MLI	G	301	-	0,6,6	0.00	-	0,7,7	0.00	-
4	FMT	G	302	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	G	303	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	G	304	-	0,2,2	0.00	-	0,1,1	0.00	-
5	P1T	H	501	-	19,21,21	3.20	4 (21%)	22,30,30	1.42	4 (18%)
4	FMT	H	503	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	H	504	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	H	505	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	H	506	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	H	507	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	H	508	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	H	509	-	0,2,2	0.00	-	0,1,1	0.00	-

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	MLI	A	301	-	-	0/0/4/4	0/0/0/0
4	FMT	A	302	-	-	0/0/0/0	0/0/0/0
5	P1T	B	501	-	-	0/10/15/15	0/1/1/1
4	FMT	B	504	-	-	0/0/0/0	0/0/0/0
4	FMT	B	505	-	-	0/0/0/0	0/0/0/0
4	FMT	B	506	-	-	0/0/0/0	0/0/0/0
4	FMT	B	507	-	-	0/0/0/0	0/0/0/0
4	FMT	B	508	-	-	0/0/0/0	0/0/0/0
3	MLI	C	301	-	-	0/0/4/4	0/0/0/0
4	FMT	C	302	-	-	0/0/0/0	0/0/0/0
5	P1T	D	501	-	-	0/10/15/15	0/1/1/1
4	FMT	D	505	-	-	0/0/0/0	0/0/0/0
4	FMT	D	506	-	-	0/0/0/0	0/0/0/0
4	FMT	D	507	-	-	0/0/0/0	0/0/0/0
4	FMT	D	508	-	-	0/0/0/0	0/0/0/0
3	MLI	E	301	-	-	0/0/4/4	0/0/0/0
5	P1T	F	501	-	-	0/10/15/15	0/1/1/1
4	FMT	F	505	-	-	0/0/0/0	0/0/0/0
4	FMT	F	506	-	-	0/0/0/0	0/0/0/0
4	FMT	F	507	-	-	0/0/0/0	0/0/0/0
4	FMT	F	508	-	-	0/0/0/0	0/0/0/0
3	MLI	G	301	-	-	0/0/4/4	0/0/0/0
4	FMT	G	302	-	-	0/0/0/0	0/0/0/0
4	FMT	G	303	-	-	0/0/0/0	0/0/0/0
4	FMT	G	304	-	-	0/0/0/0	0/0/0/0
5	P1T	H	501	-	-	0/10/15/15	0/1/1/1
4	FMT	H	503	-	-	0/0/0/0	0/0/0/0
4	FMT	H	504	-	-	0/0/0/0	0/0/0/0
4	FMT	H	505	-	-	0/0/0/0	0/0/0/0
4	FMT	H	506	-	-	0/0/0/0	0/0/0/0
4	FMT	H	507	-	-	0/0/0/0	0/0/0/0
4	FMT	H	508	-	-	0/0/0/0	0/0/0/0
4	FMT	H	509	-	-	0/0/0/0	0/0/0/0

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	501	P1T	C-CA	-4.12	1.45	1.52
5	D	501	P1T	C-CA	-4.09	1.45	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	H	501	P1T	C-CA	-4.02	1.45	1.52
5	F	501	P1T	C-CA	-3.53	1.46	1.52
5	H	501	P1T	C3-C4	4.48	1.47	1.40
5	F	501	P1T	C3-C4	4.71	1.47	1.40
5	B	501	P1T	C3-C4	5.00	1.48	1.40
5	H	501	P1T	C5-C4	5.05	1.47	1.40
5	F	501	P1T	C5-C4	5.32	1.47	1.40
5	B	501	P1T	C5-C4	5.61	1.48	1.40
5	D	501	P1T	C3-C4	5.91	1.49	1.40
5	D	501	P1T	C5-C4	6.05	1.48	1.40
5	B	501	P1T	C3-C2	10.70	1.48	1.40
5	F	501	P1T	C3-C2	11.01	1.48	1.40
5	H	501	P1T	C3-C2	11.22	1.48	1.40
5	D	501	P1T	C3-C2	11.87	1.49	1.40

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	501	P1T	C4A-C4-C5	-2.05	117.88	119.75
5	B	501	P1T	C4-C4A-N	2.07	115.57	111.13
5	H	501	P1T	C6-N1-C2	2.13	123.37	119.26
5	F	501	P1T	C6-N1-C2	2.24	123.57	119.26
5	D	501	P1T	C6-N1-C2	2.24	123.57	119.26
5	H	501	P1T	C4-C4A-N	2.25	115.96	111.13
5	B	501	P1T	O3A-C3-C2	2.31	122.62	117.78
5	H	501	P1T	O3A-C3-C2	2.39	122.79	117.78
5	F	501	P1T	C4-C4A-N	2.49	116.47	111.13
5	B	501	P1T	C6-N1-C2	2.51	124.10	119.26
5	B	501	P1T	C4A-C4-C3	3.40	123.40	119.65
5	H	501	P1T	C4A-C4-C3	3.56	123.58	119.65
5	F	501	P1T	C4A-C4-C3	3.91	123.96	119.65
5	D	501	P1T	C4A-C4-C3	4.53	124.65	119.65

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	249/276 (90%)	-0.32	4 (1%) 72 70	25, 36, 53, 80	0
1	C	249/276 (90%)	-0.14	9 (3%) 43 42	23, 42, 72, 91	0
1	E	245/276 (88%)	0.09	10 (4%) 38 36	31, 55, 80, 96	0
1	G	248/276 (89%)	-0.46	2 (0%) 86 84	22, 34, 46, 61	0
2	B	399/410 (97%)	-0.58	0 100 100	18, 25, 38, 58	0
2	D	399/410 (97%)	-0.55	5 (1%) 77 75	17, 25, 37, 68	0
2	F	399/410 (97%)	-0.56	4 (1%) 82 80	17, 26, 41, 66	0
2	H	399/410 (97%)	-0.58	2 (0%) 90 89	16, 24, 38, 59	0
All	All	2587/2744 (94%)	-0.43	36 (1%) 75 74	16, 29, 61, 96	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	8	GLU	4.7
2	D	22	PRO	4.7
1	E	23	ASN	4.7
2	F	407	LEU	4.3
1	A	196	SER	4.2
2	F	21	GLY	3.6
1	C	197	GLN	3.6
1	E	211	ASP	3.6
2	F	22	PRO	3.6
2	D	407	LEU	3.4
1	A	8	GLU	3.4
2	D	9	THR	3.4
1	C	21	ALA	3.2
1	C	9	ALA	3.0
2	D	21	GLY	2.8
1	E	21	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
2	F	9	THR	2.7
1	G	196	SER	2.6
1	E	9	ALA	2.6
1	E	248	GLU	2.6
1	A	21	ALA	2.6
1	E	20	ARG	2.6
1	C	211	ASP	2.6
1	G	197	GLN	2.6
1	A	266	LEU	2.4
2	H	22	PRO	2.4
1	E	228	GLN	2.3
1	C	20	ARG	2.3
1	E	22	ASN	2.3
1	E	198	ALA	2.3
1	C	201	GLU	2.2
1	C	23	ASN	2.2
1	C	266	LEU	2.2
2	H	277	GLU	2.1
1	E	218	LEU	2.1
2	D	23	SER	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	FMT	F	505	3/3	0.88	0.26	8.02	46,46,48,49	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	FMT	G	302	3/3	0.62	0.33	5.74	53,53,54,57	0
4	FMT	C	302	3/3	0.85	0.27	4.31	48,48,49,50	0
4	FMT	H	508	3/3	0.81	0.19	4.18	52,52,57,57	0
4	FMT	B	506	3/3	0.90	0.16	3.24	41,41,43,44	0
4	FMT	B	508	3/3	0.90	0.19	2.62	50,50,51,52	0
4	FMT	G	304	3/3	0.84	0.21	2.20	44,44,45,47	0
4	FMT	G	303	3/3	0.90	0.20	1.70	42,42,45,45	0
3	MLI	A	301	7/7	0.89	0.16	1.31	45,46,47,49	0
4	FMT	D	508	3/3	0.85	0.18	1.01	53,53,55,55	0
4	FMT	H	509	3/3	0.77	0.20	0.85	59,59,61,61	0
5	P1T	B	501	21/21	0.98	0.18	0.77	20,21,24,24	0
4	FMT	B	505	3/3	0.94	0.17	0.68	43,43,44,46	0
4	FMT	D	507	3/3	0.93	0.14	0.62	49,49,49,51	0
4	FMT	F	508	3/3	0.96	0.15	0.55	58,58,58,59	0
3	MLI	C	301	7/7	0.90	0.15	0.44	50,53,57,59	0
5	P1T	H	501	21/21	0.98	0.19	0.38	18,19,23,24	0
5	P1T	F	501	21/21	0.99	0.16	0.25	19,20,21,22	0
4	FMT	A	302	3/3	0.93	0.14	0.19	42,42,45,45	0
4	FMT	H	506	3/3	0.97	0.13	0.06	40,40,41,42	0
4	FMT	F	507	3/3	0.88	0.17	-0.02	50,50,52,54	0
5	P1T	D	501	21/21	0.99	0.15	-0.07	18,19,22,23	0
3	MLI	G	301	7/7	0.93	0.12	-0.10	48,50,53,54	0
3	MLI	E	301	7/7	0.91	0.14	-0.16	55,59,60,62	0
4	FMT	D	505	3/3	0.87	0.15	-0.37	44,44,45,45	0
4	FMT	H	503	3/3	0.95	0.10	-0.52	38,38,40,41	0
4	FMT	H	507	3/3	0.94	0.10	-0.88	53,53,53,54	0
6	CS	F	504	1/1	0.97	0.07	-0.91	61,61,61,61	1
6	CS	B	503	1/1	0.96	0.06	-1.24	36,36,36,36	1
4	FMT	B	507	3/3	0.95	0.08	-1.46	50,50,51,51	0
6	CS	D	503	1/1	0.99	0.05	-2.11	47,47,47,47	1
6	CS	H	502	1/1	0.99	0.04	-2.11	52,52,52,52	1
6	CS	D	502	1/1	0.99	0.03	-2.18	55,55,55,55	1
6	CS	D	504	1/1	0.94	0.04	-2.41	65,65,65,65	1
6	CS	B	502	1/1	0.99	0.03	-2.70	54,54,54,54	1
6	CS	F	502	1/1	0.98	0.03	-3.25	58,58,58,58	1
6	CS	F	503	1/1	0.99	0.04	-3.92	45,45,45,45	1
4	FMT	H	505	3/3	0.85	0.23	-	45,45,46,48	0
4	FMT	B	504	3/3	0.90	0.12	-	30,30,33,34	0
4	FMT	F	506	3/3	0.97	0.12	-	33,33,34,35	0
4	FMT	D	506	3/3	0.89	0.15	-	35,35,36,37	0
4	FMT	H	504	3/3	0.93	0.13	-	29,29,31,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.