



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:59 AM GMT

PDB ID : 2Z2O  
Title : Crystal Structure of apo virginiamycin B lyase from Staphylococcus aureus  
Authors : Korczynska, M.; Berghuis, A.M.  
Deposited on : 2007-05-25  
Resolution : 1.90 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](mailto: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.

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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

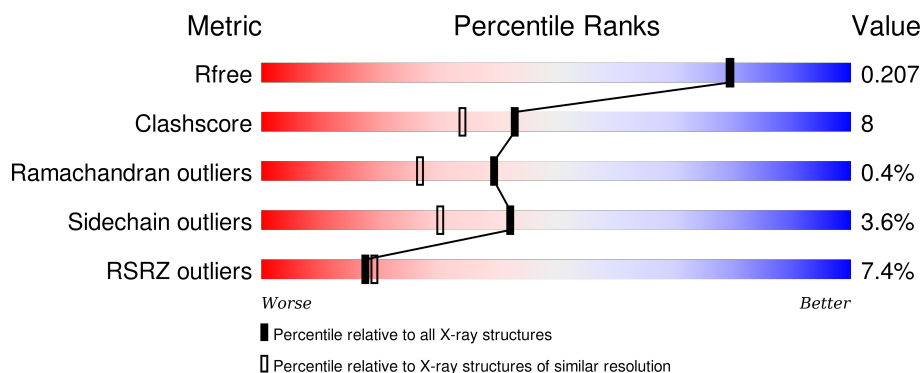
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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  $\geq 3$ , 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	299	<div> <div>7%</div> <div>82%15%.</div> </div>
1	B	299	<div> <div>6%</div> <div>78%18%..</div> </div>
1	C	299	<div> <div>8%</div> <div>81%17%.</div> </div>
1	D	299	<div> <div>8%</div> <div>84%13%..</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
2	CL	D	301	-	-	X	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10412 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 virginiamycin B lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	298	Total	C	N	O	S	0	8	0
			2346	1495	386	454	11			
1	B	293	Total	C	N	O	S	0	8	0
			2311	1471	381	449	10			
1	C	299	Total	C	N	O	S	0	5	0
			2344	1488	385	460	11			
1	D	295	Total	C	N	O	S	1	6	0
			2323	1479	383	449	12			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	211	THR	PRO	SEE REMARK 999	UNP Q53744
A	212	SER	LEU	SEE REMARK 999	UNP Q53744
B	211	THR	PRO	SEE REMARK 999	UNP Q53744
B	212	SER	LEU	SEE REMARK 999	UNP Q53744
C	211	THR	PRO	SEE REMARK 999	UNP Q53744
C	212	SER	LEU	SEE REMARK 999	UNP Q53744
D	211	THR	PRO	SEE REMARK 999	UNP Q53744
D	212	SER	LEU	SEE REMARK 999	UNP Q53744

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cl	0	0
			1	1		
2	A	1	Total	Cl	0	0
			1	1		
2	D	2	Total	Cl	0	0
			2	2		
2	C	1	Total	Cl	0	0
			1	1		

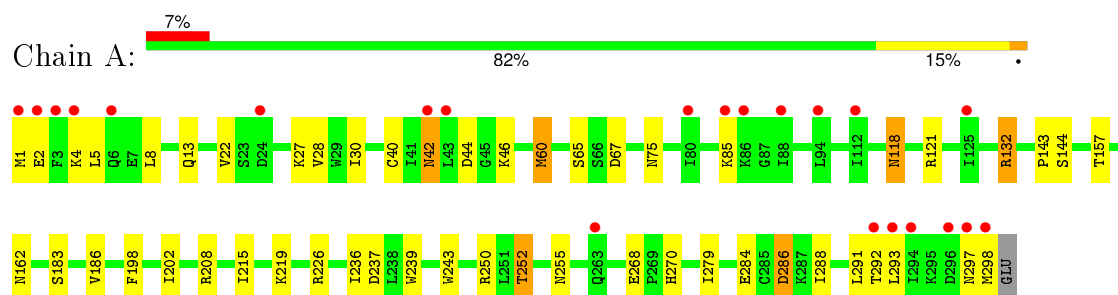
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	327	Total 327	O 327	0	0
3	B	290	Total 290	O 290	0	0
3	C	232	Total 232	O 232	0	0
3	D	234	Total 234	O 234	0	0

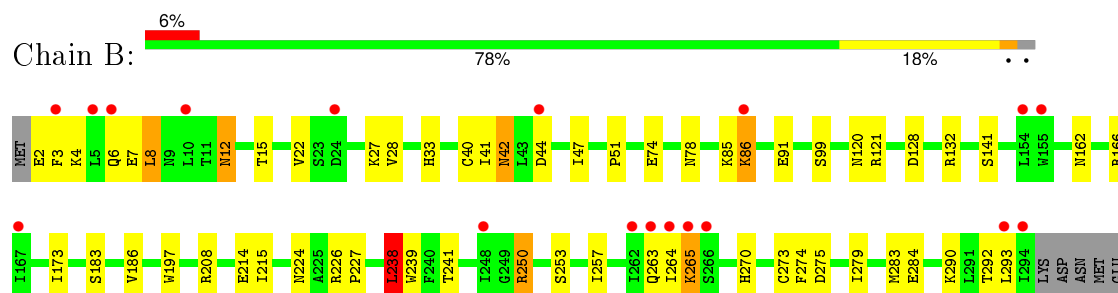
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

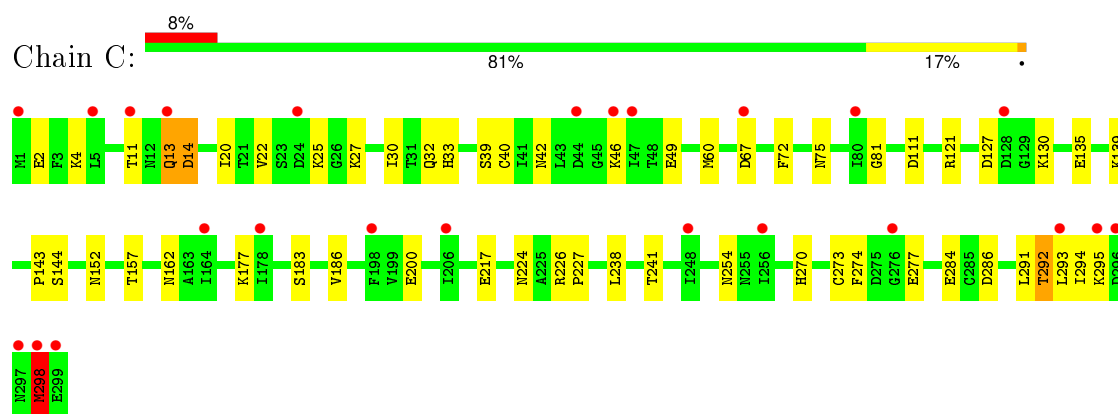
- Molecule 1: virginiamycin B lyase



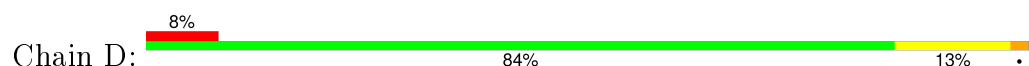
- Molecule 1: virginiamycin B lyase

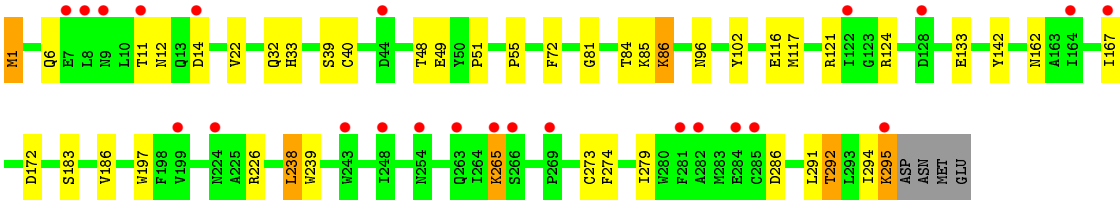


- Molecule 1: virginiamycin B lyase



- Molecule 1: virginiamycin B lyase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.54Å 79.23Å 188.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.89 – 1.90 19.88 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.7 (19.89-1.90) 99.7 (19.88-1.90)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.39 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.140 , 0.199 0.154 , 0.207	Depositor DCC
$R_{free}$ test set	8108 reflections (10.27%)	DCC
Wilson B-factor (Å <sup>2</sup> )	21.1	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 49.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 87070 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10412	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.49% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: CL

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	1.06	3/2424 (0.1%)	0.95	5/3293 (0.2%)
1	B	0.97	1/2389 (0.0%)	0.94	4/3248 (0.1%)
1	C	0.99	3/2413 (0.1%)	0.87	3/3276 (0.1%)
1	D	0.95	0/2392	0.91	3/3248 (0.1%)
All	All	0.99	7/9618 (0.1%)	0.92	15/13065 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	C	0	1
1	D	0	1
All	All	0	3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	132	ARG	CB-CG	-8.33	1.30	1.52
1	C	135	GLU	CG-CD	7.11	1.62	1.51
1	A	198	PHE	CE1-CZ	6.50	1.49	1.37
1	C	135	GLU	CD-OE1	6.50	1.32	1.25
1	B	121	ARG	CB-CG	-5.98	1.36	1.52
1	A	121	ARG	CB-CG	-5.91	1.36	1.52
1	C	121	ARG	CG-CD	5.83	1.66	1.51

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	124	ARG	NE-CZ-NH2	-8.24	116.18	120.30
1	B	250	ARG	NE-CZ-NH1	7.99	124.29	120.30
1	A	8	LEU	CA-CB-CG	7.86	133.38	115.30
1	B	8	LEU	CA-CB-CG	7.44	132.42	115.30
1	B	238	LEU	CA-CB-CG	-7.27	98.59	115.30
1	B	132	ARG	NE-CZ-NH1	6.61	123.61	120.30
1	D	238	LEU	CA-CB-CG	-6.49	100.39	115.30
1	D	121	ARG	NE-CZ-NH2	-6.34	117.13	120.30
1	A	250	ARG	NE-CZ-NH1	5.74	123.17	120.30
1	A	132	ARG	NE-CZ-NH2	-5.66	117.47	120.30
1	A	60	MET	CG-SD-CE	5.52	109.03	100.20
1	A	250	ARG	NE-CZ-NH2	-5.39	117.61	120.30
1	C	111	ASP	CB-CG-OD2	-5.38	113.46	118.30
1	C	291	LEU	CA-CB-CG	5.03	126.86	115.30
1	C	238	LEU	CB-CG-CD2	5.00	119.51	111.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	2	GLU	Peptide
1	C	298	MET	Peptide
1	D	294	ILE	Peptide

## 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	2346	0	2334	40	0
1	B	2311	0	2290	53	0
1	C	2344	0	2312	33	0
1	D	2323	0	2304	33	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	2	0	0	2	0
3	A	327	0	0	7	0
3	B	290	0	0	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	232	0	0	7	0
3	D	234	0	0	9	0
All	All	10412	0	9240	153	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 8.

All (153) 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:C:27:LYS:NZ	1:C:40:CYS:SG	2.28	1.06
1:D:22:VAL:HG12	1:D:273[B]:CYS:SG	2.06	0.95
1:C:22:VAL:HG12	1:C:273[B]:CYS:SG	2.10	0.91
1:A:243:TRP:CZ3	3:D:520:HOH:O	2.28	0.87
1:A:42[A]:ASN:HD21	1:A:46:LYS:H	1.22	0.83
1:C:13:GLN:HE21	1:C:13:GLN:HA	1.46	0.78
1:A:202:ILE:HD11	1:D:51:PRO:HD2	1.69	0.75
1:C:224:ASN:OD1	1:C:226:ARG:NH1	2.21	0.74
1:B:42:ASN:C	1:B:42:ASN:HD22	1.91	0.73
1:C:298:MET:CE	3:C:478:HOH:O	2.37	0.73
1:A:13:GLN:HG3	3:A:547:HOH:O	1.87	0.72
1:C:298:MET:HE3	3:C:478:HOH:O	1.87	0.72
1:A:226:ARG:HH11	1:D:12:ASN:HD22	1.37	0.72
1:D:292:THR:HG23	3:D:333:HOH:O	1.88	0.72
1:A:42[A]:ASN:ND2	1:A:44:ASP:OD2	2.24	0.70
1:B:270:HIS:HD2	3:B:392:HOH:O	1.75	0.70
1:D:84:THR:OG1	1:D:86:LYS:HD2	1.93	0.69
1:B:22:VAL:HG23	1:B:273[A]:CYS:SG	2.33	0.68
1:B:263:GLN:O	1:B:264:ILE:HD12	1.94	0.68
1:D:85[B]:LYS:HE3	3:D:328:HOH:O	1.94	0.67
1:A:270:HIS:CE1	1:A:284:GLU:HB2	2.29	0.67
1:B:42:ASN:HB3	3:B:562:HOH:O	1.94	0.67
1:C:224:ASN:OD1	1:C:226:ARG:CZ	2.44	0.66
1:A:42[A]:ASN:ND2	1:A:46:LYS:H	1.93	0.65
1:C:11:THR:HG21	1:C:49:GLU:OE2	1.97	0.65
1:D:85[B]:LYS:CE	3:D:328:HOH:O	2.45	0.65
1:C:32:GLN:NE2	1:C:39:SER:OG	2.26	0.64
1:C:292:THR:HG23	3:C:398:HOH:O	1.98	0.64
1:D:33:HIS:HD2	3:D:368:HOH:O	1.81	0.64
1:B:7:GLU:OE1	1:B:264:ILE:CD1	2.47	0.63
1:B:4:LYS:HG2	1:B:292[A]:THR:OG1	1.98	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:273[B]:CYS:SG	1:C:274:PHE:N	2.73	0.61
1:A:226:ARG:HH11	1:D:12:ASN:ND2	1.97	0.61
1:A:202:ILE:HD11	1:D:51:PRO:CD	2.31	0.61
1:B:226[B]:ARG:HG3	1:B:226[B]:ARG:HH11	1.66	0.60
1:A:237:ASP:OD1	1:A:252:THR:HG23	2.02	0.60
1:D:273[B]:CYS:SG	3:D:346:HOH:O	2.56	0.59
1:A:255:ASN:ND2	3:A:543:HOH:O	2.36	0.58
1:D:11:THR:OG1	1:D:49:GLU:OE2	2.21	0.58
1:B:270:HIS:HE1	1:B:284:GLU:OE1	1.87	0.57
1:A:270:HIS:HE1	1:A:284:GLU:HB2	1.69	0.56
1:B:265:LYS:HD3	3:B:551:HOH:O	2.04	0.56
1:B:162:ASN:ND2	1:B:183:SER:H	2.03	0.56
1:D:1:MET:HB2	1:D:295:LYS:O	2.06	0.56
1:D:84:THR:HG1	1:D:86:LYS:HD2	1.70	0.56
1:D:162:ASN:ND2	1:D:183:SER:H	2.04	0.56
1:B:78:ASN:HD21	1:B:99:SER:HB2	1.71	0.55
1:A:237:ASP:OD1	1:A:252:THR:CG2	2.54	0.55
1:A:1:MET:HB3	1:A:293:LEU:HB3	1.89	0.55
1:A:42[A]:ASN:HD21	1:A:46:LYS:N	2.00	0.54
1:B:42:ASN:ND2	1:B:42:ASN:C	2.60	0.54
1:A:236:ILE:C	1:A:236:ILE:HD12	2.27	0.54
1:B:7:GLU:OE1	1:B:264:ILE:HG12	2.08	0.54
1:A:243:TRP:HZ3	3:D:479:HOH:O	1.91	0.53
1:C:177:LYS:NZ	3:C:472:HOH:O	2.42	0.52
1:C:20:ILE:HD12	1:C:30:ILE:CD1	2.40	0.52
1:A:42[A]:ASN:ND2	1:A:46:LYS:N	2.58	0.52
1:B:120:ASN:ND2	1:B:141:SER:H	2.08	0.52
1:A:27:LYS:HD3	1:A:40[A]:CYS:SG	2.50	0.52
1:D:265:LYS:HA	1:D:265:LYS:CE	2.39	0.52
1:D:197:TRP:CD2	1:D:238:LEU:HD13	2.44	0.52
1:B:120:ASN:HD21	1:B:141:SER:H	1.58	0.52
1:B:78:ASN:ND2	1:B:99:SER:HB2	2.25	0.52
1:C:27:LYS:NZ	1:C:42:ASN:OD1	2.36	0.51
1:A:65:SER:OG	1:A:67:ASP:OD1	2.27	0.51
1:B:42:ASN:ND2	1:B:44:ASP:OD1	2.22	0.51
1:B:86:LYS:CG	3:B:588:HOH:O	2.58	0.51
1:D:291:LEU:HD23	1:D:291:LEU:C	2.31	0.51
1:C:20:ILE:HD12	1:C:30:ILE:HD12	1.94	0.50
1:B:85:LYS:HG2	3:B:382:HOH:O	2.12	0.50
1:B:197:TRP:CD2	1:B:238:LEU:HD13	2.48	0.49
1:B:3:PHE:HA	1:B:293:LEU:HA	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219[B]:LYS:NZ	3:A:597:HOH:O	2.23	0.49
1:A:22[B]:VAL:HG22	1:A:28:VAL:HG22	1.94	0.49
1:D:85[A]:LYS:HG3	3:D:328:HOH:O	2.12	0.49
1:B:273[B]:CYS:SG	1:B:274:PHE:N	2.86	0.49
1:B:239:TRP:CD1	1:B:279:ILE:HD11	2.48	0.49
1:A:162:ASN:ND2	1:A:183:SER:H	2.11	0.48
1:B:7:GLU:HG2	1:B:283:MET:CE	2.43	0.48
1:B:265:LYS:NZ	3:B:554:HOH:O	2.46	0.48
1:B:78:ASN:ND2	1:B:99:SER:H	2.11	0.48
1:D:116:GLU:OE2	1:D:133:GLU:OE2	2.31	0.48
1:A:297:ASN:ND2	3:A:509:HOH:O	2.43	0.48
1:A:270:HIS:HD2	3:A:592:HOH:O	1.96	0.48
1:C:60:MET:HG3	1:C:75:ASN:HB2	1.95	0.48
1:A:239:TRP:CD1	1:A:279:ILE:HD11	2.48	0.48
1:C:14:ASP:N	1:C:286:ASP:OD2	2.44	0.48
1:B:41:ILE:CD1	1:B:47:ILE:HG12	2.44	0.48
1:D:102:TYR:HB2	1:D:117[A]:MET:HG3	1.95	0.48
1:C:11:THR:CG2	3:C:400:HOH:O	2.62	0.47
1:B:6:GLN:HG2	1:B:290:LYS:HB2	1.97	0.46
1:A:118:ASN:HD22	1:A:118:ASN:N	2.12	0.46
1:D:133:GLU:OE2	2:D:301:CL:CL	2.71	0.46
1:B:51:PRO:HB3	3:B:356:HOH:O	2.14	0.46
1:A:60:MET:SD	1:D:55:PRO:HG2	2.55	0.46
1:B:162:ASN:ND2	1:B:183:SER:OG	2.49	0.46
1:B:226[B]:ARG:HG3	1:B:226[B]:ARG:NH1	2.30	0.46
1:B:214[A]:GLU:OE1	3:B:557:HOH:O	2.21	0.46
1:B:227:PRO:HA	1:B:241:THR:O	2.15	0.46
1:B:33:HIS:HD2	3:B:437:HOH:O	1.98	0.46
1:B:224:ASN:HB3	1:B:226[B]:ARG:HH21	1.80	0.46
1:B:22:VAL:HG22	1:B:28:VAL:HG22	1.98	0.46
1:A:144[B]:SER:HB3	1:A:157:THR:OG1	2.16	0.46
1:B:7:GLU:HG2	1:B:283:MET:HE3	1.98	0.45
1:A:4:LYS:HD2	1:A:5:LEU:N	2.31	0.45
1:C:224:ASN:OD1	1:C:226:ARG:NH2	2.50	0.45
1:B:7:GLU:OE1	1:B:264:ILE:HD13	2.15	0.45
1:D:167:ILE:HA	1:D:172:ASP:O	2.16	0.45
1:A:13:GLN:CD	3:A:619:HOH:O	2.54	0.45
1:A:30[B]:ILE:HD13	1:A:288:ILE:HD11	1.99	0.45
1:C:227:PRO:HA	1:C:241:THR:O	2.17	0.45
1:B:224:ASN:HB3	1:B:226[B]:ARG:NH2	2.32	0.45
1:C:270:HIS:HE1	1:C:284:GLU:HB2	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:162:ASN:ND2	1:C:183:SER:H	2.15	0.45
1:D:84:THR:OG1	1:D:86:LYS:CD	2.64	0.44
1:A:143:PRO:HA	1:A:157:THR:O	2.18	0.44
1:A:298:MET:O	1:B:8:LEU:CD2	2.65	0.44
1:B:27:LYS:HD3	1:B:40[B]:CYS:SG	2.58	0.44
1:A:268:GLU:HB2	1:A:284:GLU:HB3	2.00	0.44
1:B:208:ARG:O	1:B:215:ILE:HA	2.18	0.44
1:D:117[B]:MET:HG2	1:D:142:TYR:HD2	1.82	0.44
1:C:144[B]:SER:HB3	1:C:157:THR:OG1	2.18	0.44
1:A:291:LEU:HD23	1:A:292:THR:N	2.33	0.44
1:B:166:ARG:O	1:B:173:ILE:HA	2.18	0.43
1:B:12:ASN:HB3	1:B:15:THR:HG21	2.00	0.43
1:C:33:HIS:HD2	3:C:397:HOH:O	2.00	0.43
1:A:208:ARG:O	1:A:215:ILE:HA	2.18	0.43
1:D:72:PHE:CZ	1:D:81:GLY:HA3	2.54	0.43
1:B:42:ASN:HB2	3:B:510:HOH:O	2.19	0.43
1:C:2:GLU:HB2	1:C:294:ILE:HB	2.00	0.43
1:B:74:GLU:OE2	1:B:91:GLU:OE2	2.36	0.43
1:B:7:GLU:OE1	1:B:264:ILE:CG1	2.67	0.43
1:C:143:PRO:HA	1:C:157:THR:O	2.19	0.43
1:A:60:MET:HG3	1:A:75:ASN:HB2	1.99	0.43
1:C:4:LYS:HB2	1:C:294:ILE:HD11	2.00	0.43
1:C:152:ASN:HD22	1:C:152:ASN:HA	1.66	0.42
1:C:72:PHE:CZ	1:C:81:GLY:HA3	2.54	0.42
1:D:12:ASN:C	3:D:500:HOH:O	2.58	0.42
1:C:20:ILE:HG13	1:C:30:ILE:CD1	2.50	0.42
1:A:219[B]:LYS:CE	3:A:597:HOH:O	2.66	0.42
1:C:130:LYS:HE3	3:C:520:HOH:O	2.20	0.42
1:B:86:LYS:HG2	3:B:588:HOH:O	2.18	0.41
1:B:250:ARG:O	1:B:257:ILE:HA	2.20	0.41
1:D:273[B]:CYS:SG	1:D:274:PHE:N	2.93	0.41
1:C:200:GLU:OE2	1:C:217:GLU:OE2	2.39	0.41
1:B:42:ASN:HD21	1:B:44:ASP:CG	2.18	0.41
1:B:3:PHE:HB3	1:B:293:LEU:CD2	2.51	0.41
1:B:275:ASP:C	1:B:275:ASP:OD1	2.59	0.41
1:D:96:ASN:HB2	2:D:301:CL:CL	2.58	0.41
1:D:239:TRP:CD1	1:D:279:ILE:HD11	2.56	0.41
1:D:40[B]:CYS:SG	1:D:48:THR:HB	2.61	0.40
1:C:20:ILE:CD1	1:C:30:ILE:CD1	2.99	0.40
1:D:32:GLN:NE2	1:D:39:SER:OG	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	304/299 (102%)	288 (95%)	14 (5%)	2 (1%)	26	14
1	B	299/299 (100%)	287 (96%)	11 (4%)	1 (0%)	46	35
1	C	302/299 (101%)	288 (95%)	13 (4%)	1 (0%)	46	35
1	D	299/299 (100%)	286 (96%)	12 (4%)	1 (0%)	46	35
All	All	1204/1196 (101%)	1149 (95%)	50 (4%)	5 (0%)	39	27

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	286	ASP
1	B	186	VAL
1	A	186	VAL
1	D	186	VAL
1	C	186	VAL

### 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	259/253 (102%)	252 (97%)	7 (3%)	52	43
1	B	255/253 (101%)	247 (97%)	8 (3%)	47	37
1	C	258/253 (102%)	245 (95%)	13 (5%)	30	18
1	D	255/253 (101%)	246 (96%)	9 (4%)	43	31
All	All	1027/1012 (102%)	990 (96%)	37 (4%)	42	30

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42[A]	ASN
1	A	42[C]	ASN
1	A	85	LYS
1	A	118	ASN
1	A	132	ARG
1	A	252	THR
1	A	286	ASP
1	B	2	GLU
1	B	12	ASN
1	B	42	ASN
1	B	86	LYS
1	B	128	ASP
1	B	238	LEU
1	B	253	SER
1	B	265	LYS
1	C	13	GLN
1	C	14	ASP
1	C	25	LYS
1	C	46	LYS
1	C	67	ASP
1	C	127	ASP
1	C	139	LYS
1	C	254	ASN
1	C	277	GLU
1	C	292	THR
1	C	293	LEU
1	C	295	LYS
1	C	298	MET
1	D	1	MET
1	D	6	GLN
1	D	14	ASP
1	D	86	LYS
1	D	226	ARG
1	D	265	LYS
1	D	286	ASP
1	D	292	THR
1	D	295	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:



Mol	Chain	Res	Type
1	A	6	GLN
1	A	32	GLN
1	A	118	ASN
1	A	162	ASN
1	A	270	HIS
1	A	297	ASN
1	B	12	ASN
1	B	32	GLN
1	B	33	HIS
1	B	42	ASN
1	B	75	ASN
1	B	78	ASN
1	B	120	ASN
1	B	159	ASN
1	B	160	GLN
1	B	162	ASN
1	B	270	HIS
1	C	13	GLN
1	C	32	GLN
1	C	33	HIS
1	C	152	ASN
1	C	162	ASN
1	C	254	ASN
1	C	255	ASN
1	C	263	GLN
1	C	270	HIS
1	D	12	ASN
1	D	32	GLN
1	D	33	HIS
1	D	42	ASN
1	D	109	ASN
1	D	162	ASN
1	D	270	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

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.

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	298/299 (99%)	0.59	22 (7%) 17 19	14, 20, 36, 56	6 (2%)
1	B	293/299 (97%)	0.48	18 (6%) 25 27	15, 22, 36, 50	2 (0%)
1	C	299/299 (100%)	0.70	24 (8%) 15 17	16, 24, 39, 51	7 (2%)
1	D	295/299 (98%)	0.60	24 (8%) 15 16	16, 25, 37, 48	5 (1%)
All	All	1185/1196 (99%)	0.59	88 (7%) 17 19	14, 23, 37, 56	20 (1%)

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	6.3
1	A	298	MET	6.1
1	A	297	ASN	5.8
1	C	295	LYS	5.6
1	C	296	ASP	5.2
1	C	13	GLN	4.6
1	B	3	PHE	4.5
1	C	299	GLU	4.2
1	D	243	TRP	4.0
1	A	3	PHE	4.0
1	A	296	ASP	4.0
1	B	263	GLN	3.9
1	B	294	ILE	3.8
1	A	24	ASP	3.7
1	C	80	ILE	3.6
1	A	85	LYS	3.6
1	A	293	LEU	3.5
1	D	44	ASP	3.5
1	D	265	LYS	3.4
1	D	11	THR	3.4
1	D	266	SER	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	265	LYS	3.3
1	A	86	LYS	3.3
1	C	67	ASP	3.2
1	D	295	LYS	3.2
1	C	24	ASP	3.1
1	D	248	ILE	3.1
1	D	254	ASN	3.0
1	A	2	GLU	3.0
1	A	263	GLN	2.9
1	B	5	LEU	2.9
1	C	298	MET	2.8
1	B	44	ASP	2.8
1	D	9	ASN	2.8
1	C	297	ASN	2.7
1	A	294	ILE	2.7
1	D	7	GLU	2.7
1	C	1	MET	2.7
1	C	293	LEU	2.7
1	C	164	ILE	2.6
1	C	11	THR	2.6
1	C	248	ILE	2.6
1	B	10	LEU	2.5
1	B	266	SER	2.5
1	D	285	CYS	2.4
1	C	206	ILE	2.4
1	C	256	ILE	2.4
1	D	282	ALA	2.4
1	C	198	PHE	2.4
1	B	6	GLN	2.4
1	B	248	ILE	2.4
1	B	262	ILE	2.4
1	D	284	GLU	2.3
1	D	164	ILE	2.3
1	C	5	LEU	2.3
1	A	88	ILE	2.3
1	B	264	ILE	2.3
1	B	293	LEU	2.3
1	A	125	ILE	2.3
1	D	128	ASP	2.3
1	B	86	LYS	2.3
1	D	281	PHE	2.2
1	D	269	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	43	LEU	2.2
1	A	6	GLN	2.2
1	B	154	LEU	2.2
1	D	167	ILE	2.1
1	A	42[A]	ASN	2.1
1	C	44	ASP	2.1
1	C	276	GLY	2.1
1	A	94	LEU	2.1
1	A	80	ILE	2.1
1	D	263	GLN	2.1
1	A	112	ILE	2.0
1	C	47	ILE	2.0
1	D	8	LEU	2.0
1	D	199	VAL	2.0
1	C	128	ASP	2.0
1	D	14	ASP	2.0
1	D	224	ASN	2.0
1	D	122	ILE	2.0
1	A	292	THR	2.0
1	B	155	TRP	2.0
1	A	4	LYS	2.0
1	C	46	LYS	2.0
1	B	24	ASP	2.0
1	B	167	ILE	2.0
1	C	178	ILE	2.0

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

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
2	CL	D	301	1/1	0.98	0.23	4.73	38,38,38,38	0
2	CL	D	300	1/1	0.98	0.12	-0.02	25,25,25,25	0
2	CL	C	300	1/1	0.98	0.12	-0.14	23,23,23,23	0
2	CL	B	300	1/1	0.99	0.10	-0.80	21,21,21,21	0
2	CL	A	300	1/1	0.99	0.10	-0.94	20,20,20,20	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.