



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 10, 2020 – 10:45 AM BST

PDB ID : 6HYH  
Title : Crystal structure of MSMEG\_1712 from Mycobacterium smegmatis in complex with Beta-D-Fucofuranose  
Authors : Li, M.; Mueller, C.; Zhang, L.; Einsle, O.; Jessen-Trefzer, C.  
Deposited on : 2018-10-21  
Resolution : 2.50 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
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) : 2.13.1

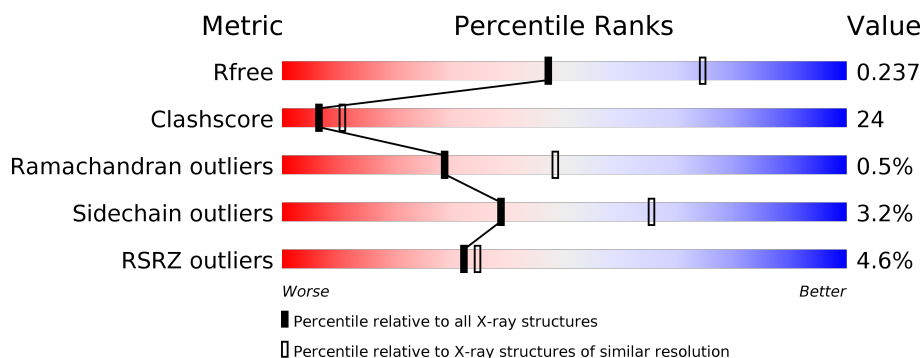
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 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	319	
1	B	319	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4652 atoms, of which 24 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 Periplasmic binding protein/LacI transcriptional regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	304	Total	C	N	O	S	0	0	0
			2283	1427	398	452	6			
1	B	304	Total	C	N	O	S	0	0	0
			2283	1427	398	452	6			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	308	LYS	-	expression tag	UNP I7G686
A	309	LEU	-	expression tag	UNP I7G686
A	310	ALA	-	expression tag	UNP I7G686
A	311	ALA	-	expression tag	UNP I7G686
A	312	ALA	-	expression tag	UNP I7G686
A	313	LEU	-	expression tag	UNP I7G686
A	314	GLU	-	expression tag	UNP I7G686
A	315	HIS	-	expression tag	UNP I7G686
A	316	HIS	-	expression tag	UNP I7G686
A	317	HIS	-	expression tag	UNP I7G686
A	318	HIS	-	expression tag	UNP I7G686
A	319	HIS	-	expression tag	UNP I7G686
A	320	HIS	-	expression tag	UNP I7G686
B	308	LYS	-	expression tag	UNP I7G686
B	309	LEU	-	expression tag	UNP I7G686
B	310	ALA	-	expression tag	UNP I7G686
B	311	ALA	-	expression tag	UNP I7G686
B	312	ALA	-	expression tag	UNP I7G686
B	313	LEU	-	expression tag	UNP I7G686
B	314	GLU	-	expression tag	UNP I7G686
B	315	HIS	-	expression tag	UNP I7G686
B	316	HIS	-	expression tag	UNP I7G686
B	317	HIS	-	expression tag	UNP I7G686
B	318	HIS	-	expression tag	UNP I7G686
B	319	HIS	-	expression tag	UNP I7G686

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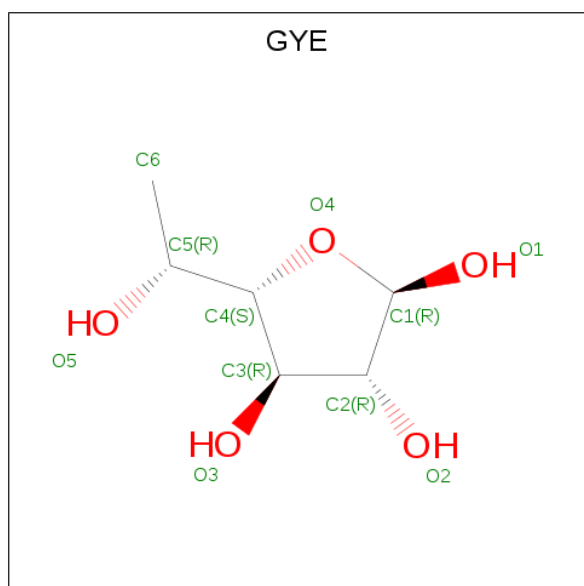
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Chain	Residue	Modelled	Actual	Comment	Reference
B	320	HIS	-	expression tag	UNP I7G686

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	5	Total	Zn	0	0
			5	5		
2	A	8	Total	Zn	0	0
			8	8		

- Molecule 3 is beta-D-fucofuranose (three-letter code: GYE) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			23	6	12	5		
3	B	1	Total	C	H	O	0	0
			23	6	12	5		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	15	Total	O	0	0
			15	15		
4	B	12	Total	O	0	0
			12	12		

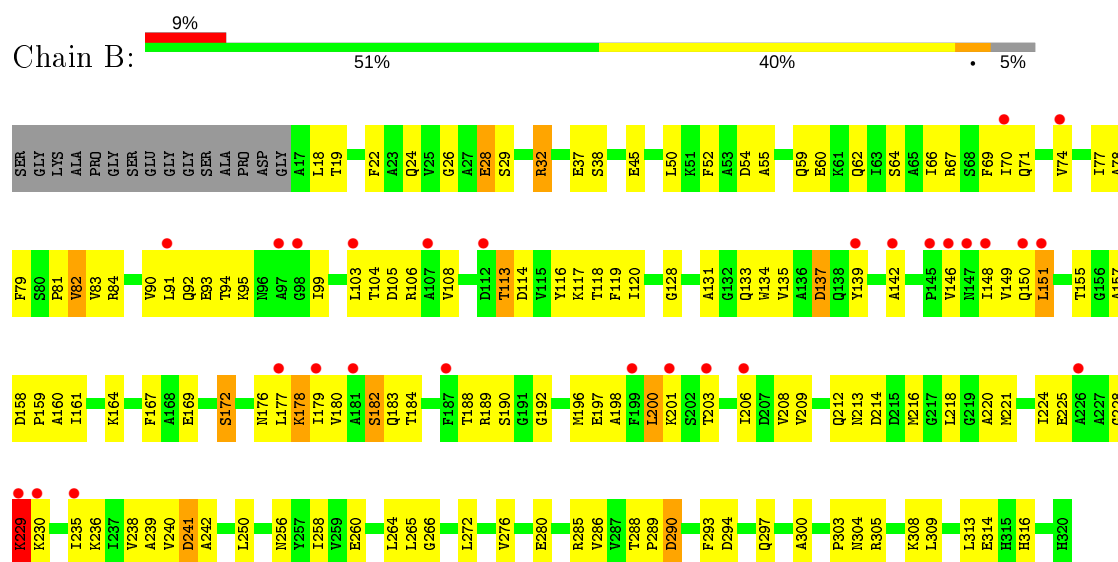
### 3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Periplasmic binding protein/LacI transcriptional regulator



- Molecule 1: Periplasmic binding protein/LacI transcriptional regulator



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	118.06Å 118.06Å 232.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	52.62 – 2.50 52.62 – 2.50	Depositor EDS
% Data completeness (in resolution range)	67.7 (52.62-2.50) 67.7 (52.62-2.50)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.11 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.14_3260, PHENIX 1.14_3260	Depositor
R, $R_{free}$	0.211 , 0.238 0.211 , 0.237	Depositor DCC
$R_{free}$ test set	936 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	63.0	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 40.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4652	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.54% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, GYE

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	0.56	0/2324	0.73	4/3152 (0.1%)
1	B	0.51	1/2324 (0.0%)	0.83	7/3152 (0.2%)
All	All	0.53	1/4648 (0.0%)	0.78	11/6304 (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	B	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	45	GLU	CB-CG	-5.50	1.41	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	200	LEU	CA-CB-CG	9.96	138.21	115.30
1	B	137	ASP	CB-CG-OD2	-9.71	109.56	118.30
1	B	45	GLU	CB-CA-C	-8.99	92.42	110.40
1	B	137	ASP	CB-CG-OD1	8.97	126.37	118.30
1	B	229	LYS	CD-CE-NZ	-6.86	95.92	111.70
1	A	305	ARG	NE-CZ-NH1	-6.55	117.02	120.30
1	A	303	PRO	C-N-CA	-5.60	107.70	121.70
1	A	202	SER	C-N-CA	-5.26	108.55	121.70
1	B	290	ASP	CB-CG-OD1	5.21	122.99	118.30
1	A	265	LEU	CB-CG-CD2	-5.18	102.19	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	B	45	GLU	N-CA-CB	5.03	119.65	110.60

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	113	THR	Peptide
1	B	229	LYS	Peptide
1	B	28	GLU	Peptide

## 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	2283	0	2234	66	0
1	B	2283	0	2233	150	0
2	A	8	0	0	0	0
2	B	5	0	0	0	0
3	A	11	12	0	0	0
3	B	11	12	0	1	0
4	A	15	0	0	1	0
4	B	12	0	0	2	0
All	All	4628	24	4467	216	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 24.

All (216) 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:150:GLN:HB3	1:B:182:SER:HB2	1.51	0.93
1:A:294:ASP:OD1	1:A:297:GLN:HG3	1.76	0.86
1:B:250:LEU:HD22	1:B:258:ILE:HG13	1.60	0.83
1:B:197:GLU:HA	1:B:200:LEU:HG	1.62	0.82
1:B:264:LEU:O	1:B:313:LEU:HD22	1.80	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:305:ARG:HG3	1:B:305:ARG:HH11	1.44	0.82
1:A:304:ASN:H	1:A:308:LYS:HE3	1.46	0.81
1:A:252:ASP:OD2	4:A:501:HOH:O	1.96	0.80
1:B:178:LYS:CE	1:B:179:ILE:N	2.43	0.80
1:B:314:GLU:OE2	4:B:501:HOH:O	1.98	0.80
1:B:92:GLN:O	1:B:95:LYS:HB3	1.81	0.79
1:B:178:LYS:HE3	1:B:179:ILE:N	1.97	0.78
1:A:238:VAL:HG13	1:A:257:TYR:HB3	1.64	0.78
1:B:183:GLN:HG2	1:B:184:THR:H	1.48	0.78
1:A:18:LEU:HD23	1:A:19:THR:N	1.99	0.77
1:B:74:VAL:HG21	1:B:77:ILE:HG13	1.66	0.76
1:B:220:ALA:O	1:B:224:ILE:HD12	1.86	0.76
1:A:149:VAL:HG11	1:A:199:PHE:CD2	2.21	0.76
1:B:28:GLU:OE1	1:B:28:GLU:N	2.14	0.75
1:B:69:PHE:HB3	1:B:77:ILE:HD11	1.69	0.74
1:B:161:ILE:O	1:B:164:LYS:HB3	1.87	0.74
1:B:180:VAL:HG21	1:B:203:THR:HG21	1.68	0.74
1:B:139:TYR:CD1	1:B:146:VAL:HG11	2.23	0.73
1:B:183:GLN:HG2	1:B:184:THR:N	2.04	0.72
1:B:303:PRO:O	1:B:308:LYS:HE3	1.89	0.72
1:A:68:SER:O	1:A:72:GLN:HG3	1.90	0.72
1:B:250:LEU:HB2	1:B:258:ILE:HD11	1.70	0.71
1:A:180:VAL:HG21	1:A:203:THR:HG21	1.71	0.71
1:B:103:LEU:HD13	1:B:116:TYR:CD2	2.26	0.71
1:B:18:LEU:HD23	1:B:19:THR:N	2.07	0.70
1:A:103:LEU:HD21	1:A:108:VAL:HG13	1.76	0.68
1:A:183:GLN:HG2	1:A:184:THR:H	1.58	0.68
1:A:183:GLN:HG2	1:A:184:THR:N	2.09	0.68
1:B:178:LYS:HE2	1:B:179:ILE:N	2.08	0.67
1:B:81:PRO:HG3	1:B:108:VAL:HG13	1.77	0.67
1:B:55:ALA:CB	1:B:62:GLN:HG3	2.24	0.67
1:B:196:MET:HG2	1:B:220:ALA:HB1	1.76	0.66
1:B:62:GLN:OE1	1:B:82:VAL:HG23	1.95	0.66
1:B:169:GLU:O	1:B:172:SER:HB3	1.96	0.65
1:A:103:LEU:HB3	1:A:119:PHE:CD1	2.31	0.65
1:B:103:LEU:HD13	1:B:116:TYR:CG	2.31	0.65
1:B:106:ARG:HG2	1:B:159:PRO:HD3	1.78	0.65
1:B:67:ARG:NH1	1:B:93:GLU:OE1	2.29	0.65
1:A:181:ALA:HB1	1:A:199:PHE:CE1	2.32	0.64
1:A:149:VAL:CG1	1:A:199:PHE:CD2	2.80	0.64
1:A:193:LYS:O	1:A:197:GLU:HG3	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:GLN:O	1:B:137:ASP:OD2	2.16	0.64
1:A:264:LEU:HD12	1:A:309:LEU:HB3	1.81	0.63
1:B:178:LYS:HE3	1:B:179:ILE:H	1.62	0.63
1:B:209:VAL:HG23	1:B:235:ILE:CG2	2.29	0.62
1:A:186:ASP:O	1:A:188:THR:HG23	1.99	0.62
1:B:178:LYS:HE3	1:B:179:ILE:CA	2.28	0.62
1:B:151:LEU:CD2	1:B:192:GLY:HA2	2.30	0.61
1:B:69:PHE:CB	1:B:77:ILE:HD11	2.31	0.61
1:B:81:PRO:HG3	1:B:108:VAL:CG1	2.30	0.61
1:B:103:LEU:CD1	1:B:116:TYR:CD2	2.84	0.61
1:A:269:LEU:O	1:A:269:LEU:HD12	2.01	0.60
1:B:188:THR:HG22	1:B:190:SER:N	2.16	0.60
1:A:62:GLN:OE1	1:A:82:VAL:HG23	2.01	0.59
1:A:213:ASN:OD1	1:A:216:MET:HG3	2.02	0.59
1:B:131:ALA:O	1:B:135:VAL:HG23	2.03	0.59
1:B:134:TRP:CE3	1:B:238:VAL:HG21	2.39	0.58
1:B:288:THR:HG22	1:B:313:LEU:HD11	1.86	0.58
1:A:105:ASP:OD1	1:A:106:ARG:N	2.37	0.57
1:B:77:ILE:HD12	1:B:99:ILE:HG21	1.86	0.57
1:B:149:VAL:HG21	1:B:206:ILE:HD13	1.87	0.57
1:B:139:TYR:CE1	1:B:146:VAL:HG11	2.39	0.56
1:B:149:VAL:HG23	1:B:206:ILE:HG23	1.87	0.56
1:B:113:THR:HG22	1:B:114:ASP:H	1.70	0.56
1:B:79:PHE:HE2	1:B:81:PRO:HB3	1.70	0.56
1:A:264:LEU:O	1:A:313:LEU:HD22	2.06	0.56
1:B:74:VAL:HG21	1:B:77:ILE:CG1	2.33	0.56
1:A:241:ASP:O	1:A:242:ALA:HB3	2.05	0.56
1:B:24:GLN:OE1	1:B:32:ARG:HD3	2.06	0.56
1:B:236:LYS:HG2	1:B:256:ASN:ND2	2.21	0.56
1:B:59:GLN:N	4:B:502:HOH:O	2.02	0.56
1:B:213:ASN:OD1	1:B:216:MET:HG3	2.06	0.56
1:B:221:MET:O	1:B:225:GLU:HG3	2.05	0.55
1:B:250:LEU:HB2	1:B:258:ILE:CD1	2.35	0.55
1:B:196:MET:O	1:B:200:LEU:HG	2.06	0.55
1:A:22:PHE:HA	1:A:78:ALA:O	2.07	0.54
1:B:70:ILE:HD11	1:B:94:THR:HG22	1.88	0.54
1:A:241:ASP:HB3	1:A:243:THR:HG23	1.89	0.53
1:B:106:ARG:HG2	1:B:159:PRO:CD	2.38	0.53
1:B:38:SER:HB2	1:B:264:LEU:HD23	1.90	0.53
1:B:104:THR:HG22	1:B:120:ILE:CG2	2.39	0.53
1:B:196:MET:HE2	1:B:224:ILE:HD11	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:LEU:HD22	1:B:192:GLY:HA2	1.91	0.53
1:B:305:ARG:HG3	1:B:305:ARG:NH1	2.18	0.53
1:B:304:ASN:O	1:B:308:LYS:HG3	2.09	0.53
1:B:84:ARG:HH21	1:B:161:ILE:CD1	2.22	0.53
1:A:149:VAL:CG1	1:A:199:PHE:CE2	2.92	0.53
1:B:117:LYS:O	1:B:118:THR:HB	2.09	0.53
1:B:134:TRP:CZ3	1:B:238:VAL:CG2	2.91	0.53
1:B:55:ALA:HB2	1:B:62:GLN:HG3	1.91	0.53
1:A:158:ASP:HB3	1:A:159:PRO:CD	2.39	0.52
1:A:304:ASN:N	1:A:308:LYS:HE3	2.21	0.52
1:B:117:LYS:HG3	1:B:276:VAL:HG13	1.92	0.52
1:A:174:ASN:HB3	1:A:177:LEU:CD1	2.39	0.52
1:A:84:ARG:HB3	1:A:109:ASP:HB2	1.92	0.52
1:B:26:GLY:HA2	1:B:54:ASP:OD1	2.10	0.52
1:B:113:THR:HG22	1:B:114:ASP:N	2.25	0.52
1:B:196:MET:HG2	1:B:220:ALA:CB	2.40	0.52
1:B:228:GLY:C	1:B:229:LYS:HG2	2.30	0.51
1:B:142:ALA:O	1:B:176:ASN:ND2	2.44	0.51
1:B:150:GLN:HB3	1:B:182:SER:CB	2.32	0.51
1:B:139:TYR:HE1	1:B:208:VAL:HG21	1.74	0.51
1:B:18:LEU:HD23	1:B:19:THR:H	1.76	0.51
1:B:149:VAL:CG2	1:B:206:ILE:HG23	2.40	0.51
1:A:26:GLY:HA2	1:A:54:ASP:OD1	2.11	0.51
1:B:288:THR:CG2	1:B:313:LEU:HD11	2.40	0.51
1:B:60:GLU:CD	1:B:60:GLU:H	2.14	0.51
1:B:135:VAL:HG12	1:B:177:LEU:CD1	2.41	0.51
1:B:84:ARG:HH21	1:B:161:ILE:HD13	1.76	0.51
1:A:77:ILE:HG13	1:A:99:ILE:HG21	1.93	0.50
1:B:66:ILE:HG21	1:B:90:VAL:HG22	1.94	0.50
1:B:91:LEU:O	1:B:95:LYS:HB2	2.11	0.50
1:A:151:LEU:HD23	1:A:199:PHE:HE2	1.77	0.49
1:A:150:GLN:HG3	1:A:210:PHE:HD2	1.77	0.49
1:A:240:VAL:O	1:A:241:ASP:HB2	2.10	0.49
1:A:89:ALA:O	1:A:93:GLU:HG3	2.13	0.49
1:B:240:VAL:O	1:B:241:ASP:HB2	2.12	0.49
1:B:272:LEU:O	1:B:276:VAL:HG23	2.13	0.49
1:B:70:ILE:HD11	1:B:94:THR:CG2	2.42	0.49
1:A:120:ILE:O	1:A:120:ILE:HG23	2.12	0.49
1:B:250:LEU:O	1:B:250:LEU:HD12	2.12	0.49
1:B:150:GLN:O	1:B:182:SER:HA	2.13	0.49
1:A:158:ASP:HB3	1:A:159:PRO:HD3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:241:ASP:O	1:B:242:ALA:HB3	2.12	0.48
1:B:134:TRP:CZ3	1:B:238:VAL:HG21	2.48	0.48
1:B:240:VAL:O	1:B:241:ASP:CB	2.61	0.48
1:B:37:GLU:HA	1:B:37:GLU:OE1	2.14	0.48
1:A:74:VAL:HG21	1:A:77:ILE:HG12	1.96	0.48
1:B:228:GLY:O	1:B:229:LYS:HG2	2.14	0.48
1:B:55:ALA:HB3	1:B:62:GLN:HG3	1.96	0.48
1:A:150:GLN:HG3	1:A:210:PHE:CD2	2.49	0.48
1:B:83:VAL:HA	1:B:157:ALA:HA	1.95	0.48
1:A:150:GLN:HB3	1:A:182:SER:OG	2.15	0.47
1:B:264:LEU:HD12	1:B:309:LEU:HB3	1.95	0.47
1:B:103:LEU:CD1	1:B:116:TYR:CG	2.98	0.47
1:B:188:THR:HG22	1:B:190:SER:H	1.79	0.47
1:B:229:LYS:C	1:B:230:LYS:HE2	2.34	0.47
1:A:24:GLN:HG3	1:A:25:VAL:N	2.28	0.47
1:B:92:GLN:HA	1:B:95:LYS:CB	2.45	0.47
1:B:214:ASP:O	1:B:218:LEU:HG	2.14	0.47
1:A:180:VAL:CG2	1:A:203:THR:HG21	2.44	0.47
1:B:294:ASP:H	1:B:297:GLN:HE21	1.63	0.46
1:B:148:ILE:CD1	1:B:208:VAL:HB	2.46	0.46
1:B:158:ASP:N	1:B:159:PRO:HD2	2.30	0.46
1:B:22:PHE:HA	1:B:78:ALA:O	2.16	0.46
1:A:163:ARG:NH1	1:A:212:GLN:O	2.49	0.45
1:B:197:GLU:HA	1:B:200:LEU:CG	2.41	0.45
1:B:289:PRO:HD3	1:B:316:HIS:HA	1.97	0.45
1:B:66:ILE:HG21	1:B:90:VAL:CG2	2.46	0.45
1:B:29:SER:OG	3:B:406:GYE:O2	2.27	0.45
1:B:135:VAL:HG12	1:B:177:LEU:HD13	1.99	0.45
1:B:158:ASP:HB3	1:B:159:PRO:HD3	1.98	0.45
1:B:103:LEU:HB3	1:B:119:PHE:CD1	2.51	0.45
1:B:29:SER:HB3	1:B:32:ARG:H	1.81	0.45
1:A:38:SER:HB2	1:A:264:LEU:HD22	1.98	0.45
1:A:74:VAL:HG21	1:A:77:ILE:CG1	2.47	0.45
1:B:158:ASP:HB3	1:B:159:PRO:CD	2.47	0.45
1:A:251:ALA:O	1:A:299:LYS:HE3	2.17	0.45
1:B:50:LEU:HD21	1:B:52:PHE:HD2	1.82	0.45
1:A:188:THR:O	1:A:216:MET:HG2	2.16	0.45
1:B:155:THR:HG23	1:B:184:THR:CB	2.47	0.45
1:B:293:PHE:HA	1:B:297:GLN:HE21	1.82	0.45
1:B:28:GLU:HG2	1:B:29:SER:HB2	1.99	0.44
1:B:139:TYR:HB2	1:B:177:LEU:HD21	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:198:ALA:O	1:B:201:LYS:HB3	2.17	0.44
1:A:257:TYR:CZ	1:A:292:ALA:HB1	2.52	0.44
1:B:84:ARG:HH21	1:B:161:ILE:HB	1.82	0.44
1:B:67:ARG:O	1:B:71:GLN:HB2	2.18	0.44
1:B:300:ALA:O	1:B:303:PRO:HD2	2.18	0.43
1:A:151:LEU:CD2	1:A:199:PHE:HE2	2.31	0.43
1:B:196:MET:CE	1:B:224:ILE:HD11	2.49	0.43
1:B:69:PHE:HB3	1:B:77:ILE:CD1	2.44	0.43
1:B:151:LEU:HD12	1:B:209:VAL:CG1	2.48	0.43
1:B:285:ARG:HG2	1:B:286:VAL:N	2.33	0.43
1:A:240:VAL:O	1:A:241:ASP:CB	2.66	0.43
1:B:160:ALA:HA	1:B:212:GLN:OE1	2.19	0.43
1:B:224:ILE:HG22	1:B:229:LYS:O	2.19	0.43
1:B:117:LYS:HG3	1:B:276:VAL:CG1	2.48	0.43
1:B:94:THR:HB	1:B:99:ILE:HB	2.00	0.43
1:A:271:ASP:O	1:A:275:LYS:HG2	2.18	0.43
1:B:209:VAL:HG23	1:B:235:ILE:HG21	1.99	0.43
1:A:250:LEU:HD22	1:A:258:ILE:HG13	2.00	0.43
1:A:304:ASN:O	1:A:308:LYS:HG3	2.18	0.43
1:B:74:VAL:CG2	1:B:77:ILE:HG13	2.41	0.43
1:A:195:VAL:HG12	1:A:199:PHE:CE2	2.54	0.43
1:B:265:LEU:O	1:B:266:GLY:C	2.56	0.43
1:A:264:LEU:HD12	1:A:309:LEU:CB	2.47	0.42
1:B:139:TYR:CB	1:B:177:LEU:HD21	2.49	0.42
1:B:104:THR:O	1:B:105:ASP:HB3	2.19	0.42
1:B:148:ILE:HB	1:B:179:ILE:HA	1.99	0.42
1:A:242:ALA:HA	1:A:258:ILE:HG21	2.01	0.42
1:A:289:PRO:HG2	1:A:312:ALA:O	2.19	0.42
1:A:29:SER:HB2	1:A:215:ASP:OD2	2.19	0.42
1:B:305:ARG:CG	1:B:305:ARG:NH1	2.80	0.42
1:A:159:PRO:O	1:A:163:ARG:HG3	2.20	0.42
1:A:147:ASN:OD1	1:A:180:VAL:HG11	2.20	0.42
1:B:120:ILE:O	1:B:120:ILE:HG23	2.20	0.42
1:B:139:TYR:CG	1:B:146:VAL:HG11	2.53	0.41
1:A:275:LYS:CD	1:A:282:VAL:HG12	2.50	0.41
1:B:260:GLU:O	1:B:290:ASP:HB3	2.19	0.41
1:B:70:ILE:CG1	1:B:94:THR:HG22	2.50	0.41
1:A:169:GLU:HA	1:A:172:SER:HB3	2.03	0.41
1:B:214:ASP:HB3	1:B:239:ALA:CB	2.49	0.41
1:B:160:ALA:O	1:B:164:LYS:HB2	2.20	0.41
1:B:133:GLN:HG3	1:B:137:ASP:OD2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:LEU:HD11	1:A:220:ALA:HB2	2.01	0.41
1:B:208:VAL:HG22	1:B:236:LYS:HB2	2.03	0.41
1:A:265:LEU:O	1:A:266:GLY:C	2.58	0.40
1:B:128:GLY:O	1:B:167:PHE:HA	2.21	0.40
1:B:236:LYS:HG2	1:B:256:ASN:HD21	1.85	0.40
1:A:77:ILE:O	1:A:101:VAL:HA	2.21	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	302/319 (95%)	289 (96%)	12 (4%)	1 (0%)	41	61
1	B	302/319 (95%)	286 (95%)	14 (5%)	2 (1%)	22	39
All	All	604/638 (95%)	575 (95%)	26 (4%)	3 (0%)	29	48

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	241	ASP
1	B	241	ASP
1	B	82	VAL

### 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	233/241 (97%)	226 (97%)	7 (3%)	41	68
1	B	233/241 (97%)	225 (97%)	8 (3%)	37	63
All	All	466/482 (97%)	451 (97%)	15 (3%)	39	65

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	SER
1	A	45	GLU
1	A	79	PHE
1	A	80	SER
1	A	105	ASP
1	A	130	ARG
1	A	202	SER
1	B	32	ARG
1	B	64	SER
1	B	151	LEU
1	B	172	SER
1	B	178	LYS
1	B	182	SER
1	B	189	ARG
1	B	280	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	297	GLN

### 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 ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry

Of 15 ligands modelled in this entry, 13 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	GYE	B	406	-	11,11,11	0.26	0	15,16,16	0.95	0
3	GYE	A	409	-	11,11,11	0.33	0	15,16,16	0.91	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
3	GYE	B	406	-	-	1/4/20/20	0/1/1/1
3	GYE	A	409	-	-	0/4/20/20	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	406	GYE	C3-C4-C5-C6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	406	GYE	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, 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	304/319 (95%)	-0.15	0 100 100	27, 52, 83, 109	0
1	B	304/319 (95%)	0.34	28 (9%) 9 9	30, 88, 129, 164	0
All	All	608/638 (95%)	0.10	28 (4%) 32 34	27, 67, 120, 164	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	199	PHE	5.2
1	B	147	ASN	3.9
1	B	201	LYS	3.9
1	B	177	LEU	3.7
1	B	91	LEU	3.7
1	B	146	VAL	3.5
1	B	70	ILE	3.3
1	B	97	ALA	3.1
1	B	229	LYS	3.1
1	B	142	ALA	3.0
1	B	145	PRO	2.8
1	B	206	ILE	2.7
1	B	151	LEU	2.7
1	B	181	ALA	2.7
1	B	98	GLY	2.6
1	B	112	ASP	2.5
1	B	187	PHE	2.5
1	B	179	ILE	2.4
1	B	150	GLN	2.4
1	B	148	ILE	2.4
1	B	235	ILE	2.3
1	B	226	ALA	2.2
1	B	139	TYR	2.2
1	B	107	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	103	LEU	2.2
1	B	74	VAL	2.2
1	B	203	THR	2.1
1	B	230	LYS	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 monosaccharides 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, 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	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ZN	A	404	1/1	0.91	0.14	114,114,114,114	0
3	GYE	B	406	11/11	0.96	0.18	20,20,55,58	23
3	GYE	A	409	11/11	0.97	0.14	20,20,32,33	0
2	ZN	A	407	1/1	0.98	0.20	53,53,53,53	0
2	ZN	A	405	1/1	0.98	0.19	51,51,51,51	0
2	ZN	A	406	1/1	0.98	0.14	70,70,70,70	0
2	ZN	B	404	1/1	0.98	0.17	53,53,53,53	0
2	ZN	A	403	1/1	0.99	0.15	63,63,63,63	0
2	ZN	B	405	1/1	0.99	0.21	60,60,60,60	0
2	ZN	B	403	1/1	0.99	0.18	49,49,49,49	0
2	ZN	A	402	1/1	0.99	0.17	35,35,35,35	0
2	ZN	A	408	1/1	0.99	0.21	60,60,60,60	0
2	ZN	A	401	1/1	0.99	0.15	33,33,33,33	0
2	ZN	B	401	1/1	0.99	0.17	46,46,46,46	0
2	ZN	B	402	1/1	1.00	0.18	45,45,45,45	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.