



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:06 PM GMT

PDB ID : 5AGV
Title : The sliding clamp of Mycobacterium tuberculosis in complex with a natural product.
Authors : Lukat, P.; Kling, A.; Heinz, D.W.; Mueller, R.
Deposited on : 2015-02-03
Resolution : 1.93 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

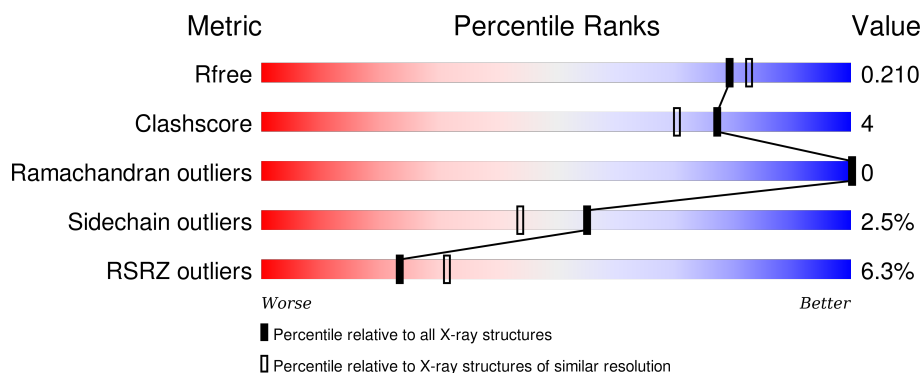
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.93 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	2910 (1.96-1.92)
Clashscore	102246	3095 (1.96-1.92)
Ramachandran outliers	100387	3062 (1.96-1.92)
Sidechain outliers	100360	3062 (1.96-1.92)
RSRZ outliers	91569	2915 (1.96-1.92)

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	406	<div> <div>6%</div> <div>88%</div> <div>8%</div> <div>.</div> </div>
1	B	406	<div> <div>6%</div> <div>89%</div> <div>7%</div> <div>..</div> </div>
2	C	11	<div> <div>45%</div> <div>36%</div> <div>18%</div> </div>
2	D	11	<div> <div>82%</div> <div>9%</div> <div>9%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	CA	B	1407	-	-	-	X
4	NA	A	1409	-	-	-	X
4	NA	B	1410	-	-	-	X
5	BU3	A	1413	-	-	-	X
5	BU3	B	1413	-	-	-	X
5	BU3	B	1414	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7271 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 DNA POLYMERASE III SUBUNIT BETA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	390	Total	C	N	O	S	0	27	0
			3038	1926	508	598	6			
1	B	393	Total	C	N	O	S	0	23	0
			3040	1924	520	590	6			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	EXPRESSION TAG	UNP I6XU56
A	-2	GLY	-	EXPRESSION TAG	UNP I6XU56
A	-1	GLY	-	EXPRESSION TAG	UNP I6XU56
A	0	ARG	-	EXPRESSION TAG	UNP I6XU56
B	-3	GLY	-	EXPRESSION TAG	UNP I6XU56
B	-2	GLY	-	EXPRESSION TAG	UNP I6XU56
B	-1	GLY	-	EXPRESSION TAG	UNP I6XU56
B	0	ARG	-	EXPRESSION TAG	UNP I6XU56

- Molecule 2 is a protein called CYCLOHEXYL GRISELIMYCIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	11	Total	C	N	O	0	0	0
			85	63	10	12			
2	D	11	Total	C	N	O	0	0	0
			85	63	10	12			

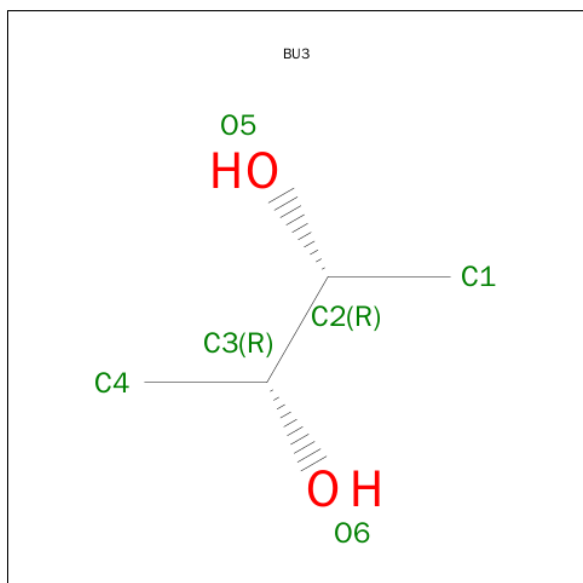
- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	5	Total	Ca	0	0
			5	5		
3	A	1	Total	Ca	0	0
			1	1		

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	5	Total	Na	0	0
			5	5		
4	A	7	Total	Na	0	0
			7	7		

- Molecule 5 is (R,R)-2,3-BUTANEDIOL (three-letter code: BU3) (formula: C₄H₁₀O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	4	2		
5	A	1	Total	C	O	0	0
			6	4	2		
5	B	1	Total	C	O	0	0
			6	4	2		
5	B	1	Total	C	O	0	0
			6	4	2		
5	B	1	Total	C	O	0	0
			6	4	2		

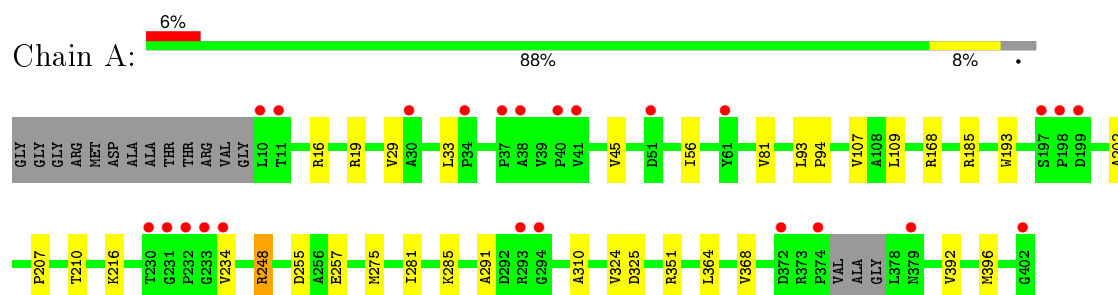
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	475	Total	O	0	0
			475	475		
6	B	500	Total	O	0	0
			500	500		

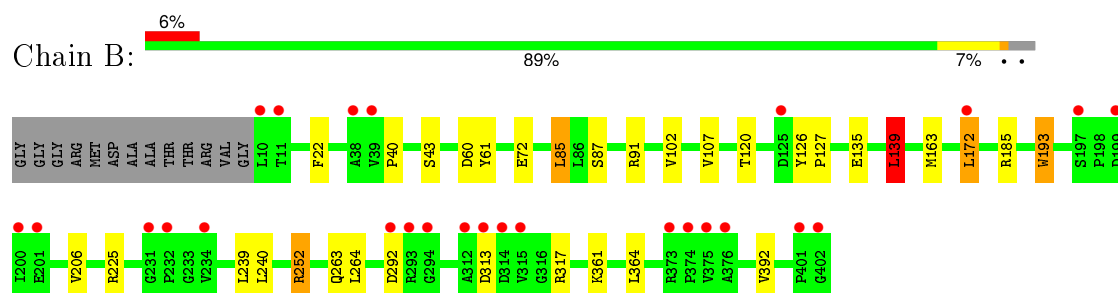
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

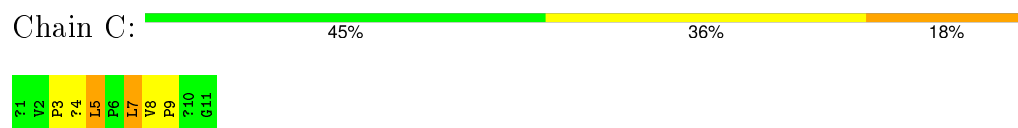
- Molecule 1: DNA POLYMERASE III SUBUNIT BETA



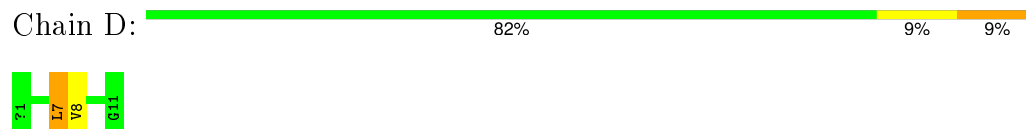
- Molecule 1: DNA POLYMERASE III SUBUNIT BETA



- Molecule 2: CYCLOHEXYL GRISELIMYCIN



- Molecule 2: CYCLOHEXYL GRISELIMYCIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	72.87Å 89.76Å 81.93Å 90.00° 95.14° 90.00°	Depositor
Resolution (Å)	40.80 – 1.93 40.80 – 1.93	Depositor EDS
% Data completeness (in resolution range)	100.0 (40.80-1.93) 100.0 (40.80-1.93)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 1.94Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.169 , 0.209 0.173 , 0.210	Depositor DCC
R_{free} test set	3846 reflections (5.12%)	DCC
Wilson B-factor (Å ²)	26.4	Xtriage
Anisotropy	0.143	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 49.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 78978 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7271	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.93% 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: ACE, BU3, CA, MLU, NZC, MVA, NA, PH6, MP8

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	0.36	0/3107	0.52	0/4245
1	B	0.37	0/3106	0.55	1/4243 (0.0%)
2	C	0.49	0/14	1.39	0/16
2	D	0.47	0/14	0.84	0/16
All	All	0.37	0/6241	0.54	1/8520 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	139	LEU	CA-CB-CG	7.31	132.11	115.30

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	3038	0	3063	23	0
1	B	3040	0	3077	22	0
2	C	85	0	103	5	0
2	D	85	0	103	1	0
3	A	1	0	0	0	0
3	B	5	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	7	0	0	0	0
4	B	5	0	0	0	0
5	A	12	0	18	1	0
5	B	18	0	27	0	0
6	A	475	0	0	4	1
6	B	500	0	0	3	0
All	All	7271	0	6391	47	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 (47) 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:135:GLU:HB2	1:B:225:ARG:HH11	1.54	0.72
1:A:325[A]:ASP:OD2	6:A:2333:HOH:O	2.12	0.67
1:A:216:LYS:NZ	6:A:2226:HOH:O	2.31	0.64
1:B:139:LEU:HD13	1:B:193:TRP:HB3	1.80	0.64
1:A:16:ARG:NH1	1:B:72:GLU:OE1	2.31	0.63
1:A:396:MET:HB2	2:C:5:LEU:HD23	1.82	0.62
1:A:33:LEU:HD11	1:A:45[A]:VAL:HG23	1.82	0.62
1:B:163:MET:O	6:B:2246:HOH:O	2.16	0.61
1:B:102:VAL:HG22	1:B:107[A]:VAL:HG22	1.82	0.60
1:A:81[A]:VAL:HG11	1:A:107:VAL:HG11	1.83	0.59
1:B:206:VAL:HG21	1:B:240[B]:LEU:HD12	1.83	0.59
1:A:248:ARG:NH1	6:A:2310:HOH:O	2.36	0.57
1:A:281:ILE:O	1:A:285[A]:LYS:HG2	2.06	0.56
1:A:275:MET:HG2	1:A:324[B]:VAL:HG12	1.88	0.56
1:B:313:ASP:OD1	1:B:317[B]:ARG:HD3	2.09	0.53
1:B:135:GLU:OE2	6:B:2205:HOH:O	2.17	0.53
1:A:16:ARG:HH12	1:B:72:GLU:CD	2.13	0.52
1:B:135:GLU:HB2	1:B:225:ARG:HD2	1.92	0.50
1:B:139:LEU:HD13	1:B:193:TRP:CB	2.41	0.50
1:B:107[A]:VAL:HG23	1:B:120:THR:HG22	1.95	0.49
1:B:60:ASP:O	1:B:61:TYR:HB2	2.13	0.49
1:B:364[B]:LEU:HD11	1:B:392[B]:VAL:HG22	1.95	0.48
2:C:7:LEU:HA	2:C:8:MVA:HN1	1.62	0.47
1:A:257:GLU:HB3	2:C:9:PH6:H51	1.97	0.47
1:A:19:ARG:HD3	1:A:94:PRO:O	2.15	0.46
1:A:29:VAL:HG11	1:A:56:ILE:HG22	1.97	0.46
1:A:168:ARG:NH1	6:A:2254:HOH:O	2.35	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:351:ARG:HB2	1:A:368[A]:VAL:HG23	1.99	0.44
1:B:40:PRO:O	1:B:43[B]:SER:OG	2.24	0.43
1:B:126:TYR:HA	1:B:127:PRO:HD3	1.86	0.43
1:A:93:LEU:HD11	1:A:109:LEU:HD21	2.00	0.43
1:A:392[B]:VAL:HG23	5:A:1412:BU3:H11	1.99	0.43
1:B:85:LEU:HA	1:B:85:LEU:HD12	1.90	0.43
1:A:364:LEU:HD21	1:A:392[A]:VAL:HG12	1.99	0.42
1:A:33:LEU:HD11	1:A:45[B]:VAL:HG13	2.01	0.42
1:B:263:GLN:HG3	1:B:264:LEU:HD12	2.02	0.42
2:C:3:MP8:HA	2:C:4:NZC:H40	1.75	0.42
1:A:202:ALA:HB2	1:A:234:VAL:HG13	2.01	0.42
1:B:87:SER:OG	1:B:91[A]:ARG:NH2	2.52	0.41
1:B:172:LEU:HB2	6:B:2255:HOH:O	2.20	0.41
1:A:207:PRO:HB2	1:A:210[B]:THR:HG22	2.03	0.41
2:D:7:LEU:HA	2:D:8:MVA:HN1	1.68	0.41
1:B:361:LYS:HA	1:B:361:LYS:HD3	1.88	0.41
2:C:8:MVA:HA	2:C:9:PH6:HA	1.94	0.41
1:B:239[A]:LEU:CD2	1:B:252:ARG:HG2	2.51	0.41
1:A:351:ARG:CB	1:A:368[A]:VAL:HG23	2.51	0.40
1:A:291:ALA:HB2	1:A:310:ALA:HB1	2.03	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)
6:A:2063:HOH:O	6:A:2396:HOH:O[1_556]	2.18	0.02

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	413/406 (102%)	405 (98%)	8 (2%)	0	100 100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	414/406 (102%)	408 (99%)	6 (1%)	0	100	100
2	C	2/11 (18%)	2 (100%)	0	0	100	100
2	D	2/11 (18%)	2 (100%)	0	0	100	100
All	All	831/834 (100%)	817 (98%)	14 (2%)	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	330/317 (104%)	325 (98%)	5 (2%)	72	66
1	B	326/317 (103%)	318 (98%)	8 (2%)	55	44
2	C	2/2 (100%)	0	2 (100%)	0	0
2	D	2/2 (100%)	1 (50%)	1 (50%)	0	0
All	All	660/638 (103%)	644 (98%)	16 (2%)	55	46

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	185	ARG
1	A	193	TRP
1	A	248	ARG
1	A	255[A]	ASP
1	A	255[B]	ASP
1	B	22	PHE
1	B	85	LEU
1	B	139	LEU
1	B	172	LEU
1	B	185	ARG
1	B	193	TRP
1	B	252	ARG
1	B	292	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	C	5	LEU
2	C	7	LEU
2	D	7	LEU

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

14 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$
2	MLU	C	10	2	7,8,9	0.43	0	4,9,11	1.05	1 (25%)
2	MVA	C	2	2	6,7,8	0.62	0	6,8,10	1.41	1 (16%)
2	MP8	C	3	2	7,8,9	0.45	0	5,10,12	1.15	1 (20%)
2	NZC	C	4	2	6,7,8	0.41	0	6,8,10	1.46	1 (16%)
2	MP8	C	6	2	7,8,9	0.33	0	5,10,12	1.30	0
2	MVA	C	8	2	6,7,8	0.66	0	6,8,10	1.62	1 (16%)
2	PH6	C	9	2	12,14,15	0.42	0	10,18,20	1.14	1 (10%)
2	MLU	D	10	2	7,8,9	0.40	0	4,9,11	0.89	0
2	MVA	D	2	2	6,7,8	0.52	0	6,8,10	1.22	1 (16%)
2	MP8	D	3	2	7,8,9	0.37	0	5,10,12	1.14	1 (20%)
2	NZC	D	4	2	6,7,8	0.51	0	6,8,10	1.57	2 (33%)
2	MP8	D	6	2	7,8,9	0.32	0	5,10,12	1.01	0
2	MVA	D	8	2	6,7,8	0.51	0	6,8,10	1.60	2 (33%)
2	PH6	D	9	2	12,14,15	0.39	0	10,18,20	1.18	1 (10%)

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	MLU	C	10	2	-	0/4/8/10	0/0/0/0
2	MVA	C	2	2	-	0/5/8/10	0/0/0/0
2	MP8	C	3	2	-	0/0/11/13	0/1/1/1
2	NZC	C	4	2	-	0/5/8/10	0/0/0/0
2	MP8	C	6	2	-	0/0/11/13	0/1/1/1
2	MVA	C	8	2	-	0/5/8/10	0/0/0/0
2	PH6	C	9	2	-	0/4/23/25	0/2/2/2
2	MLU	D	10	2	-	0/4/8/10	0/0/0/0
2	MVA	D	2	2	-	0/5/8/10	0/0/0/0
2	MP8	D	3	2	-	0/0/11/13	0/1/1/1
2	NZC	D	4	2	-	0/5/8/10	0/0/0/0
2	MP8	D	6	2	-	0/0/11/13	0/1/1/1
2	MVA	D	8	2	-	0/5/8/10	0/0/0/0
2	PH6	D	9	2	-	0/4/23/25	0/2/2/2

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	8	MVA	CB-CA-N	-3.35	105.78	111.57
2	D	9	PH6	O-C-CA	-2.84	117.94	125.44
2	C	2	MVA	CB-CA-N	-2.73	106.84	111.57
2	C	9	PH6	O-C-CA	-2.49	118.86	125.44
2	C	3	MP8	O-C-CA	-2.27	119.44	125.44
2	D	2	MVA	CB-CA-N	-2.26	107.65	111.57
2	D	8	MVA	CB-CA-N	-2.23	107.70	111.57
2	D	8	MVA	O-C-CA	-2.12	119.22	125.74
2	D	3	MP8	O-C-CA	-2.09	119.92	125.44
2	C	10	MLU	O-C-CA	-2.07	119.96	125.44
2	D	4	NZC	O-C-CA	-2.06	119.43	125.74
2	D	4	NZC	C40-N-CA	2.80	122.29	113.65
2	C	4	NZC	C40-N-CA	2.81	122.30	113.65

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	3	MP8	1	0
2	C	4	NZC	1	0
2	C	8	MVA	2	0
2	C	9	PH6	2	0
2	D	8	MVA	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 23 ligands modelled in this entry, 18 are monoatomic - leaving 5 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	$\# Z > 2$
5	BU3	A	1412	3	4,5,5	0.67	0	6,6,6	0.32	0
5	BU3	A	1413	-	4,5,5	0.61	0	6,6,6	0.30	0
5	BU3	B	1413	-	4,5,5	0.77	0	6,6,6	0.43	0
5	BU3	B	1414	3	4,5,5	0.52	0	6,6,6	0.31	0
5	BU3	B	1415	3	4,5,5	0.68	0	6,6,6	0.32	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
5	BU3	A	1412	3	-	0/4/4/4	0/0/0/0
5	BU3	A	1413	-	-	0/4/4/4	0/0/0/0
5	BU3	B	1413	-	-	0/4/4/4	0/0/0/0
5	BU3	B	1414	3	-	0/4/4/4	0/0/0/0
5	BU3	B	1415	3	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1412	BU3	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	390/406 (96%)	0.01	24 (6%) 24 32	17, 30, 64, 106	0
1	B	393/406 (96%)	0.02	26 (6%) 22 29	15, 27, 69, 84	0
2	C	3/11 (27%)	-0.37	0 100 100	22, 22, 25, 25	0
2	D	3/11 (27%)	-0.37	0 100 100	18, 18, 20, 21	0
All	All	789/834 (94%)	0.01	50 (6%) 23 31	15, 29, 67, 106	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	10	LEU	6.5
1	B	10	LEU	6.1
1	A	374	PRO	5.5
1	A	402	GLY	5.5
1	B	200	ILE	5.2
1	A	30	ALA	4.5
1	B	375	VAL	4.5
1	B	292	ASP	4.2
1	B	402	GLY	4.2
1	A	198	PRO	4.1
1	B	294	GLY	3.9
1	B	199	ASP	3.9
1	A	34	PRO	3.9
1	B	373	ARG	3.8
1	A	11	THR	3.6
1	B	231	GLY	3.6
1	A	232	PRO	3.3
1	B	197	SER	3.3
1	A	40	PRO	3.2
1	A	234	VAL	3.1
1	B	232	PRO	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	374	PRO	2.9
1	A	294	GLY	2.8
1	A	199	ASP	2.8
1	A	38	ALA	2.8
1	A	372	ASP	2.8
1	B	315	VAL	2.7
1	B	376	ALA	2.7
1	A	61[A]	TYR	2.6
1	B	11	THR	2.6
1	A	41	VAL	2.6
1	A	379	ASN	2.5
1	B	201	GLU	2.5
1	B	38	ALA	2.5
1	A	231	GLY	2.5
1	A	293	ARG	2.4
1	B	234	VAL	2.4
1	B	401	PRO	2.4
1	A	197	SER	2.4
1	B	312	ALA	2.2
1	A	233	GLY	2.2
1	B	125	ASP	2.1
1	B	293	ARG	2.1
1	B	39	VAL	2.1
1	B	313	ASP	2.1
1	A	37	PRO	2.1
1	A	230	THR	2.1
1	A	51	ASP	2.1
1	B	172	LEU	2.0
1	B	314	ASP	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	MLU	D	10	9/10	0.98	0.08	-	19,27,33,35	0
2	MVA	C	2	8/9	0.97	0.09	-	22,25,29,29	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	MP8	D	3	8/9	0.99	0.06	-	17,19,20,21	0
2	NZC	D	4	8/9	0.97	0.08	-	17,20,24,24	0
2	PH6	D	9	13/14	0.95	0.09	-	22,28,44,45	0
2	MVA	D	8	8/9	0.97	0.09	-	22,37,38,42	0
2	MVA	C	8	8/9	0.95	0.08	-	29,35,38,40	0
2	PH6	C	9	13/14	0.96	0.08	-	25,28,36,36	0
2	MP8	C	3	8/9	0.97	0.11	-	20,22,26,27	0
2	NZC	C	4	8/9	0.97	0.07	-	20,23,27,28	0
2	MLU	C	10	9/10	0.98	0.08	-	26,32,36,36	0
2	MP8	D	6	8/9	0.97	0.07	-	18,21,24,25	0
2	MVA	D	2	8/9	0.97	0.07	-	22,25,28,30	0
2	MP8	C	6	8/9	0.97	0.06	-	21,25,28,28	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. 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(\AA^2)	Q<0.9
4	NA	A	1409	1/1	0.98	0.23	8.66	50,50,50,50	0
4	NA	B	1410	1/1	0.94	0.27	7.48	61,61,61,61	0
3	CA	B	1407	1/1	0.83	0.13	4.48	70,70,70,70	0
5	BU3	B	1414	6/6	0.93	0.16	2.28	41,43,44,46	0
5	BU3	B	1413	6/6	0.95	0.14	2.13	43,47,48,50	0
5	BU3	A	1413	6/6	0.81	0.16	2.04	53,57,59,61	0
4	NA	A	1408	1/1	0.95	0.12	0.32	53,53,53,53	0
4	NA	B	1411	1/1	0.90	0.16	0.05	68,68,68,68	0
5	BU3	A	1412	6/6	0.95	0.09	-0.50	28,32,33,36	0
3	CA	B	1405	1/1	0.99	0.09	-0.76	24,24,24,24	0
4	NA	A	1405	1/1	0.97	0.08	-1.05	43,43,43,43	0
5	BU3	B	1415	6/6	0.98	0.07	-2.44	29,31,34,36	0
4	NA	A	1406	1/1	0.99	0.04	-4.16	38,38,38,38	0
4	NA	A	1407	1/1	0.90	0.15	-	62,62,62,62	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	CA	B	1404	1/1	1.00	0.04	-	30,30,30,30	0
4	NA	B	1409	1/1	0.85	0.16	-	67,67,67,67	0
3	CA	A	1404	1/1	0.98	0.04	-	40,40,40,40	0
4	NA	A	1411	1/1	0.97	0.06	-	46,46,46,46	0
3	CA	B	1406	1/1	0.99	0.13	-	46,46,46,46	0
3	CA	B	1403	1/1	0.95	0.04	-	50,50,50,50	0
4	NA	B	1408	1/1	0.79	0.16	-	69,69,69,69	0
4	NA	B	1412	1/1	0.80	0.26	-	86,86,86,86	0
4	NA	A	1410	1/1	0.59	0.58	-	111,111,111,111	0

6.5 Other polymers [i](#)

There are no such residues in this entry.