



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:56 PM GMT

PDB ID : 1MR7  
Title : Crystal Structure of Streptogramin A Acetyltransferase  
Authors : Kehoe, L.E.; Snidwongse, J.; Courvalin, P.; Rafferty, J.B.; Murray, I.A.  
Deposited on : 2002-09-18  
Resolution : 1.80 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

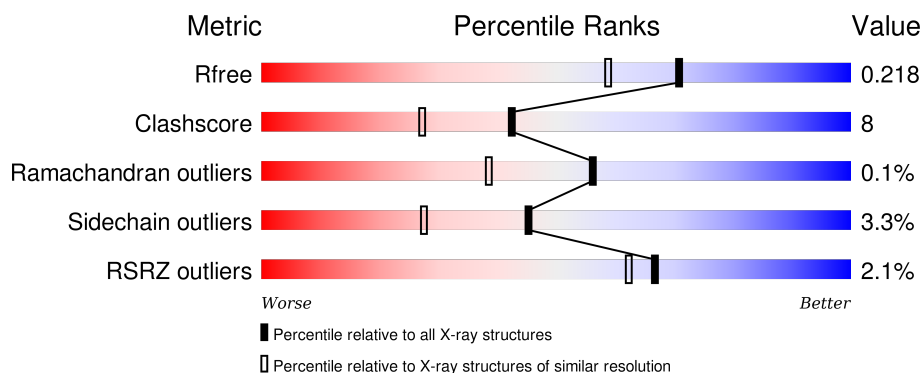
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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	209	<div> <div></div> <div>84% 12% ..</div> </div>
1	B	209	<div> <div></div> <div>81% 15% ..</div> </div>
1	C	209	<div> <div>2%</div> <div>80% 15% ..</div> </div>
1	X	209	<div> <div></div> <div>79% 17% ..</div> </div>
1	Y	209	<div> <div>3%</div> <div>81% 14% ..</div> </div>

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Mol	Chain	Length	Quality of chain
1	Z	209	<div><div></div><div>4%</div><div></div><div>78%</div><div></div><div>18%</div><div></div><div>• •</div></div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10313 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 Streptogramin A Acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	203	Total	C	N	O	S	0	1	0
			1599	1031	260	298	10			
1	B	203	Total	C	N	O	S	0	4	0
			1619	1045	264	297	13			
1	C	202	Total	C	N	O	S	0	3	0
			1612	1042	262	296	12			
1	X	203	Total	C	N	O	S	0	2	0
			1611	1040	263	297	11			
1	Y	203	Total	C	N	O	S	0	1	0
			1591	1026	259	296	10			
1	Z	203	Total	C	N	O	S	0	1	0
			1587	1025	260	292	10			

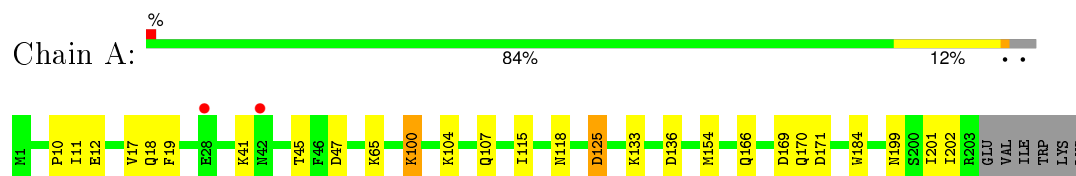
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	103	Total	O	0	0
			103	103		
2	B	129	Total	O	0	0
			129	129		
2	C	101	Total	O	0	0
			101	101		
2	X	124	Total	O	0	0
			124	124		
2	Y	129	Total	O	0	0
			129	129		
2	Z	108	Total	O	0	0
			108	108		

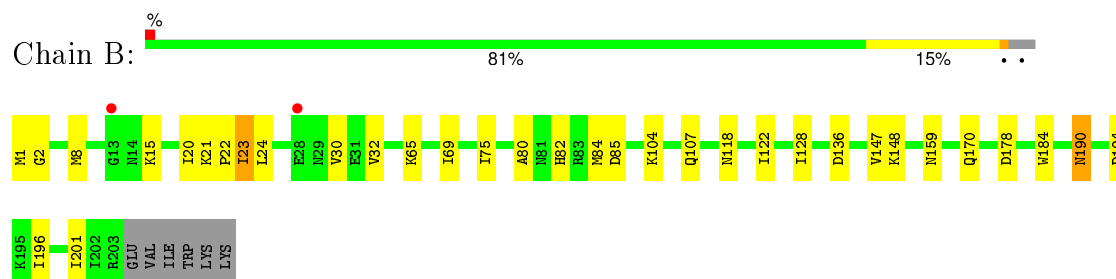
### 3 Residue-property plots [i](#)

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.

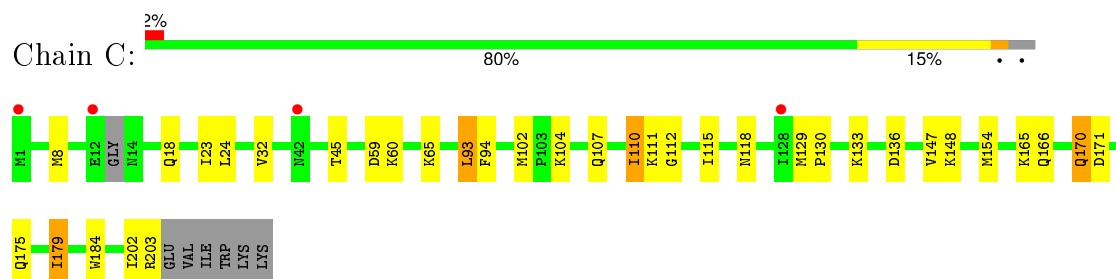
- Molecule 1: Streptogramin A Acetyltransferase



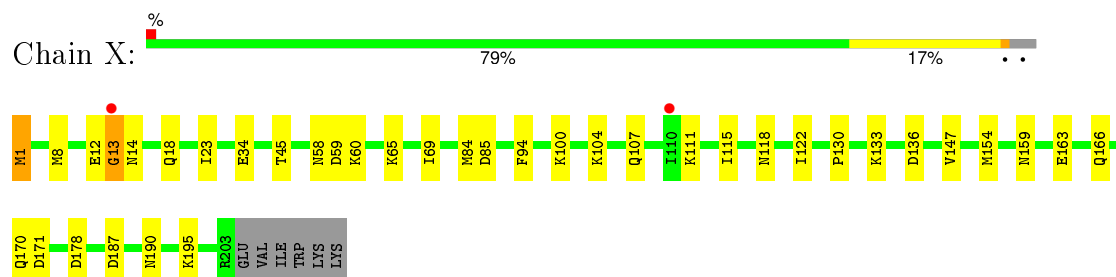
- Molecule 1: Streptogramin A Acetyltransferase



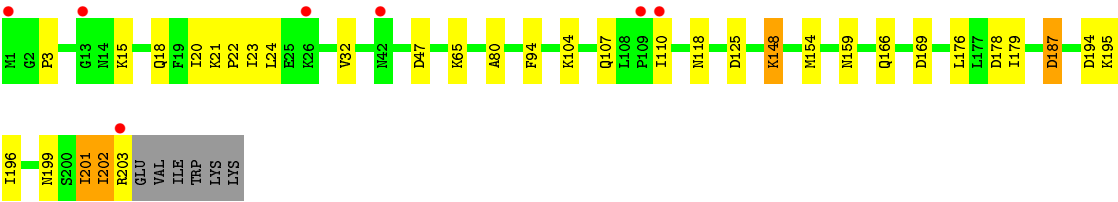
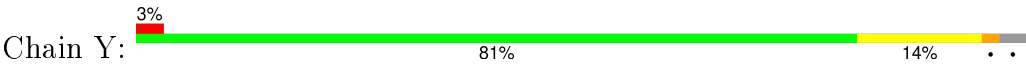
- Molecule 1: Streptogramin A Acetyltransferase



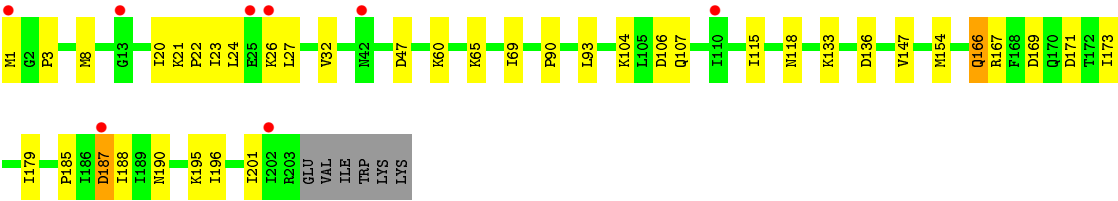
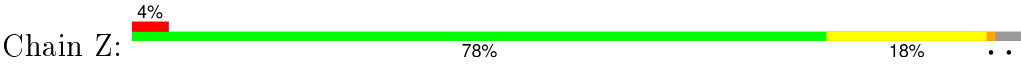
- Molecule 1: Streptogramin A Acetyltransferase



- Molecule 1: Streptogramin A Acetyltransferase



● Molecule 1: Streptogramin A Acetyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	F 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	181.39Å 183.00Å 184.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	21.22 – 1.80 21.21 – 1.80	Depositor EDS
% Data completeness (in resolution range)	97.3 (21.22-1.80) 97.3 (21.21-1.80)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.24 (at 1.80Å)	Xtriage
Refinement program	REFMAC 5	Depositor
R, $R_{free}$	0.176 , 0.207 0.188 , 0.218	Depositor DCC
$R_{free}$ test set	6864 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	23.7	Xtriage
Anisotropy	0.059	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 52.0	EDS
Estimated twinning fraction	0.008 for -h,l,k 0.010 for l,-k,h 0.008 for -k,-h,l 0.003 for -k,-l,h 0.003 for l,-h,-k	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 136323 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	10313	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% 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

### 5.1 Standard geometry

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	0/1638	0.72	5/2225 (0.2%)
1	B	0.54	0/1658	0.77	3/2249 (0.1%)
1	C	0.53	0/1650	0.73	2/2236 (0.1%)
1	X	0.57	0/1650	0.77	5/2238 (0.2%)
1	Y	0.55	0/1630	0.75	5/2216 (0.2%)
1	Z	0.55	0/1626	0.77	5/2210 (0.2%)
All	All	0.55	0/9852	0.75	25/13374 (0.2%)

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	X	0	2
1	Y	0	1
All	All	0	3

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Z	187	ASP	CB-CG-OD2	6.70	124.33	118.30
1	B	136	ASP	CB-CG-OD2	6.68	124.31	118.30
1	Y	125	ASP	CB-CG-OD2	6.46	124.12	118.30
1	C	136	ASP	CB-CG-OD2	6.19	123.87	118.30
1	B	85	ASP	CB-CG-OD2	6.09	123.78	118.30
1	Z	106	ASP	CB-CG-OD2	6.08	123.77	118.30
1	Z	167	ARG	NE-CZ-NH2	-6.07	117.27	120.30
1	Y	169	ASP	CB-CG-OD2	5.99	123.69	118.30
1	X	171	ASP	CB-CG-OD2	5.90	123.61	118.30
1	X	136	ASP	CB-CG-OD2	5.88	123.59	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	47	ASP	CB-CG-OD2	5.88	123.59	118.30
1	A	169	ASP	CB-CG-OD2	5.80	123.52	118.30
1	Z	136	ASP	CB-CG-OD2	5.78	123.50	118.30
1	A	136	ASP	CB-CG-OD2	5.69	123.42	118.30
1	Z	47	ASP	CB-CG-OD2	5.67	123.40	118.30
1	Y	47	ASP	CB-CG-OD2	5.57	123.31	118.30
1	B	178	ASP	CB-CG-OD2	5.54	123.28	118.30
1	C	59	ASP	CB-CG-OD2	5.51	123.26	118.30
1	Y	187	ASP	CB-CG-OD2	5.49	123.24	118.30
1	X	59	ASP	CB-CG-OD2	5.39	123.15	118.30
1	A	171	ASP	CB-CG-OD2	5.26	123.03	118.30
1	Y	178	ASP	CB-CG-OD2	5.20	122.98	118.30
1	X	187	ASP	CB-CG-OD2	5.16	122.94	118.30
1	A	125	ASP	CB-CG-OD2	5.11	122.90	118.30
1	X	85	ASP	CB-CG-OD2	5.03	122.83	118.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	X	12	GLU	Peptide
1	X	13	GLY	Peptide
1	Y	202	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	1599	0	1577	17	0
1	B	1619	0	1604	28	0
1	C	1612	0	1603	27	0
1	X	1611	0	1603	28	0
1	Y	1591	0	1562	29	0
1	Z	1587	0	1565	33	0
2	A	103	0	0	0	0
2	B	129	0	0	4	0
2	C	101	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	X	124	0	0	6	0
2	Y	129	0	0	2	0
2	Z	108	0	0	5	0
All	All	10313	0	9514	158	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 (158) 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:Y:18:GLN:HB3	1:Y:23:ILE:CD1	1.91	1.01
1:B:24:LEU:HD12	1:B:32:VAL:HG21	1.53	0.87
1:Y:202:ILE:O	1:Y:203:ARG:O	1.92	0.85
1:Y:154[A]:MET:SD	1:Y:166:GLN:HG2	2.19	0.82
1:A:65:LYS:H	1:A:118:ASN:HD22	1.28	0.82
1:Z:195:LYS:C	1:Z:201:ILE:HG23	2.01	0.80
1:Y:20:ILE:HG13	1:Y:32:VAL:HG11	1.65	0.79
1:B:196:ILE:HG13	1:B:201:ILE:HD12	1.65	0.79
1:Z:65:LYS:H	1:Z:118:ASN:HD22	1.29	0.77
1:C:110:ILE:HD12	1:C:112:GLY:H	1.49	0.77
1:Z:154[A]:MET:SD	1:Z:166:GLN:HG3	2.24	0.77
1:X:34:GLU:CD	2:X:306:HOH:O	2.22	0.77
1:X:65:LYS:H	1:X:118:ASN:HD22	1.34	0.75
1:Y:18:GLN:HE21	1:Y:23:ILE:HD13	1.50	0.75
1:Y:179:ILE:HD11	1:Y:201:ILE:HD13	1.69	0.73
1:B:104:LYS:H	1:B:107:GLN:HE21	1.34	0.73
1:Z:24:LEU:HD12	1:Z:32:VAL:HG21	1.71	0.72
1:X:34:GLU:CG	2:X:306:HOH:O	2.38	0.72
1:C:154[A]:MET:SD	1:C:166:GLN:HG2	2.28	0.72
1:Y:18:GLN:HB3	1:Y:23:ILE:HD12	1.72	0.71
1:Z:147:VAL:HG22	2:Z:226:HOH:O	1.91	0.70
1:Y:65:LYS:H	1:Y:118:ASN:HD22	1.37	0.70
1:A:154[A]:MET:SD	1:A:166:GLN:HG2	2.31	0.69
1:B:65:LYS:H	1:B:118:ASN:HD22	1.40	0.69
1:Z:147:VAL:CG2	2:Z:246:HOH:O	2.40	0.69
1:A:41:LYS:NZ	1:A:125:ASP:OD2	2.23	0.69
1:Y:196:ILE:HG13	1:Y:201:ILE:HG21	1.74	0.69
1:A:65:LYS:H	1:A:118:ASN:ND2	1.89	0.69
1:Z:65:LYS:H	1:Z:118:ASN:ND2	1.90	0.68
1:A:104:LYS:H	1:A:107:GLN:HE21	1.40	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:104:LYS:H	1:C:107:GLN:HE21	1.41	0.68
1:C:65:LYS:H	1:C:118:ASN:HD22	1.38	0.68
1:Y:18:GLN:HB3	1:Y:23:ILE:HD11	1.77	0.66
1:C:202:ILE:O	1:C:203:ARG:CB	2.43	0.66
1:C:179:ILE:HD12	1:C:184:TRP:CZ2	2.31	0.65
1:B:184:TRP:CH2	1:B:201:ILE:HD11	2.33	0.64
1:Z:196:ILE:HG13	1:Z:201:ILE:HG21	1.81	0.63
1:A:18:GLN:HE22	1:A:45:THR:HA	1.63	0.62
1:X:154[A]:MET:SD	1:X:166:GLN:HG2	2.40	0.62
1:Y:195:LYS:C	1:Y:201:ILE:HG23	2.20	0.62
1:A:17:VAL:HG11	1:C:93:LEU:HD21	1.80	0.62
1:Y:202:ILE:HG13	1:Y:202:ILE:O	1.99	0.61
1:C:93:LEU:HD23	1:C:93:LEU:C	2.20	0.61
1:X:104:LYS:H	1:X:107:GLN:HE21	1.46	0.61
1:B:65:LYS:H	1:B:118:ASN:ND2	1.98	0.61
1:C:65:LYS:H	1:C:118:ASN:ND2	1.97	0.61
1:X:65:LYS:H	1:X:118:ASN:ND2	1.99	0.61
1:Y:65:LYS:H	1:Y:118:ASN:ND2	1.98	0.61
1:C:104:LYS:H	1:C:107:GLN:NE2	1.98	0.60
1:C:179:ILE:CD1	1:C:184:TRP:CZ2	2.84	0.60
1:B:194:ASP:HB2	2:B:273:HOH:O	2.00	0.60
1:Z:190:ASN:ND2	2:Z:260:HOH:O	2.34	0.59
1:X:1:MET:HB3	1:X:190:ASN:ND2	2.18	0.59
1:Z:26:LYS:O	1:Z:27:LEU:HD23	2.02	0.59
1:B:147:VAL:CG2	2:B:305:HOH:O	2.51	0.59
1:X:1:MET:CB	1:X:190:ASN:ND2	2.66	0.58
1:C:175:GLN:O	1:C:179:ILE:HG23	2.03	0.58
1:C:110:ILE:HD12	1:C:111:LYS:N	2.19	0.58
1:B:184:TRP:CZ3	1:B:201:ILE:HD11	2.39	0.57
1:C:110:ILE:CD1	1:C:112:GLY:H	2.16	0.57
1:Z:187:ASP:OD1	1:Z:188:ILE:N	2.38	0.57
1:B:1:MET:HB3	1:B:190:ASN:ND2	2.18	0.57
1:X:34:GLU:HG3	2:X:306:HOH:O	2.01	0.57
1:B:20:ILE:HG23	1:B:24:LEU:HG	1.88	0.56
1:Y:18:GLN:NE2	1:Y:23:ILE:HD13	2.18	0.56
1:A:104:LYS:H	1:A:107:GLN:NE2	2.03	0.56
1:Y:23:ILE:HG22	1:Y:24:LEU:HD23	1.88	0.56
1:Z:179:ILE:HD11	1:Z:201:ILE:HD12	1.89	0.56
1:Y:179:ILE:HD11	1:Y:201:ILE:CD1	2.36	0.56
1:Y:24:LEU:HD12	1:Y:32:VAL:HG21	1.88	0.55
1:Z:104:LYS:H	1:Z:107:GLN:HE21	1.53	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:58:ASN:ND2	1:X:60:LYS:NZ	2.55	0.55
1:B:196:ILE:CG1	1:B:201:ILE:HD12	2.36	0.55
1:Z:169:ASP:OD1	1:Z:171:ASP:HB3	2.07	0.55
1:C:93:LEU:HD22	1:C:94:PHE:CE2	2.41	0.54
1:B:147:VAL:HG22	2:B:301:HOH:O	2.05	0.54
1:X:34:GLU:H	1:X:34:GLU:CD	2.10	0.53
1:X:147:VAL:HG22	2:X:227:HOH:O	2.09	0.53
1:Y:20:ILE:HG13	1:Y:32:VAL:CG1	2.37	0.53
1:X:34:GLU:OE1	1:X:65:LYS:HE2	2.08	0.52
1:B:104:LYS:H	1:B:107:GLN:NE2	2.06	0.52
1:Z:166:GLN:HG2	1:Z:173:ILE:HD11	1.89	0.52
1:Z:147:VAL:HG23	2:Z:246:HOH:O	2.08	0.52
1:X:111:LYS:HD3	1:X:130:PRO:HB2	1.91	0.52
1:B:122:ILE:HD13	1:B:128:ILE:HD11	1.93	0.51
1:X:154[A]:MET:SD	1:X:166:GLN:CG	2.98	0.51
1:Y:20:ILE:CG1	1:Y:32:VAL:HG11	2.40	0.50
1:X:8[A]:MET:HG3	1:X:23:ILE:HD11	1.93	0.50
1:X:104:LYS:H	1:X:107:GLN:NE2	2.09	0.50
1:B:147:VAL:HG23	1:B:148:LYS:HG3	1.94	0.49
1:B:82:HIS:O	1:B:84[A]:MET:HG3	2.13	0.49
1:Y:148:LYS:HD3	2:Y:293:HOH:O	2.12	0.49
1:C:93:LEU:O	1:C:93:LEU:HD23	2.12	0.49
1:B:21:LYS:HB3	1:B:22:PRO:HD3	1.95	0.49
1:Z:196:ILE:N	1:Z:201:ILE:HG23	2.27	0.48
1:Z:104:LYS:H	1:Z:107:GLN:NE2	2.11	0.48
1:A:11:ILE:HD11	1:C:93:LEU:HD23	1.96	0.48
1:Z:195:LYS:O	1:Z:201:ILE:HG23	2.14	0.48
1:Z:187:ASP:OD1	1:Z:188:ILE:HG13	2.13	0.48
1:X:154[B]:MET:CE	1:X:166:GLN:HG2	2.44	0.48
1:Z:115:ILE:HB	1:Z:133:LYS:HD2	1.96	0.48
1:Z:20:ILE:CD1	1:Z:69:ILE:HD12	2.44	0.48
1:A:104:LYS:N	1:A:107:GLN:HE21	2.10	0.47
1:X:147:VAL:CG2	2:X:292:HOH:O	2.63	0.47
1:C:110:ILE:C	1:C:110:ILE:HD12	2.35	0.47
1:Z:21:LYS:HE2	1:Z:32:VAL:O	2.14	0.47
1:Z:166:GLN:HG2	1:Z:173:ILE:CD1	2.45	0.47
1:Y:199:ASN:O	1:Y:202:ILE:HG12	2.15	0.46
1:Y:176:LEU:CD2	1:Y:201:ILE:HD11	2.44	0.46
1:B:21:LYS:HB3	1:B:22:PRO:CD	2.46	0.46
1:B:69:ILE:HD12	1:B:122:ILE:HD12	1.97	0.46
1:Y:21:LYS:HB3	1:Y:22:PRO:CD	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:80:ALA:HB3	2:B:234:HOH:O	2.16	0.46
1:C:93:LEU:HD22	1:C:94:PHE:CZ	2.51	0.46
1:Y:80:ALA:HB3	2:Y:278:HOH:O	2.16	0.45
1:A:184:TRP:CH2	1:A:201:ILE:HD11	2.51	0.45
1:C:166:GLN:HE22	1:C:170:GLN:HG2	1.82	0.45
1:Z:21:LYS:HB3	1:Z:22:PRO:HD3	1.98	0.45
1:X:115:ILE:HB	1:X:133:LYS:HD2	1.98	0.45
1:B:1:MET:HB3	1:B:190:ASN:CG	2.36	0.45
1:X:18:GLN:HE22	1:X:45:THR:HA	1.81	0.45
1:B:20:ILE:CG2	1:B:24:LEU:HG	2.46	0.45
1:Z:154[B]:MET:HE3	2:Z:311:HOH:O	2.17	0.45
1:Y:21:LYS:HB3	1:Y:22:PRO:HD3	1.98	0.45
1:X:94:PHE:HB3	1:Y:3:PRO:HG3	1.98	0.45
1:X:69:ILE:HD12	1:X:122:ILE:HD12	1.98	0.44
1:Z:196:ILE:CG1	1:Z:201:ILE:HG21	2.46	0.44
1:B:24:LEU:HD13	1:B:30:VAL:HG11	1.99	0.44
1:A:10:PRO:HB3	1:A:19:PHE:CE1	2.52	0.44
1:Y:195:LYS:CB	1:Y:201:ILE:HG23	2.47	0.43
1:Z:26:LYS:O	1:Z:27:LEU:CD2	2.66	0.43
1:A:115:ILE:HB	1:A:133:LYS:HD2	1.99	0.43
1:C:8[A]:MET:HG3	1:C:23:ILE:HD11	2.00	0.43
1:Y:94:PHE:HB3	1:Z:3:PRO:HG3	1.99	0.43
1:X:1:MET:HB2	1:X:190:ASN:ND2	2.33	0.43
1:A:18:GLN:NE2	1:A:45:THR:HA	2.31	0.43
1:B:8[B]:MET:HG3	1:B:23:ILE:CD1	2.48	0.43
1:C:115:ILE:HB	1:C:133:LYS:HD2	2.01	0.42
1:C:171:ASP:O	1:C:175:GLN:HG3	2.20	0.42
1:A:199:ASN:O	1:A:202:ILE:HG12	2.18	0.42
1:Z:21:LYS:HB3	1:Z:22:PRO:CD	2.50	0.42
1:Z:185:PRO:HD2	1:Z:188:ILE:HD12	2.02	0.42
1:B:75:ILE:HG12	1:B:128:ILE:HD12	2.02	0.42
1:C:129:MET:HB3	1:C:130:PRO:CD	2.50	0.42
1:C:24:LEU:HD12	1:C:32:VAL:HG21	2.02	0.42
1:X:195:LYS:HG3	2:X:304:HOH:O	2.20	0.41
1:C:147:VAL:HG22	2:C:244:HOH:O	2.20	0.41
1:Z:8:MET:HE3	1:Z:8:MET:HB3	1.84	0.41
1:A:184:TRP:HH2	1:A:201:ILE:HD11	1.85	0.41
1:A:100:LYS:HB3	1:A:100:LYS:HE3	1.72	0.41
1:Y:104:LYS:H	1:Y:107:GLN:HE21	1.67	0.41
1:X:8[B]:MET:O	1:X:18:GLN:HA	2.21	0.41
1:B:8[B]:MET:HG3	1:B:23:ILE:HD11	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:8[A]:MET:O	1:X:18:GLN:HA	2.21	0.40
1:X:154[B]:MET:HG2	1:X:163:GLU:HG2	2.03	0.40
1:C:18:GLN:HE22	1:C:45:THR:HA	1.86	0.40
1:B:1:MET:HG2	1:B:2:GLY:N	2.35	0.40
1:Z:90:PRO:HB2	1:Z:93:LEU:HG	2.04	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	202/209 (97%)	197 (98%)	5 (2%)	0	100	100
1	B	205/209 (98%)	200 (98%)	5 (2%)	0	100	100
1	C	201/209 (96%)	195 (97%)	6 (3%)	0	100	100
1	X	203/209 (97%)	198 (98%)	4 (2%)	1 (0%)	34	17
1	Y	202/209 (97%)	196 (97%)	6 (3%)	0	100	100
1	Z	202/209 (97%)	197 (98%)	5 (2%)	0	100	100
All	All	1215/1254 (97%)	1183 (97%)	31 (3%)	1 (0%)	56	38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	X	13	GLY

### 5.3.2 Protein sidechains [i](#)

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	174/183 (95%)	171 (98%)	3 (2%)	68	57
1	B	176/183 (96%)	171 (97%)	5 (3%)	51	35
1	C	176/183 (96%)	168 (96%)	8 (4%)	34	16
1	X	176/183 (96%)	169 (96%)	7 (4%)	38	20
1	Y	172/183 (94%)	165 (96%)	7 (4%)	37	19
1	Z	171/183 (93%)	167 (98%)	4 (2%)	58	42
All	All	1045/1098 (95%)	1011 (97%)	34 (3%)	45	27

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	GLU
1	A	100	LYS
1	A	170	GLN
1	B	15	LYS
1	B	23	ILE
1	B	159	ASN
1	B	170	GLN
1	B	190	ASN
1	C	60	LYS
1	C	93	LEU
1	C	102	MET
1	C	110	ILE
1	C	148	LYS
1	C	165	LYS
1	C	170	GLN
1	C	179	ILE
1	X	1	MET
1	X	14	ASN
1	X	84	MET
1	X	100	LYS
1	X	159	ASN
1	X	170	GLN
1	X	178	ASP
1	Y	15	LYS
1	Y	110	ILE
1	Y	148	LYS
1	Y	159	ASN

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Mol	Chain	Res	Type
1	Y	187	ASP
1	Y	194	ASP
1	Y	201	ILE
1	Z	1	MET
1	Z	23	ILE
1	Z	60	LYS
1	Z	166	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	58	ASN
1	A	107	GLN
1	A	118	ASN
1	A	166	GLN
1	A	170	GLN
1	B	18	GLN
1	B	107	GLN
1	B	118	ASN
1	B	159	ASN
1	B	162	ASN
1	B	170	GLN
1	C	18	GLN
1	C	107	GLN
1	C	118	ASN
1	C	166	GLN
1	C	175	GLN
1	X	14	ASN
1	X	18	GLN
1	X	58	ASN
1	X	107	GLN
1	X	118	ASN
1	X	159	ASN
1	X	166	GLN
1	Y	18	GLN
1	Y	58	ASN
1	Y	107	GLN
1	Y	118	ASN
1	Y	159	ASN
1	Y	166	GLN
1	Z	18	GLN

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Mol	Chain	Res	Type
1	Z	107	GLN
1	Z	118	ASN
1	Z	162	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

### 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	203/209 (97%)	-0.21	2 (0%) 84 82	10, 17, 27, 36	0
1	B	203/209 (97%)	-0.13	2 (0%) 84 82	11, 16, 30, 40	0
1	C	202/209 (96%)	-0.14	4 (1%) 68 64	9, 17, 28, 41	0
1	X	203/209 (97%)	-0.21	2 (0%) 84 82	10, 17, 30, 46	0
1	Y	203/209 (97%)	-0.15	7 (3%) 49 43	9, 17, 30, 43	0
1	Z	203/209 (97%)	-0.05	8 (3%) 43 37	10, 17, 32, 50	0
All	All	1217/1254 (97%)	-0.15	25 (2%) 67 62	9, 17, 30, 50	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	Z	26	LYS	5.8
1	X	13	GLY	4.7
1	B	13	GLY	4.3
1	X	110	ILE	3.7
1	Z	13	GLY	3.4
1	Y	110	ILE	3.3
1	Z	187	ASP	2.7
1	Z	1	MET	2.7
1	Z	42	ASN	2.6
1	Z	202	ILE	2.6
1	Z	110	ILE	2.5
1	Y	109	PRO	2.5
1	C	12	GLU	2.4
1	Y	42	ASN	2.4
1	A	28	GLU	2.4
1	Y	203	ARG	2.4
1	A	42	ASN	2.2
1	C	42	ASN	2.2
1	Y	13	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	28	GLU	2.1
1	Y	1	MET	2.1
1	Z	25	GLU	2.1
1	C	1	MET	2.1
1	Y	26	LYS	2.1
1	C	128	ILE	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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.