



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 06:41 pm BST

PDB ID : 6CD0
Title : Crystal structure of Medicago truncatula serine hydroxymethyltransferase 3 (MtSHMT3), PLP-internal aldimine and apo form
Authors : Ruszkowski, M.; Sekula, B.; Ruszkowska, A.; Dauter, Z.
Deposited on : 2018-02-07
Resolution : 1.74 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 : 2.11
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) : 2.11

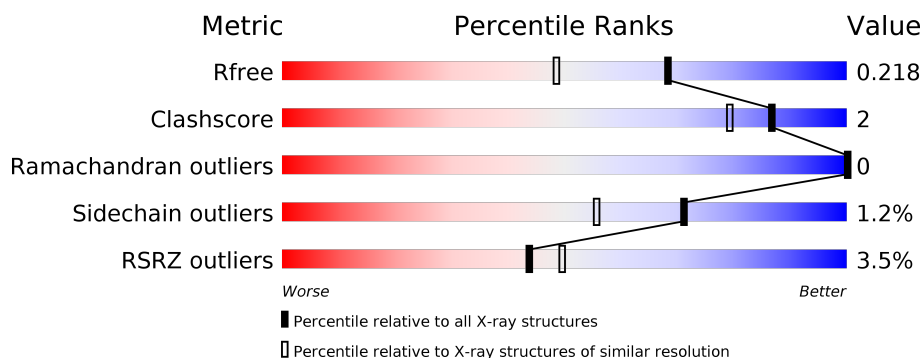
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.74 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	3764 (1.76-1.72)
Clashscore	141614	3923 (1.76-1.72)
Ramachandran outliers	138981	3878 (1.76-1.72)
Sidechain outliers	138945	3878 (1.76-1.72)
RSRZ outliers	127900	3705 (1.76-1.72)

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	455	<div> <div>3%</div> <div>95%</div> <div>6%</div> <div>• •</div> </div>
1	B	455	<div> <div>3%</div> <div>93%</div> <div>6%</div> <div>•</div> </div>
1	D	455	<div> <div>4%</div> <div>96%</div> <div>• •</div> </div>
2	C	455	<div> <div>4%</div> <div>95%</div> <div>• •</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15607 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 Serine hydroxymethyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	452	Total	C	N	O	P	S	0	1	0
			3492	2212	608	656	1	15			
1	B	452	Total	C	N	O	P	S	0	1	0
			3493	2213	609	655	1	15			
1	D	452	Total	C	N	O	P	S	0	6	0
			3527	2232	613	666	1	15			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	79	SER	-	expression tag	UNP G7ILW0
A	80	ASN	-	expression tag	UNP G7ILW0
A	81	ALA	-	expression tag	UNP G7ILW0
B	79	SER	-	expression tag	UNP G7ILW0
B	80	ASN	-	expression tag	UNP G7ILW0
B	81	ALA	-	expression tag	UNP G7ILW0
D	79	SER	-	expression tag	UNP G7ILW0
D	80	ASN	-	expression tag	UNP G7ILW0
D	81	ALA	-	expression tag	UNP G7ILW0

- Molecule 2 is a protein called Serine hydroxymethyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	452	Total	C	N	O	S	0	24	0
			3662	2318	640	689	15			

There are 3 discrepancies between the modelled and reference sequences:

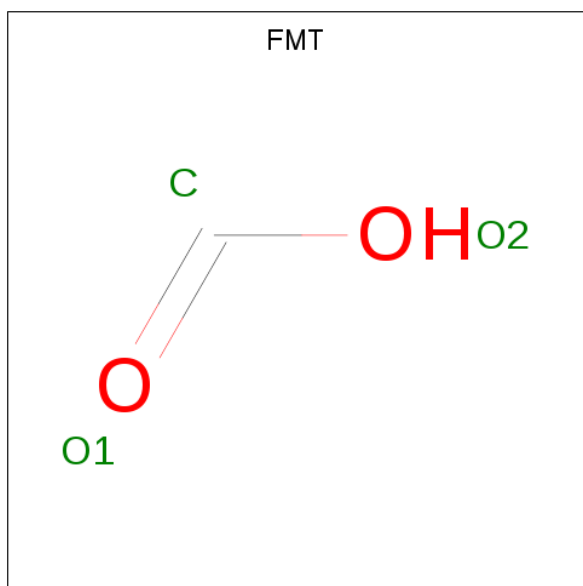
Chain	Residue	Modelled	Actual	Comment	Reference
C	79	SER	-	expression tag	UNP G7ILW0
C	80	ASN	-	expression tag	UNP G7ILW0
C	81	ALA	-	expression tag	UNP G7ILW0

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 4 is FORMIC ACID (three-letter code: FMT) (formula: CH_2O_2).



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

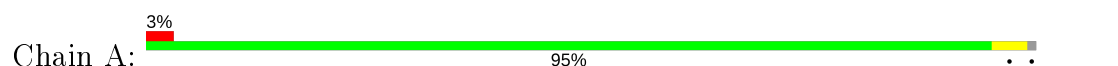
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	318	Total O 318 318	0	0
5	B	313	Total O 313 313	0	0
5	C	365	Total O 365 365	0	0
5	D	413	Total O 413 413	0	0

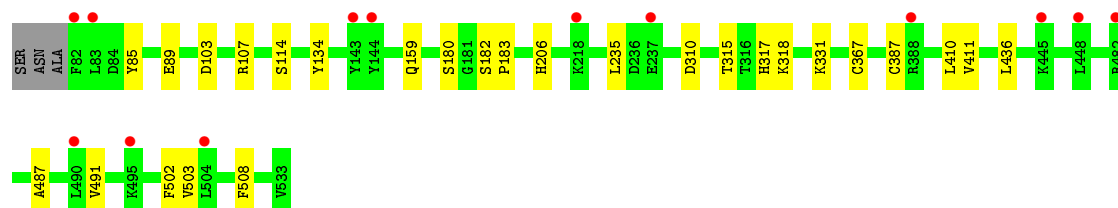
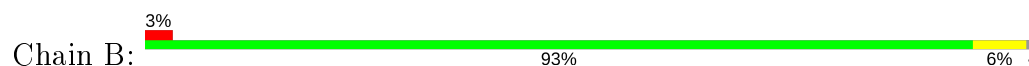
3 Residue-property plots [i](#)

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.

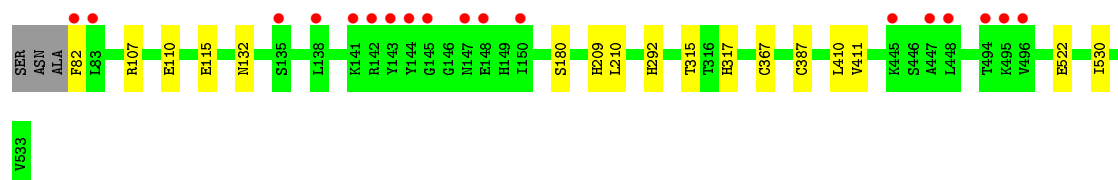
- Molecule 1: Serine hydroxymethyltransferase



- Molecule 1: Serine hydroxymethyltransferase



- Molecule 1: Serine hydroxymethyltransferase



- Molecule 2: Serine hydroxymethyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	151.66 Å 201.56 Å 64.78 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	83.94 – 1.74 83.94 – 1.74	Depositor EDS
% Data completeness (in resolution range)	99.6 (83.94-1.74) 99.6 (83.94-1.74)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 1.74 Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.175 , 0.219 0.184 , 0.218	Depositor DCC
R_{free} test set	1017 reflections (0.50%)	wwPDB-VP
Wilson B-factor (Å ²)	26.1	Xtriage
Anisotropy	0.140	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 39.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	15607	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 41.95 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.1892e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: FMT, LLP, ACT

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	0.52	0/3539	0.64	0/4786
1	B	0.51	0/3540	0.63	0/4786
1	D	0.53	0/3577	0.65	0/4836
2	C	0.52	0/3742	0.62	0/5058
All	All	0.52	0/14398	0.63	0/19466

There are no bond length outliers.

There are no bond angle outliers.

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	3492	0	3478	10	0
1	B	3493	0	3478	15	0
1	D	3527	0	3503	14	0
2	C	3662	0	3641	28	0
3	A	4	0	3	1	0
3	B	4	0	3	0	0
3	C	4	0	3	0	0
4	B	3	0	1	1	0
4	C	3	0	1	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	6	0	2	0	0
5	A	318	0	0	1	0
5	B	313	0	0	1	0
5	C	365	0	0	2	0
5	D	413	0	0	3	0
All	All	15607	0	14113	60	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 2.

All (60) 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:C:138[B]:LEU:HD21	1:D:210:LEU:HB2	1.46	0.98
2:C:138[B]:LEU:HD21	1:D:210:LEU:CB	2.02	0.89
2:C:141[B]:LYS:CD	2:C:142[B]:ARG:H	1.89	0.86
2:C:141[B]:LYS:HD3	2:C:142[B]:ARG:H	1.42	0.84
2:C:141[B]:LYS:HD3	2:C:142[B]:ARG:N	1.94	0.81
2:C:138[B]:LEU:HD13	1:D:209:HIS:CE1	2.18	0.79
2:C:141[B]:LYS:N	2:C:141[B]:LYS:HD2	2.01	0.74
2:C:141[B]:LYS:H	2:C:141[B]:LYS:HD2	1.59	0.66
2:C:138[B]:LEU:HD13	1:D:209:HIS:NE2	2.09	0.66
1:A:487:ALA:O	1:A:491:VAL:HG23	1.97	0.64
2:C:530:ILE:HG21	5:C:729:HOH:O	2.01	0.61
2:C:141[B]:LYS:HD2	2:C:142[B]:ARG:H	1.64	0.61
1:B:103:ASP:OD2	1:B:107:ARG:NH2	2.33	0.61
2:C:141[B]:LYS:CD	2:C:142[B]:ARG:N	2.57	0.60
2:C:138[B]:LEU:HD22	1:D:209:HIS:CD2	2.42	0.55
1:A:491:VAL:HG22	1:A:502:PHE:CD2	2.42	0.55
1:B:487:ALA:O	1:B:491:VAL:HG23	2.07	0.54
1:B:491:VAL:HG22	1:B:502:PHE:CD2	2.44	0.53
2:C:317:HIS:HB2	5:C:1004:HOH:O	2.07	0.53
2:C:315:THR:HB	2:C:317:HIS:CE1	2.43	0.53
2:C:139[B]:PRO:O	2:C:349:PHE:CZ	2.64	0.51
1:B:310:ASP:HA	1:B:331:LYS:HD2	1.92	0.51
2:C:138[B]:LEU:HD21	1:D:210:LEU:HB3	1.90	0.50
2:C:140[B]:GLY:O	2:C:144[B]:TYR:CD2	2.65	0.50
1:A:318:LLP:C4'	3:A:601:ACT:H3	2.43	0.48
2:C:134[B]:TYR:OH	2:C:137[B]:GLY:HA3	2.14	0.47
2:C:349:PHE:CD1	2:C:350:PRO:HA	2.50	0.47
1:A:435:THR:HG22	5:A:704:HOH:O	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:491:VAL:HG22	1:B:502:PHE:CG	2.50	0.46
2:C:138[B]:LEU:HG	2:C:139[B]:PRO:HA	1.97	0.46
1:A:491:VAL:HG22	1:A:502:PHE:CG	2.50	0.46
2:C:141[B]:LYS:CD	2:C:141[B]:LYS:N	2.70	0.46
1:D:132:ASN:ND2	5:D:706:HOH:O	2.48	0.46
1:D:315:THR:HB	1:D:317[A]:HIS:CE1	2.51	0.45
1:B:387:CYS:SG	1:B:411:VAL:HG13	2.57	0.45
1:B:206:HIS:HB3	1:B:235:LEU:HG	1.97	0.45
1:B:182:SER:HB2	1:B:183:PRO:HD3	1.97	0.45
1:B:85:TYR:HB3	1:B:89:GLU:HB2	1.98	0.45
1:D:530:ILE:HG21	5:D:773:HOH:O	2.17	0.45
1:D:110[A]:GLU:CG	1:D:115[A]:GLU:OE2	2.65	0.44
1:B:503:VAL:HA	1:B:508:PHE:CD1	2.52	0.44
1:A:301:VAL:HG11	1:A:384:VAL:HG21	2.00	0.44
2:C:139[B]:PRO:O	2:C:349:PHE:HZ	2.01	0.44
2:C:138[B]:LEU:CD1	1:D:209:HIS:CE1	2.95	0.44
1:B:159:GLN:NE2	5:B:704:HOH:O	2.47	0.43
1:B:317[B]:HIS:N	4:B:601:FMT:O2	2.51	0.43
1:B:410:LEU:C	1:B:410:LEU:HD12	2.39	0.43
1:D:107:ARG:O	1:D:522:GLU:HG2	2.19	0.43
2:C:142[B]:ARG:HE	2:C:142[B]:ARG:HB2	1.72	0.43
1:A:410:LEU:C	1:A:410:LEU:HD12	2.40	0.42
2:C:301:VAL:HG11	2:C:384:VAL:HG21	2.02	0.42
2:C:140[B]:GLY:O	2:C:144[B]:TYR:HD2	2.03	0.42
1:A:182:SER:HB2	1:A:183:PRO:HD3	2.02	0.41
1:A:457:SER:N	1:A:458:PRO:CD	2.82	0.41
2:C:138[B]:LEU:CD2	5:D:1019:HOH:O	2.67	0.41
1:B:114:SER:HB2	1:B:318:LLP:HE3	2.03	0.41
1:B:315:THR:HB	1:B:317[B]:HIS:CE1	2.54	0.41
1:A:206:HIS:HB3	1:A:235:LEU:HG	2.02	0.41
1:D:387:CYS:SG	1:D:411:VAL:HG13	2.61	0.41
1:D:410:LEU:C	1:D:410:LEU:HD12	2.42	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	450/455 (99%)	439 (98%)	11 (2%)	0	100	100
1	B	450/455 (99%)	437 (97%)	13 (3%)	0	100	100
1	D	455/455 (100%)	444 (98%)	11 (2%)	0	100	100
2	C	474/455 (104%)	461 (97%)	13 (3%)	0	100	100
All	All	1829/1820 (100%)	1781 (97%)	48 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	370/371 (100%)	365 (99%)	5 (1%)	67	50
1	B	370/371 (100%)	366 (99%)	4 (1%)	73	59
1	D	375/371 (101%)	371 (99%)	4 (1%)	73	59
2	C	390/372 (105%)	384 (98%)	6 (2%)	65	47
All	All	1505/1485 (101%)	1486 (99%)	19 (1%)	71	52

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	134	TYR
1	A	180	SER
1	A	367	CYS
1	A	411	VAL
1	A	436	LEU
1	B	134	TYR
1	B	180	SER
1	B	367	CYS

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Mol	Chain	Res	Type
1	B	436	LEU
2	C	134[A]	TYR
2	C	134[B]	TYR
2	C	367	CYS
2	C	441	VAL
2	C	507	GLU
2	C	519	ARG
1	D	82	PHE
1	D	180	SER
1	D	292	HIS
1	D	367	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	99	ASN
1	A	292	HIS
1	B	99	ASN
1	B	159	GLN
1	D	358	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 non-standard protein/DNA/RNA residues 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$
1	LLP	A	318	1	23,24,25	1.24	3 (13%)	25,32,34	2.02	10 (40%)
1	LLP	B	318	1	23,24,25	1.29	3 (13%)	25,32,34	1.79	6 (24%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	LLP	D	318	1	23,24,25	1.37	3 (13%)	25,32,34	1.95	6 (24%)

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
1	LLP	A	318	1	-	5/16/17/19	0/1/1/1
1	LLP	B	318	1	-	7/16/17/19	0/1/1/1
1	LLP	D	318	1	-	7/16/17/19	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	318	LLP	C4-C4'	3.36	1.53	1.46
1	B	318	LLP	CB-CA	2.87	1.57	1.53
1	D	318	LLP	C4-C5	-2.85	1.38	1.42
1	A	318	LLP	C4-C4'	2.81	1.52	1.46
1	B	318	LLP	C4-C4'	2.81	1.52	1.46
1	A	318	LLP	C4-C5	-2.31	1.39	1.42
1	D	318	LLP	C2'-C2	2.31	1.54	1.50
1	B	318	LLP	C4-C5	-2.27	1.39	1.42
1	A	318	LLP	CB-CA	2.08	1.56	1.53

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	318	LLP	OP4-C5'-C5	5.12	119.11	109.35
1	D	318	LLP	C5'-C5-C6	-4.47	112.02	119.37
1	D	318	LLP	OP4-C5'-C5	4.24	117.43	109.35
1	B	318	LLP	OP4-C5'-C5	4.22	117.39	109.35
1	A	318	LLP	CE-NZ-C4'	4.05	131.35	118.90
1	B	318	LLP	CE-NZ-C4'	4.03	131.27	118.90
1	D	318	LLP	CE-NZ-C4'	3.84	130.68	118.90
1	A	318	LLP	C5'-C5-C6	-3.22	114.08	119.37
1	A	318	LLP	C4-C4'-NZ	-3.08	110.19	124.31
1	B	318	LLP	C4-C4'-NZ	-3.03	110.40	124.31
1	D	318	LLP	C4-C4'-NZ	-2.90	110.98	124.31
1	B	318	LLP	C5'-C5-C6	-2.72	114.90	119.37
1	D	318	LLP	CG-CD-CE	-2.51	104.82	113.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	318	LLP	OP4-P-OP1	-2.25	100.15	106.47
1	A	318	LLP	C6-N1-C2	2.24	123.32	119.17
1	D	318	LLP	C3-C4-C4'	-2.24	116.23	120.41
1	B	318	LLP	C5-C6-N1	-2.24	120.09	123.82
1	A	318	LLP	C2'-C2-C3	2.19	123.59	120.89
1	A	318	LLP	C3-C2-N1	-2.17	117.97	120.77
1	A	318	LLP	C3-C4-C5	2.09	119.86	118.26
1	B	318	LLP	C3-C4-C5	2.07	119.85	118.26
1	A	318	LLP	C5-C6-N1	-2.03	120.44	123.82

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	318	LLP	O-C-CA-CB
1	B	318	LLP	C6-C5-C5'-OP4
1	B	318	LLP	O-C-CA-CB
1	D	318	LLP	C6-C5-C5'-OP4
1	D	318	LLP	O-C-CA-CB
1	D	318	LLP	CG-CD-CE-NZ
1	B	318	LLP	CG-CD-CE-NZ
1	A	318	LLP	CG-CD-CE-NZ
1	D	318	LLP	C5-C4-C4'-NZ
1	D	318	LLP	C3-C4-C4'-NZ
1	A	318	LLP	C6-C5-C5'-OP4
1	B	318	LLP	CD-CE-NZ-C4'
1	B	318	LLP	N-CA-CB-CG
1	D	318	LLP	N-CA-CB-CG
1	A	318	LLP	CD-CE-NZ-C4'
1	D	318	LLP	CD-CE-NZ-C4'
1	A	318	LLP	C3-C4-C4'-NZ
1	B	318	LLP	C3-C4-C4'-NZ
1	B	318	LLP	C5-C4-C4'-NZ

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	318	LLP	1	0
1	B	318	LLP	1	0

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

7 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
4	FMT	D	602	-	0,2,2	0.00	-	0,1,1	0.00	-
3	ACT	C	602	-	1,3,3	2.54	1 (100%)	0,3,3	0.00	-
4	FMT	C	601	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	D	601	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	B	601	-	0,2,2	0.00	-	0,1,1	0.00	-
3	ACT	A	601	-	1,3,3	2.61	1 (100%)	0,3,3	0.00	-
3	ACT	B	602	-	1,3,3	2.52	1 (100%)	0,3,3	0.00	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	ACT	CH3-C	2.61	1.52	1.48
3	C	602	ACT	CH3-C	2.54	1.52	1.48
3	B	602	ACT	CH3-C	2.52	1.51	1.48

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	601	FMT	1	0
3	A	601	ACT	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	451/455 (99%)	0.13	14 (3%) 49 55	19, 34, 57, 85	0
1	B	451/455 (99%)	0.14	13 (2%) 51 57	19, 34, 58, 82	0
1	D	451/455 (99%)	0.03	18 (3%) 38 43	18, 27, 51, 84	0
2	C	452/455 (99%)	0.03	18 (3%) 38 43	19, 28, 45, 68	0
All	All	1805/1820 (99%)	0.08	63 (3%) 44 49	18, 30, 55, 85	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	82	PHE	11.3
2	C	82	PHE	9.7
1	A	82	PHE	8.8
1	D	82	PHE	8.8
1	B	448	LEU	7.9
1	A	448	LEU	7.0
1	A	445	LYS	5.7
1	D	141	LYS	4.3
2	C	532	GLY	4.3
2	C	143[A]	TYR	4.3
1	A	490	LEU	4.1
1	B	143	TYR	4.1
1	B	218	LYS	4.1
2	C	533	VAL	4.0
2	C	146[A]	GLY	4.0
1	B	445	LYS	3.9
2	C	144[A]	TYR	3.9
1	D	142	ARG	3.8
1	D	494	THR	3.7
1	A	446	SER	3.6
2	C	134[A]	TYR	3.5

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Mol	Chain	Res	Type	RSRZ
1	D	147	ASN	3.4
1	A	143	TYR	3.4
1	D	145	GLY	3.4
2	C	138[A]	LEU	3.2
2	C	507	GLU	3.2
2	C	83	LEU	3.0
2	C	139[A]	PRO	3.0
1	A	494	THR	2.9
1	D	143	TYR	2.9
1	D	83	LEU	2.8
1	D	148	GLU	2.8
1	A	502	PHE	2.7
1	D	448	LEU	2.6
1	B	482	ARG	2.6
1	D	150	ILE	2.5
1	D	135	SER	2.5
1	A	218	LYS	2.5
1	B	83	LEU	2.4
2	C	145[A]	GLY	2.4
1	A	509	PRO	2.3
1	B	237	GLU	2.3
2	C	147[A]	ASN	2.3
2	C	218	LYS	2.3
2	C	150[A]	ILE	2.3
1	A	491	VAL	2.2
1	A	506	PRO	2.2
2	C	135[A]	SER	2.2
1	D	445	LYS	2.2
1	D	496	VAL	2.2
1	B	490	LEU	2.2
1	D	138	LEU	2.2
1	D	495	LYS	2.2
2	C	506	PRO	2.1
1	D	144	TYR	2.1
1	A	482	ARG	2.1
1	D	447	ALA	2.1
1	A	441	VAL	2.1
1	B	388	ARG	2.1
2	C	149[A]	HIS	2.1
1	B	504	LEU	2.1
1	B	495	LYS	2.0
1	B	144	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	LLP	D	318	24/25	0.85	0.17	18,22,26,27	15
1	LLP	B	318	24/25	0.91	0.12	30,32,34,36	0
1	LLP	A	318	24/25	0.96	0.08	26,29,30,31	0

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. 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(Å ²)	Q<0.9
3	ACT	C	602	4/4	0.69	0.20	41,45,47,50	0
3	ACT	A	601	4/4	0.79	0.16	43,48,49,54	0
4	FMT	D	601	3/3	0.81	0.27	47,47,48,49	0
3	ACT	B	602	4/4	0.86	0.18	56,57,58,60	0
4	FMT	C	601	3/3	0.91	0.22	37,37,37,37	0
4	FMT	D	602	3/3	0.92	0.23	28,28,28,29	3
4	FMT	B	601	3/3	0.92	0.17	25,25,25,26	3

6.5 Other polymers [i](#)

There are no such residues in this entry.