



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:30 PM GMT

PDB ID : 4IKF  
Title : PFV intasome with inhibitor MB-76  
Authors : Taltynov, O.; Demeulemeester, J.; Desimmie, B.A.; Suchaud, V.; Billamboz, M.; Lion, C.; Bailly, F.; Debyser, Z.; Cotellet, P.; Christ, F.; Strelkov, S.V.  
Deposited on : 2012-12-26  
Resolution : 3.40 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

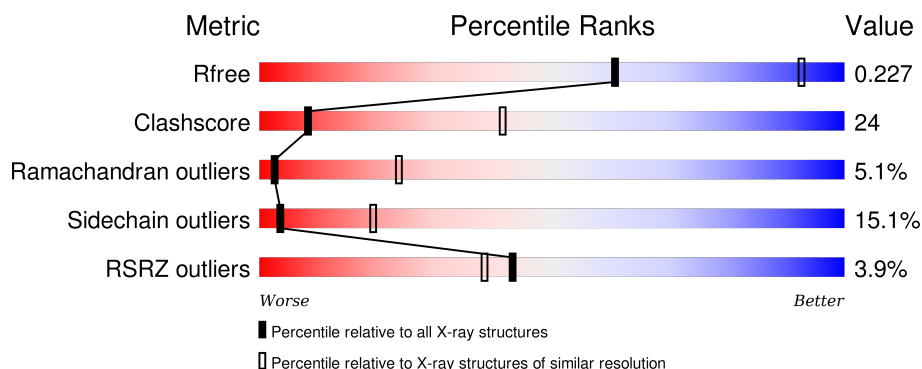
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.40 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	1476 (3.50-3.30)
Clashscore	102246	1611 (3.50-3.30)
Ramachandran outliers	100387	1571 (3.50-3.30)
Sidechain outliers	100360	1571 (3.50-3.30)
RSRZ outliers	91569	1485 (3.50-3.30)

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	395	<div> <div>3%</div> <div>43%</div> <div>43%</div> <div>8%</div> <div>7%</div> </div>
1	B	395	<div> <div>3%</div> <div>22%</div> <div>18%</div> <div>5%</div> <div>53%</div> </div>
2	C	19	<div> <div>5%</div> <div>5%</div> <div>32%</div> <div>63%</div> </div>
3	D	17	<div> <div>6%</div> <div>53%</div> <div>41%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
8	GOL	A	408	-	-	-	X
8	GOL	B	401	-	-	-	X
9	HEZ	B	403	-	-	-	X

## 2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 5177 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 Integrase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	368	Total	C	N	O	S	0	1	0
			2934	1881	517	532	4			
1	B	184	Total	C	N	O	S	0	0	0
			1441	935	234	271	1			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP P14350
A	-1	PRO	-	EXPRESSION TAG	UNP P14350
A	0	GLY	-	EXPRESSION TAG	UNP P14350
A	217	SER	GLY	CONFLICT	UNP P14350
A	218	GLY	SER	CONFLICT	UNP P14350
B	-2	GLY	-	EXPRESSION TAG	UNP P14350
B	-1	PRO	-	EXPRESSION TAG	UNP P14350
B	0	GLY	-	EXPRESSION TAG	UNP P14350
B	217	SER	GLY	CONFLICT	UNP P14350
B	218	GLY	SER	CONFLICT	UNP P14350

- Molecule 2 is a DNA chain called 5'-D(\*AP\*TP\*TP\*GP\*TP\*CP\*AP\*TP\*GP\*GP\*AP\*A P\*TP\*TP\*TP\*CP\*GP\*CP\*A)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	19	Total	C	N	O	P	0	0	0
			387	187	68	114	18			

- Molecule 3 is a DNA chain called 5'-D(\*TP\*GP\*CP\*GP\*AP\*AP\*AP\*TP\*TP\*CP\*CP\*A P\*TP\*GP\*AP\*CP\*A)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	17	Total	C	N	O	P	0	0	0
			345	166	65	98	16			

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

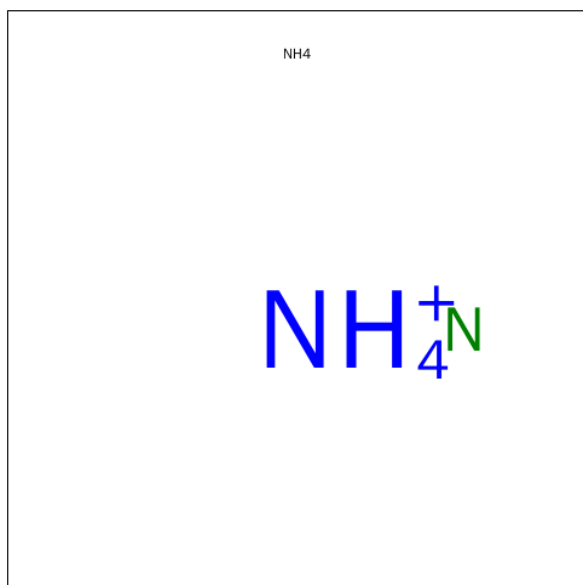
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is AMMONIUM ION (three-letter code: NH4) (formula: H<sub>4</sub>N).

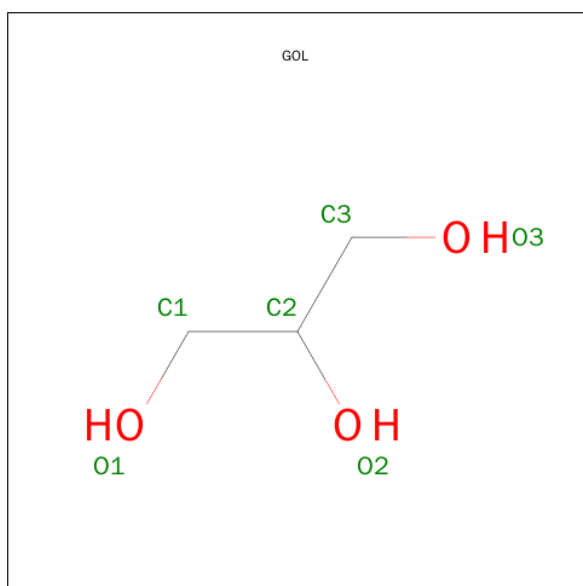


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total N 1 1	0	0

- Molecule 7 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

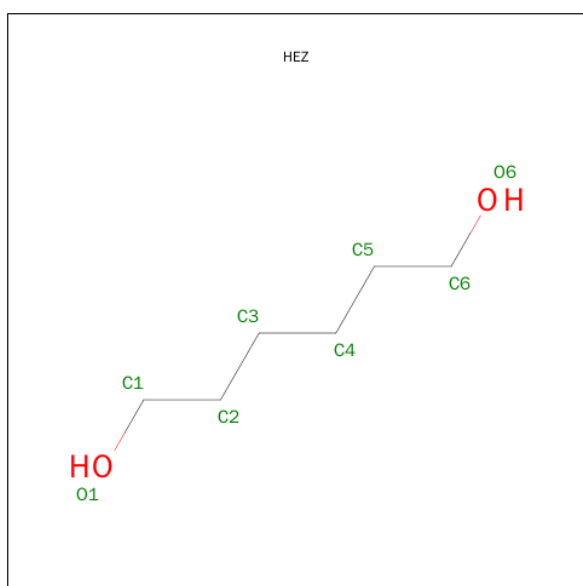
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	2	Total Mg 2 2	0	0

- Molecule 8 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



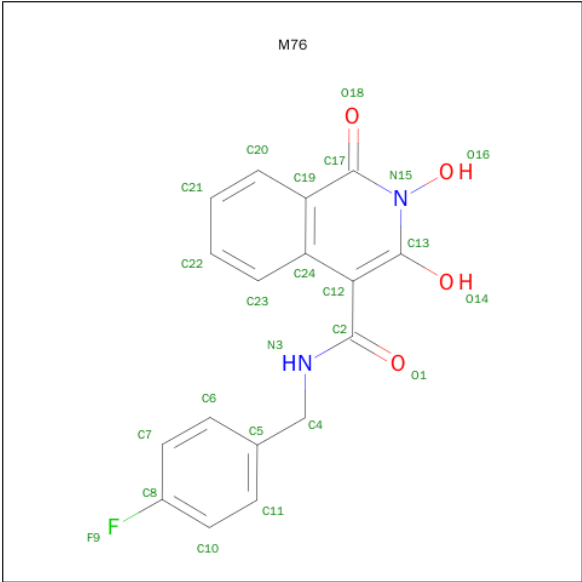
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			6	3	3		
8	A	1	Total	C	O	0	0
			6	3	3		
8	A	1	Total	C	O	0	0
			6	3	3		
8	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 9 is HEXANE-1,6-DIOL (three-letter code: HEZ) (formula:  $C_6H_{14}O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	B	1	Total	C	O	0	0
			8	6	2		

- Molecule 10 is N-(4-FLUOROBENZYL)-2,3-DIHYDROXY-1-OXO-1,2-DIHYDROISOQUINOLINE-4-CARBOXAMIDE (three-letter code: M76) (formula:  $C_{17}H_{13}FN_2O_4$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
10	D	1	Total	C	F	N	O	0	0
			24	17	1	2	4		







● Molecule 2: 5'-D(\*AP\*TP\*TP\*GP\*TP\*CP\*AP\*TP\*GP\*GP\*AP\*AP\*TP\*TP\*TP\*CP\*GP\*CP\*A)-3'



● Molecule 3: 5'-D(\*TP\*GP\*CP\*GP\*AP\*AP\*AP\*TP\*TP\*CP\*CP\*AP\*TP\*GP\*AP\*CP\*A)-3'



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	159.07Å 159.07Å 123.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.85 – 3.40 48.85 – 3.40	Depositor EDS
% Data completeness (in resolution range)	93.8 (48.85-3.40) 93.8 (48.85-3.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.41 (at 3.40Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.2_869)	Depositor
R, $R_{free}$	0.180 , 0.230 0.178 , 0.227	Depositor DCC
$R_{free}$ test set	1080 reflections (5.41%)	DCC
Wilson B-factor (Å <sup>2</sup> )	114.1	Xtriage
Anisotropy	0.002	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 94.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 21042 reflections (0.005%)	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5177	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	115.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.28% 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: GOL, MG, M76, ZN, HEZ, NH4, SO4

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	0.82	1/3013 (0.0%)	0.95	2/4109 (0.0%)
1	B	0.74	0/1481	0.84	0/2025
2	C	1.59	3/433 (0.7%)	2.73	46/667 (6.9%)
3	D	1.32	1/387 (0.3%)	2.24	19/595 (3.2%)
All	All	0.93	5/5314 (0.1%)	1.33	67/7396 (0.9%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	6	DC	C3'-O3'	-8.32	1.33	1.44
2	C	7	DA	N3-C4	-7.01	1.30	1.34
2	C	1	DA	C3'-O3'	5.62	1.51	1.44
3	D	11	DC	N1-C6	-5.44	1.33	1.37
1	A	228	LYS	CD-CE	5.05	1.63	1.51

All (67) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	12	DA	O4'-C1'-N9	-14.35	97.96	108.00
3	D	13	DT	O4'-C1'-N1	-11.48	99.96	108.00
2	C	1	DA	O4'-C1'-N9	-11.15	100.19	108.00
2	C	5	DT	N3-C4-O4	10.87	126.42	119.90
2	C	7	DA	O4'-C1'-N9	-10.49	100.66	108.00
2	C	18	DC	O4'-C4'-C3'	-9.32	100.41	106.00
2	C	8	DT	N3-C4-O4	9.20	125.42	119.90
3	D	11	DC	O4'-C1'-N1	9.07	114.35	108.00
2	C	4	DG	O4'-C1'-N9	8.88	114.22	108.00
2	C	18	DC	C1'-O4'-C4'	-8.86	101.24	110.10
2	C	9	DG	N1-C6-O6	8.53	125.02	119.90
2	C	2	DT	N3-C2-O2	-8.39	117.27	122.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	9	DG	O4'-C1'-N9	-8.31	102.19	108.00
2	C	11	DA	N1-C6-N6	8.28	123.57	118.60
2	C	11	DA	C2-N3-C4	-8.22	106.49	110.60
2	C	6	DC	O4'-C4'-C3'	-8.03	101.18	106.00
2	C	13	DT	N3-C4-O4	7.95	124.67	119.90
2	C	12	DA	P-O5'-C5'	-7.80	108.41	120.90
3	D	17	DA	C1'-O4'-C4'	-7.69	102.41	110.10
2	C	13	DT	C5-C4-O4	-7.61	119.57	124.90
2	C	17	DG	O5'-P-OP2	-7.53	98.92	105.70
2	C	18	DC	O4'-C1'-N1	-7.33	102.87	108.00
2	C	9	DG	N3-C2-N2	-7.32	114.78	119.90
2	C	5	DT	C4-C5-C6	7.31	122.39	118.00
3	D	9	DT	P-O5'-C5'	-7.26	109.28	120.90
3	D	8	DT	O4'-C1'-N1	-7.23	102.94	108.00
3	D	3	DC	P-O5'-C5'	-7.22	109.35	120.90
2	C	9	DG	C5-C6-N1	-7.20	107.90	111.50
2	C	8	DT	C5-C4-O4	-7.14	119.90	124.90
2	C	4	DG	C8-N9-C4	-6.69	103.73	106.40
2	C	14	DT	N3-C4-O4	6.67	123.90	119.90
1	A	336	ARG	NE-CZ-NH2	6.62	123.61	120.30
3	D	14	DG	O4'-C1'-N9	-6.61	103.37	108.00
2	C	12	DA	O5'-P-OP2	-6.56	99.80	105.70
2	C	5	DT	C4-C5-C7	-6.50	115.10	119.00
3	D	17	DA	O4'-C1'-C2'	-6.43	100.76	105.90
3	D	17	DA	N1-C6-N6	6.33	122.40	118.60
3	D	16	DC	O4'-C1'-N1	6.31	112.42	108.00
2	C	14	DT	O4'-C4'-C3'	6.24	109.75	106.00
2	C	13	DT	C1'-O4'-C4'	-6.20	103.91	110.10
2	C	17	DG	O4'-C1'-N9	-6.15	103.70	108.00
2	C	7	DA	C6-N1-C2	-6.14	114.92	118.60
2	C	1	DA	N9-C1'-C2'	6.11	124.21	112.60
3	D	15	DA	O4'-C4'-C3'	6.06	109.64	106.00
2	C	14	DT	C5-C4-O4	-5.89	120.77	124.90
2	C	16	DC	O4'-C1'-N1	-5.84	103.91	108.00
2	C	3	DT	N1-C2-N3	5.77	118.06	114.60
1	A	185	ASP	CB-CG-OD2	5.74	123.47	118.30
2	C	17	DG	OP1-P-OP2	5.72	128.18	119.60
3	D	17	DA	O4'-C1'-N9	5.64	111.95	108.00
2	C	3	DT	N1-C2-O2	-5.58	118.63	123.10
3	D	11	DC	N3-C4-C5	-5.52	119.69	121.90
2	C	4	DG	O5'-P-OP1	-5.51	100.74	105.70
3	D	8	DT	P-O5'-C5'	-5.41	112.24	120.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	5	DT	N1-C2-O2	-5.36	118.81	123.10
2	C	13	DT	P-O5'-C5'	-5.34	112.35	120.90
2	C	5	DT	N3-C4-C5	-5.34	112.00	115.20
3	D	13	DT	N3-C4-O4	5.29	123.08	119.90
2	C	19	DA	N9-C1'-C2'	5.22	122.52	112.60
2	C	17	DG	C5-C6-O6	-5.22	125.47	128.60
3	D	14	DG	P-O5'-C5'	-5.22	112.55	120.90
3	D	17	DA	P-O5'-C5'	-5.21	112.56	120.90
3	D	11	DC	N3-C4-N4	5.21	121.65	118.00
2	C	3	DT	P-O5'-C5'	-5.17	112.63	120.90
2	C	7	DA	C5-C6-N1	5.12	120.26	117.70
2	C	13	DT	O4'-C1'-N1	-5.11	104.42	108.00
2	C	5	DT	O4'-C4'-C3'	5.08	109.05	106.00

There are no chirality outliers.

There are no planarity outliers.

## 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	2934	0	2978	157	1
1	B	1441	0	1421	70	0
2	C	387	0	218	12	0
3	D	345	0	193	14	1
4	A	1	0	0	0	0
5	A	5	0	0	0	0
5	B	5	0	0	0	0
6	A	1	0	0	0	0
7	A	2	0	0	0	0
8	A	18	0	24	1	0
8	B	6	0	8	1	0
9	B	8	0	14	0	0
10	D	24	0	11	6	0
All	All	5177	0	4867	243	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 24.

All (243) 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:170:LEU:O	1:B:174:THR:OG1	1.85	0.93
1:A:296:GLN:O	1:A:300:THR:HG23	1.84	0.76
1:A:81:TRP:CE3	1:A:82:TRP:HA	2.21	0.76
1:A:18:GLY:O	1:A:26:LYS:NZ	2.19	0.75
1:A:292:LEU:HD22	1:B:271:GLY:HA3	1.69	0.75
1:B:235:LEU:HD13	1:B:235:LEU:O	1.86	0.74
2:C:18:DC:H2''	2:C:19:DA:O5'	1.87	0.74
1:A:210:THR:HG23	2:C:3:DT:OP1	1.89	0.72
1:B:133:LEU:HD12	1:B:242:TRP:CZ2	2.24	0.72
1:B:127:ILE:HD11	1:B:182:ILE:CG2	2.22	0.69
1:A:113:LEU:HG	1:A:350:ARG:HG3	1.75	0.68
1:A:341:SER:HB3	1:A:356:ASP:HA	1.76	0.67
1:B:285:ASP:O	1:B:287:THR:N	2.28	0.67
1:B:121:PRO:O	1:B:122:PHE:HB2	1.94	0.67
3:D:17:DA:C2'	10:D:101:M76:C2	2.73	0.67
1:A:373:SER:O	1:A:374:HIS:HB2	1.95	0.66
1:B:263:TYR:CE1	8:B:401:GOL:H32	2.31	0.66
1:B:290:GLU:O	1:B:294:LEU:CB	2.44	0.66
1:A:360:ASN:OD1	1:A:360:ASN:N	2.29	0.65
3:D:17:DA:H2''	10:D:101:M76:C2	2.26	0.65
1:A:108:ALA:O	1:A:314:SER:HA	1.98	0.64
1:A:220:VAL:O	1:A:221:GLU:C	2.36	0.64
1:A:230:LEU:O	1:A:234:LEU:HG	1.97	0.64
1:A:315:TRP:CE2	1:A:371:PRO:HD3	2.33	0.64
1:B:169:SER:O	1:B:172:VAL:HB	1.97	0.64
1:A:253:LEU:HD23	1:A:253:LEU:N	2.13	0.63
1:A:32:LEU:HD23	1:A:33:GLU:N	2.13	0.63
1:A:57:ILE:HG22	1:A:58:VAL:N	2.14	0.62
1:B:287:THR:O	1:B:289:GLU:N	2.33	0.62
1:A:292:LEU:HD22	1:B:271:GLY:CA	2.30	0.62
2:C:1:DA:C2'	2:C:2:DT:H5'	2.30	0.61
1:A:139:TYR:CD2	1:A:243:TYR:CE1	2.88	0.61
1:B:125:PHE:CD2	1:B:179:PRO:HB3	2.36	0.61
1:A:82:TRP:HB2	1:A:83:PRO:HD2	1.83	0.61
1:B:202:ARG:HH11	1:B:202:ARG:HG3	1.66	0.61
1:A:102:THR:HG23	1:A:335:PRO:HA	1.82	0.61
1:A:260:VAL:O	1:A:260:VAL:HG22	2.01	0.61
1:A:111:PRO:O	1:A:350:ARG:HD3	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:HIS:CE1	1:A:215[A]:GLN:HB2	2.36	0.60
1:A:12:LEU:HD22	1:A:37:VAL:HG21	1.84	0.60
1:A:82:TRP:HB2	1:A:83:PRO:CD	2.32	0.60
1:A:217:SER:O	1:A:220:VAL:HG12	2.02	0.60
1:A:331:ALA:O	1:A:336:ARG:NH1	2.34	0.59
1:A:128:ASP:OD1	1:A:185:ASP:OD1	2.18	0.59
1:A:373:SER:O	1:A:374:HIS:CB	2.51	0.59
1:B:160:ALA:HB1	1:B:161:PRO:HD2	1.85	0.59
1:A:344:LEU:HD11	1:A:355:LEU:HB2	1.85	0.58
2:C:1:DA:N3	2:C:1:DA:H5'	2.18	0.58
1:B:127:ILE:HD11	1:B:182:ILE:HG21	1.86	0.57
1:B:202:ARG:NH1	1:B:202:ARG:HG3	2.19	0.57
1:B:131:GLY:HA2	1:B:132:PRO:O	2.04	0.57
3:D:3:DC:H2'	3:D:4:DG:C8	2.40	0.56
1:A:132:PRO:C	1:A:133:LEU:HD23	2.26	0.56
1:B:121:PRO:O	1:B:122:PHE:CB	2.52	0.56
1:A:185:ASP:C	1:A:187:GLY:H	2.09	0.56
1:A:121:PRO:O	1:A:122:PHE:HB2	2.05	0.56
1:B:232:THR:HA	1:B:235:LEU:HD12	1.88	0.55
1:A:319:VAL:HG13	1:A:344:LEU:O	2.06	0.55
1:A:40:SER:HA	1:A:45:VAL:HG22	1.86	0.55
1:A:150:THR:HG21	1:A:269:LEU:HB2	1.88	0.55
1:A:17:GLN:HG2	1:A:18:GLY:N	2.21	0.55
1:A:15:LEU:HD21	1:A:30:TYR:CD1	2.42	0.55
1:A:131:GLY:HA3	1:A:141:TYR:CD1	2.43	0.54
1:A:129:TYR:HD2	1:A:142:VAL:O	1.91	0.54
1:B:255:ASN:C	1:B:264:THR:HG21	2.28	0.54
1:A:374:HIS:O	1:A:375:GLN:HB2	2.07	0.54
1:A:258:SER:OG	1:A:260:VAL:HG12	2.07	0.54
1:A:160:ALA:HB1	1:A:161:PRO:CD	2.37	0.54
1:A:222:ARG:NH2	2:C:6:DC:O2	2.36	0.54
1:A:374:HIS:O	1:A:375:GLN:CB	2.56	0.54
1:A:40:SER:HA	1:A:45:VAL:CG2	2.37	0.54
1:A:230:LEU:O	1:A:230:LEU:HG	2.08	0.54
1:A:73:LEU:HD23	1:A:73:LEU:C	2.28	0.54
1:A:170:LEU:O	1:A:174:THR:HG23	2.08	0.54
1:B:255:ASN:HA	1:B:264:THR:HG21	1.88	0.53
1:A:220:VAL:O	1:A:223:LYS:N	2.41	0.53
1:A:327:VAL:HG23	1:A:336:ARG:C	2.29	0.53
1:A:132:PRO:O	1:A:133:LEU:HD23	2.09	0.53
1:A:123:ASP:O	1:A:181:VAL:HG23	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:254:ASN:HA	1:B:265:PRO:HG2	1.91	0.53
1:A:260:VAL:HG13	1:A:261:LEU:HG	1.90	0.52
1:B:256:THR:HG22	1:B:256:THR:O	2.10	0.52
1:A:251:LEU:HD21	1:B:178:ILE:HD11	1.91	0.52
1:B:229:ARG:HB3	1:B:284:LEU:HD11	1.91	0.52
1:A:202:ARG:HH11	1:A:202:ARG:HG3	1.74	0.52
1:A:281:GLN:O	1:A:282:ASP:C	2.48	0.52
1:A:347:LEU:N	1:A:347:LEU:HD23	2.23	0.52
3:D:17:DA:H2'	10:D:101:M76:C2	2.40	0.52
1:A:327:VAL:HG23	1:A:337:TRP:N	2.25	0.52
1:A:134:PRO:O	1:A:135:PRO:C	2.48	0.52
1:A:167:VAL:O	1:A:168:LYS:C	2.48	0.52
1:B:249:VAL:HG12	1:B:253:LEU:HD11	1.91	0.51
1:A:228:LYS:HD2	3:D:17:DA:OP2	2.11	0.51
1:A:154:TRP:HA	1:A:250:GLN:OE1	2.09	0.51
1:A:296:GLN:O	1:A:300:THR:CG2	2.55	0.51
1:A:327:VAL:HG21	1:A:335:PRO:O	2.11	0.51
1:B:217:SER:O	1:B:218:GLY:C	2.49	0.51
1:B:235:LEU:O	1:B:235:LEU:CD1	2.56	0.51
1:A:139:TYR:CE1	1:A:159:LYS:HG3	2.46	0.51
1:A:304:HIS:CG	1:A:305:PRO:HD2	2.46	0.51
1:A:137:GLN:H	8:A:407:GOL:H2	1.76	0.50
1:B:167:VAL:HG22	1:B:198:TRP:CD1	2.46	0.50
1:A:81:TRP:CE3	1:A:82:TRP:CA	2.94	0.50
1:A:47:ILE:CG2	1:A:48:ILE:N	2.71	0.50
1:A:288:ARG:HB2	1:B:149:MET:CE	2.41	0.50
1:A:14:GLN:H	1:A:14:GLN:CD	2.15	0.50
1:A:358:LEU:N	1:A:359:GLY:HA2	2.26	0.50
3:D:17:DA:H2''	10:D:101:M76:O1	2.12	0.50
1:B:152:PHE:HA	1:B:265:PRO:HB3	1.94	0.50
2:C:17:DG:H2''	2:C:18:DC:O4'	2.11	0.49
1:A:133:LEU:HB3	1:A:134:PRO:HD2	1.94	0.49
1:A:73:LEU:CD1	1:A:86:ARG:HG3	2.43	0.49
1:A:91:LYS:O	1:A:95:ARG:HG3	2.12	0.49
1:B:246:LEU:N	1:B:247:PRO:CD	2.76	0.49
1:A:39:VAL:O	1:A:39:VAL:HG13	2.13	0.49
1:A:286:LEU:HB3	1:A:290:GLU:HB2	1.95	0.49
1:B:256:THR:C	1:B:264:THR:HG22	2.33	0.48
1:A:214:PRO:HB2	1:A:221:GLU:HG3	1.96	0.48
1:A:161:PRO:O	1:A:189:ALA:HB2	2.12	0.48
1:A:318:VAL:CG1	1:A:319:VAL:N	2.76	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:LYS:HE2	3:D:9:DT:OP2	2.14	0.48
1:A:174:THR:HA	1:A:177:ALA:O	2.12	0.48
1:B:230:LEU:O	1:B:230:LEU:HD12	2.14	0.48
1:A:129:TYR:CE1	1:A:161:PRO:HA	2.49	0.48
1:A:245:LEU:O	1:A:246:LEU:C	2.52	0.48
1:A:304:HIS:CD2	1:A:305:PRO:HD2	2.49	0.48
1:B:235:LEU:C	1:B:235:LEU:CD1	2.82	0.47
1:A:86:ARG:O	1:A:90:VAL:HG23	2.14	0.47
1:A:272:ILE:CG1	1:A:273:ASP:O	2.62	0.47
2:C:13:DT:H2''	2:C:14:DT:O5'	2.15	0.47
1:A:327:VAL:HG23	1:A:337:TRP:CA	2.45	0.47
1:A:358:LEU:N	1:A:359:GLY:CA	2.77	0.47
1:B:283:THR:C	1:B:285:ASP:H	2.18	0.47
1:B:233:LYS:HD2	1:B:284:LEU:HA	1.96	0.47
2:C:11:DA:H2''	2:C:12:DA:OP2	2.15	0.47
1:A:12:LEU:CD2	1:A:37:VAL:HG21	2.44	0.47
10:D:101:M76:O1	10:D:101:M76:H13	2.15	0.46
1:B:163:THR:HG23	1:B:195:PHE:HB2	1.96	0.46
1:A:318:VAL:HG13	1:A:319:VAL:N	2.31	0.46
1:A:238:ARG:NH1	1:A:238:ARG:HB2	2.30	0.46
1:A:129:TYR:CZ	1:A:143:LEU:HD13	2.51	0.46
1:A:294:LEU:O	1:A:297:GLU:HB3	2.16	0.46
1:A:139:TYR:CD1	1:A:159:LYS:HG3	2.51	0.46
1:A:26:LYS:O	1:A:28:TYR:N	2.45	0.46
1:A:272:ILE:HG12	1:A:273:ASP:O	2.16	0.46
1:A:315:TRP:CZ2	1:A:371:PRO:HD3	2.51	0.46
1:B:249:VAL:O	1:B:252:ALA:HB3	2.16	0.46
1:A:286:LEU:CB	1:A:290:GLU:HB2	2.45	0.46
1:A:316:SER:HB3	1:A:317:PRO:CD	2.46	0.46
1:B:280:ASN:O	1:B:282:ASP:N	2.48	0.46
1:A:81:TRP:CZ3	1:A:82:TRP:HA	2.50	0.45
1:A:16:LEU:HD21	1:A:31:PHE:C	2.37	0.45
1:A:339:LYS:HD2	1:A:357:HIS:CE1	2.52	0.45
1:A:167:VAL:O	1:A:169:SER:N	2.48	0.45
1:A:348:ASN:O	1:A:350:ARG:N	2.49	0.45
1:A:319:VAL:HG12	1:A:320:GLY:N	2.31	0.45
1:B:181:VAL:CG1	1:B:182:ILE:N	2.79	0.45
1:A:133:LEU:HB3	1:A:134:PRO:CD	2.47	0.45
1:B:123:ASP:OD1	1:B:180:LYS:HE3	2.17	0.45
1:B:232:THR:HA	1:B:235:LEU:CD1	2.46	0.44
1:A:183:HIS:HA	1:A:207:GLU:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:221:GLU:O	3:D:16:DC:H1'	2.17	0.44
1:A:287:THR:OG1	1:A:290:GLU:HG3	2.18	0.44
1:A:152:PHE:CE1	1:A:254:ASN:HB3	2.52	0.44
1:B:288:ARG:HB3	1:B:288:ARG:HH11	1.82	0.44
1:A:21:ILE:HG22	1:A:22:LYS:N	2.32	0.44
1:B:135:PRO:HG3	1:B:140:LEU:HD21	1.98	0.44
1:A:307:THR:HA	1:A:308:PRO:HD2	1.81	0.44
1:A:41:ARG:HB3	1:A:42:PRO:HD2	2.00	0.44
3:D:1:DT:H2''	3:D:2:DG:OP2	2.17	0.44
1:B:245:LEU:C	1:B:247:PRO:HD2	2.38	0.44
1:B:254:ASN:O	1:B:264:THR:HB	2.17	0.43
1:B:149:MET:SD	1:B:149:MET:C	2.97	0.43
1:A:113:LEU:O	2:C:4:DG:H5'	2.18	0.43
1:B:141:TYR:CZ	1:B:161:PRO:HD3	2.53	0.43
1:A:256:THR:O	1:A:256:THR:HG22	2.17	0.43
1:A:47:ILE:HG22	1:A:48:ILE:N	2.34	0.43
1:B:240:THR:O	1:B:240:THR:OG1	2.34	0.43
1:A:111:PRO:HD3	1:A:310:ALA:CB	2.49	0.43
1:A:358:LEU:CB	1:A:359:GLY:HA2	2.49	0.43
1:A:20:TYR:CD2	1:A:20:TYR:C	2.91	0.43
1:B:289:GLU:O	1:B:292:LEU:CB	2.66	0.43
1:A:57:ILE:HD11	1:A:80:TYR:CE1	2.54	0.43
1:A:333:LEU:O	1:A:334:ARG:HD3	2.19	0.43
1:A:220:VAL:O	1:A:222:ARG:N	2.52	0.43
1:B:293:SER:OG	1:B:293:SER:O	2.34	0.43
1:A:121:PRO:O	1:A:122:PHE:CB	2.65	0.43
1:B:222:ARG:O	1:B:223:LYS:C	2.57	0.43
2:C:17:DG:H2''	2:C:18:DC:O5'	2.19	0.42
1:A:356:ASP:OD1	1:A:358:LEU:HB2	2.18	0.42
1:A:185:ASP:O	1:A:187:GLY:N	2.51	0.42
1:A:28:TYR:O	1:A:30:TYR:CD2	2.72	0.42
1:B:136:SER:O	1:B:137:GLN:C	2.56	0.42
1:B:183:HIS:CD2	1:B:184:SER:N	2.87	0.42
3:D:13:DT:H2''	3:D:14:DG:O5'	2.19	0.42
1:A:231:LEU:O	1:A:234:LEU:N	2.51	0.42
1:B:238:ARG:HB2	1:B:241:LYS:HB2	2.01	0.42
1:A:16:LEU:HD11	1:A:32:LEU:HB2	2.01	0.42
1:A:202:ARG:HH11	1:A:202:ARG:CG	2.32	0.42
1:B:209:SER:O	1:B:210:THR:C	2.58	0.42
1:A:142:VAL:HG12	1:A:143:LEU:N	2.35	0.42
1:A:343:VAL:HA	1:A:354:ILE:HG22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:181:VAL:HG12	1:B:182:ILE:N	2.33	0.42
1:A:332:SER:HA	1:A:336:ARG:HH12	1.85	0.42
2:C:14:DT:H2"	2:C:15:DT:OP2	2.20	0.41
3:D:15:DA:O5'	3:D:15:DA:H2'	2.20	0.41
1:A:316:SER:HB3	1:A:317:PRO:HD3	2.02	0.41
1:B:181:VAL:HG22	1:B:205:HIS:HB3	2.03	0.41
1:A:158:THR:OG1	1:A:160:ALA:O	2.31	0.41
1:B:255:ASN:CA	1:B:264:THR:HG21	2.50	0.41
1:A:246:LEU:HB2	1:A:247:PRO:HD3	2.03	0.41
1:B:183:HIS:HA	1:B:207:GLU:O	2.20	0.41
1:A:126:PHE:C	1:A:127:ILE:HG23	2.40	0.41
1:A:227:ILE:O	1:A:228:LYS:C	2.58	0.41
1:A:139:TYR:CD2	1:A:243:TYR:HE1	2.36	0.41
1:A:374:HIS:O	1:A:375:GLN:HG3	2.21	0.41
3:D:9:DT:H1'	3:D:10:DC:H5'	2.02	0.41
3:D:16:DC:H2'	10:D:101:M76:H3	2.02	0.41
1:A:365:SER:C	1:A:367:ASP:N	2.74	0.41
1:A:185:ASP:C	1:A:187:GLY:N	2.74	0.41
1:B:185:ASP:HA	1:B:209:SER:OG	2.20	0.41
1:A:322:LEU:HD23	1:A:322:LEU:HA	1.85	0.41
1:A:40:SER:CA	1:A:45:VAL:HG22	2.51	0.41
1:B:195:PHE:CD2	1:B:206:LEU:HD11	2.56	0.41
3:D:6:DA:H2"	3:D:7:DA:H5'	2.02	0.41
1:B:130:ILE:HG22	1:B:133:LEU:HD21	2.04	0.40
1:A:180:LYS:HE3	1:B:275:ASN:OD1	2.20	0.40
1:A:208:PHE:CD2	1:A:208:PHE:N	2.89	0.40
1:A:269:LEU:O	1:A:269:LEU:HD12	2.22	0.40
1:A:154:TRP:HB3	1:A:156:TYR:HE2	1.85	0.40
1:A:295:LEU:HD12	1:A:295:LEU:HA	1.84	0.40
1:A:65:ALA:O	1:A:66:HIS:C	2.58	0.40
2:C:1:DA:O5'	2:C:2:DT:H72	2.21	0.40
1:A:126:PHE:N	1:A:126:PHE:CD2	2.89	0.40
1:A:82:TRP:CD2	1:A:85:MET:HE3	2.56	0.40
1:B:245:LEU:O	1:B:249:VAL:HG23	2.22	0.40
1:B:136:SER:HB3	1:B:243:TYR:HB2	2.03	0.40
1:B:257:TYR:CD1	1:B:257:TYR:C	2.93	0.40
1:A:76:ILE:O	1:A:77:ALA:C	2.60	0.40
1:A:37:VAL:HG12	1:A:37:VAL:O	2.22	0.40
1:B:230:LEU:CD1	1:B:230:LEU:O	2.69	0.40
1:A:206:LEU:HB3	1:A:208:PHE:HE2	1.85	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:SER:OG	3:D:13:DT:OP1[8_554]	2.05	0.15

## 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	366/395 (93%)	301 (82%)	50 (14%)	15 (4%)	3	30
1	B	182/395 (46%)	140 (77%)	29 (16%)	13 (7%)	1	15
All	All	548/790 (69%)	441 (80%)	79 (14%)	28 (5%)	2	24

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	282	ASP
1	A	328	ALA
1	B	216	SER
1	B	286	LEU
1	B	288	ARG
1	A	366	ILE
1	B	202	ARG
1	B	281	GLN
1	A	27	GLN
1	A	51	GLN
1	A	85	MET
1	A	161	PRO
1	A	186	GLN
1	B	137	GLN
1	B	198	TRP
1	A	33	GLU
1	A	77	ALA
1	A	168	LYS

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Mol	Chain	Res	Type
1	A	188	ALA
1	A	283	THR
1	B	201	GLU
1	A	317	PRO
1	B	192	SER
1	B	218	GLY
1	B	251	LEU
1	A	201	GLU
1	B	287	THR
1	B	210	THR

### 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	330/354 (93%)	285 (86%)	45 (14%)	5	24
1	B	159/354 (45%)	130 (82%)	29 (18%)	2	11
All	All	489/708 (69%)	415 (85%)	74 (15%)	3	19

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ASP
1	A	17	GLN
1	A	20	TYR
1	A	27	GLN
1	A	29	THR
1	A	34	ASP
1	A	57	ILE
1	A	102	THR
1	A	107	LYS
1	A	119	GLN
1	A	124	LYS
1	A	140	LEU
1	A	146	VAL

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Mol	Chain	Res	Type
1	A	158	THR
1	A	178	ILE
1	A	181	VAL
1	A	193	SER
1	A	194	THR
1	A	200	LYS
1	A	206	LEU
1	A	207	GLU
1	A	216	SER
1	A	219	LYS
1	A	226	ASP
1	A	228	LYS
1	A	236	VAL
1	A	248	VAL
1	A	262	LYS
1	A	268	LEU
1	A	272	ILE
1	A	276	THR
1	A	286	LEU
1	A	301	SER
1	A	307	THR
1	A	312	SER
1	A	318	VAL
1	A	332	SER
1	A	341	SER
1	A	342	THR
1	A	357	HIS
1	A	358	LEU
1	A	360	ASN
1	A	365	SER
1	A	370	LYS
1	A	373	SER
1	B	117	ARG
1	B	119	GLN
1	B	125	PHE
1	B	127	ILE
1	B	133	LEU
1	B	137	GLN
1	B	149	MET
1	B	159	LYS
1	B	167	VAL
1	B	168	LYS

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Mol	Chain	Res	Type
1	B	174	THR
1	B	192	SER
1	B	193	SER
1	B	194	THR
1	B	215	GLN
1	B	223	LYS
1	B	230	LEU
1	B	235	LEU
1	B	238	ARG
1	B	240	THR
1	B	253	LEU
1	B	256	THR
1	B	258	SER
1	B	274	SER
1	B	280	ASN
1	B	284	LEU
1	B	286	LEU
1	B	287	THR
1	B	288	ARG

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

Mol	Chain	Res	Type
1	A	357	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 ⓘ

There are no carbohydrates in this entry.



## 5.6 Ligand geometry

Of 12 ligands modelled in this entry, 3 are monoatomic and 1 is modelled with single atom - leaving 8 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	SO4	A	402	-	4,4,4	0.16	0	6,6,6	0.22	0
8	GOL	A	406	-	5,5,5	0.40	0	5,5,5	0.41	0
8	GOL	A	407	-	5,5,5	0.54	0	5,5,5	0.99	0
8	GOL	A	408	-	5,5,5	0.43	0	5,5,5	0.59	0
8	GOL	B	401	-	5,5,5	0.43	0	5,5,5	0.87	0
5	SO4	B	402	-	4,4,4	0.26	0	6,6,6	0.34	0
9	HEZ	B	403	-	7,7,7	0.32	0	6,6,6	1.14	0
10	M76	D	101	7	23,26,26	1.56	3 (13%)	25,37,37	1.04	3 (12%)

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	SO4	A	402	-	-	0/0/0/0	0/0/0/0
8	GOL	A	406	-	-	0/4/4/4	0/0/0/0
8	GOL	A	407	-	-	0/4/4/4	0/0/0/0
8	GOL	A	408	-	-	0/4/4/4	0/0/0/0
8	GOL	B	401	-	-	0/4/4/4	0/0/0/0
5	SO4	B	402	-	-	0/0/0/0	0/0/0/0
9	HEZ	B	403	-	-	0/5/5/5	0/0/0/0
10	M76	D	101	7	-	0/9/9/9	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	D	101	M76	C12-C2	2.74	1.55	1.51
10	D	101	M76	C12-C24	3.67	1.50	1.43
10	D	101	M76	C17-C19	4.07	1.48	1.41

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	D	101	M76	C23-C24-C19	-2.17	116.20	119.39
10	D	101	M76	O1-C2-C12	2.01	123.72	120.80
10	D	101	M76	C20-C19-C24	2.39	121.61	118.59

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	407	GOL	1	0
8	B	401	GOL	1	0
10	D	101	M76	6	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, 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	368/395 (93%)	0.16	11 (2%) 54 49	80, 104, 154, 194	1 (0%)
1	B	184/395 (46%)	0.13	11 (5%) 25 23	90, 120, 207, 230	0
2	C	19/19 (100%)	-0.09	1 (5%) 30 27	85, 107, 136, 158	0
3	D	17/17 (100%)	-0.11	0 100 100	91, 103, 142, 166	0
All	All	588/826 (71%)	0.14	23 (3%) 43 38	80, 107, 178, 230	1 (0%)

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	8	LEU	7.7
1	B	298	ILE	4.5
1	B	299	ARG	4.4
1	A	37	VAL	3.8
1	A	36	LYS	3.5
1	B	286	LEU	3.4
1	B	287	THR	3.3
1	A	15	LEU	3.3
1	A	20	TYR	3.3
1	B	296	GLN	3.2
1	B	293	SER	3.2
1	B	290	GLU	3.0
1	B	295	LEU	3.0
1	A	21	ILE	3.0
1	B	288	ARG	2.5
1	A	318	VAL	2.5
1	A	38	LYS	2.4
1	A	32	LEU	2.3
1	A	12	LEU	2.2
1	A	9	ASP	2.2
1	B	211	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	294	LEU	2.1
2	C	1	DA	2.1

## 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](#)

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(Å <sup>2</sup> )	Q<0.9
8	GOL	B	401	6/6	0.82	0.97	12.99	116,124,135,142	0
9	HEZ	B	403	8/8	0.72	0.40	5.97	124,132,141,142	0
8	GOL	A	408	6/6	0.70	0.36	2.64	124,132,143,151	0
4	ZN	A	401	1/1	0.99	0.17	0.32	94,94,94,94	0
10	M76	D	101	24/24	0.95	0.22	0.25	64,81,90,100	0
5	SO4	B	402	5/5	0.91	0.12	-0.57	141,147,160,169	0
8	GOL	A	406	6/6	0.89	0.15	-0.91	109,124,132,133	0
7	MG	A	405	1/1	0.99	0.16	-1.27	77,77,77,77	0
7	MG	A	404	1/1	0.96	0.23	-	95,95,95,95	0
8	GOL	A	407	6/6	0.93	0.15	-	109,116,126,151	0
5	SO4	A	402	5/5	0.93	0.17	-	135,136,148,156	0
6	NH4	A	403	1/1	0.49	0.46	-	57,57,57,57	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.