



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 07:14 PM GMT

PDB ID : 4RLD  
Title : Crystal structure of kkf mutant of bla G 2 protein  
Authors : Li, M.; Gustchina, A.; Pomes, A.; Wlodawer, A.  
Deposited on : 2014-10-16  
Resolution : 2.90 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 : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026982  
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) : rb-20026982

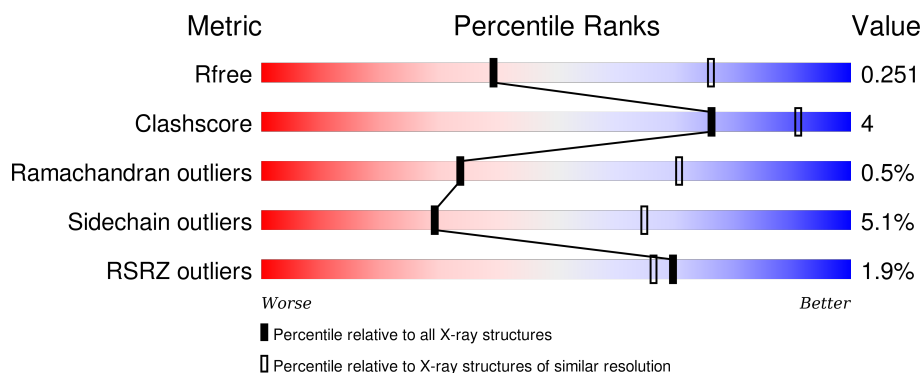
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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	330	<div> <div>2%</div> <div>87%</div> <div>12%</div> </div>
1	B	330	<div> <div>2%</div> <div>89%</div> <div>10%</div> </div>
1	C	330	<div> <div>%</div> <div>90%</div> <div>10%</div> </div>
1	D	330	<div> <div>3%</div> <div>79%</div> <div>19%</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	NAG	D	401	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10320 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 Aspartic protease Bla g 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	3	0
			2542	1620	418	492	12			
1	B	330	Total	C	N	O	S	0	3	0
			2542	1620	418	492	12			
1	C	330	Total	C	N	O	S	0	3	0
			2542	1620	418	492	12			
1	D	330	Total	C	N	O	S	0	3	0
			2542	1620	418	492	12			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	GLY	-	EXPRESSION TAG	UNP P54958
A	-7	ALA	-	EXPRESSION TAG	UNP P54958
A	-6	SER	-	EXPRESSION TAG	UNP P54958
A	-5	ILE	-	EXPRESSION TAG	UNP P54958
A	93	GLN	ASN	CONFLICT	UNP P54958
A	132	ALA	LYS	ENGINEERED MUTATION	UNP P54958
A	162	TYR	PHE	ENGINEERED MUTATION	UNP P54958
A	251	ALA	LYS	ENGINEERED MUTATION	UNP P54958
B	-8	GLY	-	EXPRESSION TAG	UNP P54958
B	-7	ALA	-	EXPRESSION TAG	UNP P54958
B	-6	SER	-	EXPRESSION TAG	UNP P54958
B	-5	ILE	-	EXPRESSION TAG	UNP P54958
B	93	GLN	ASN	ENGINEERED MUTATION	UNP P54958
B	132	ALA	LYS	ENGINEERED MUTATION	UNP P54958
B	162	TYR	PHE	ENGINEERED MUTATION	UNP P54958
B	251	ALA	LYS	ENGINEERED MUTATION	UNP P54958
C	-8	GLY	-	EXPRESSION TAG	UNP P54958
C	-7	ALA	-	EXPRESSION TAG	UNP P54958
C	-6	SER	-	EXPRESSION TAG	UNP P54958
C	-5	ILE	-	EXPRESSION TAG	UNP P54958
C	93	GLN	ASN	CONFLICT	UNP P54958

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Chain	Residue	Modelled	Actual	Comment	Reference
C	132	ALA	LYS	ENGINEERED MUTATION	UNP P54958
C	162	TYR	PHE	ENGINEERED MUTATION	UNP P54958
C	251	ALA	LYS	ENGINEERED MUTATION	UNP P54958
D	-8	GLY	-	EXPRESSION TAG	UNP P54958
D	-7	ALA	-	EXPRESSION TAG	UNP P54958
D	-6	SER	-	EXPRESSION TAG	UNP P54958
D	-5	ILE	-	EXPRESSION TAG	UNP P54958
D	93	GLN	ASN	CONFLICT	UNP P54958
D	132	ALA	LYS	ENGINEERED MUTATION	UNP P54958
D	162	TYR	PHE	ENGINEERED MUTATION	UNP P54958
D	251	ALA	LYS	ENGINEERED MUTATION	UNP P54958

- Molecule 2 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	2	Total	C	N	O	0	0
			28	16	2	10		
2	B	2	Total	C	N	O	0	0
			28	16	2	10		
2	C	2	Total	C	N	O	0	0
			28	16	2	10		
2	D	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		
3	A	1	Total	Zn	0	0
			1	1		
3	D	1	Total	Zn	0	0
			1	1		
3	C	1	Total	Zn	0	0
			1	1		

- Molecule 4 is water.

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

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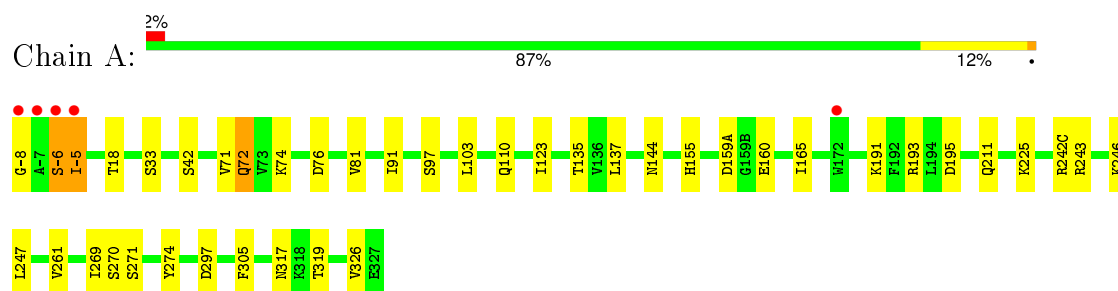
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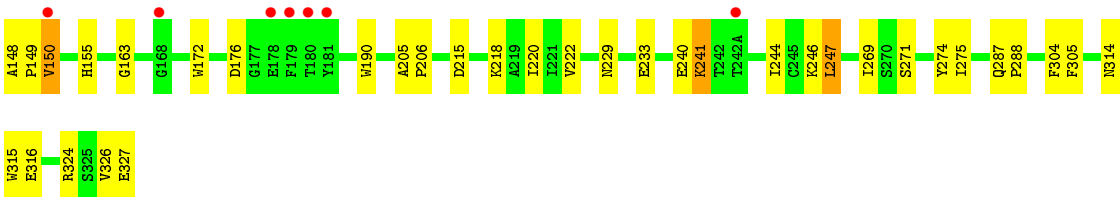
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	13	Total 13	O 13	0	0
4	C	7	Total 7	O 7	0	0
4	D	4	Total 4	O 4	0	0

### 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: Aspartic protease Bla g 2







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.82Å 75.41Å 339.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.90 49.48 – 2.90	Depositor EDS
% Data completeness (in resolution range)	88.0 (50.00-2.90) 88.0 (49.48-2.90)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.72 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.196 , 0.257 0.191 , 0.251	Depositor DCC
$R_{free}$ test set	1074 reflections (3.21%)	DCC
Wilson B-factor (Å <sup>2</sup> )	50.5	Xtriage
Anisotropy	0.233	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 41.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 34481 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	10320	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.57% 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

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

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	0.48	1/2618 (0.0%)	0.58	0/3559
1	B	0.46	0/2618	0.58	0/3559
1	C	0.46	1/2618 (0.0%)	0.57	0/3559
1	D	0.52	2/2618 (0.1%)	0.58	0/3559
All	All	0.48	4/10472 (0.0%)	0.58	0/14236

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	138	GLU	CB-CG	6.40	1.64	1.52
1	D	172	TRP	CD2-CE3	6.20	1.49	1.40
1	C	-6	SER	CB-OG	5.12	1.49	1.42
1	A	-8	GLY	N-CA	5.10	1.53	1.46

There are no bond angle outliers.

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	2542	0	2454	13	0
1	B	2542	0	2454	20	0
1	C	2542	0	2454	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2542	0	2454	36	0
2	A	28	0	25	0	0
2	B	28	0	25	0	0
2	C	28	0	25	0	0
2	D	28	0	25	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	12	0	0	1	0
4	B	13	0	0	2	0
4	C	7	0	0	0	0
4	D	4	0	0	0	0
All	All	10320	0	9916	76	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 4.

All (76) 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:D:150:VAL:HG12	1:D:314:ASN:HA	1.53	0.89
1:B:29:THR:HG21	1:B:121:VAL:HG23	1.62	0.81
1:D:229:ASN:O	1:D:233:GLU:HG2	1.88	0.73
1:D:149:PRO:HG2	1:D:316:GLU:HB2	1.71	0.72
1:D:72:GLN:HG2	1:D:81:VAL:HG22	1.73	0.71
1:B:29:THR:HG23	1:B:119:VAL:HG12	1.73	0.70
1:A:-5:ILE:HD13	1:A:144:ASN:HD21	1.58	0.68
1:A:269:ILE:HD11	1:A:305:PHE:HD1	1.61	0.64
1:B:29:THR:HG21	1:B:121:VAL:CG2	2.27	0.62
1:D:326:VAL:O	1:D:327:GLU:HB2	1.99	0.61
1:B:242:THR:HG23	4:B:511:HOH:O	2.00	0.61
1:D:269:ILE:HD11	1:D:305:PHE:HD1	1.64	0.61
1:A:317:ASN:O	1:A:319:THR:HG23	2.01	0.60
1:B:269:ILE:HD11	1:B:305:PHE:HD1	1.67	0.59
1:A:195:ASP:HB2	1:A:261:VAL:HG12	1.86	0.58
1:B:191:LYS:HD2	1:B:211:GLN:HE21	1.69	0.57
1:A:193:ARG:HD2	4:A:601:HOH:O	2.04	0.56
1:D:51:VAL:O	1:D:51(B):PRO:HD3	2.05	0.56
1:D:215:ASP:OD2	1:D:218:LYS:HE2	2.07	0.55
1:C:222:VAL:HG22	1:C:287:GLN:HB3	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:4:VAL:HB	1:C:165:ILE:HG23	1.89	0.55
1:B:233:GLU:HA	1:B:233:GLU:OE1	2.07	0.54
1:C:33:SER:HA	1:C:123:ILE:HG13	1.90	0.53
1:D:74:LYS:HE2	1:D:76:ASP:O	2.09	0.53
1:D:73:VAL:HG11	1:D:102:VAL:HG21	1.90	0.53
1:D:121:VAL:HG13	1:D:136:VAL:HG21	1.91	0.53
1:B:29:THR:CG2	1:B:121:VAL:HG23	2.36	0.52
1:C:72:GLN:HG2	1:C:81:VAL:HG22	1.92	0.52
1:A:159(A):ASP:OD2	1:A:160:GLU:OE1	2.29	0.52
1:D:33:SER:HA	1:D:123:ILE:HG13	1.94	0.50
1:D:13:GLN:O	1:D:14:TYR:HB2	2.13	0.49
1:D:72:GLN:CG	1:D:81:VAL:HG22	2.42	0.48
1:D:32:ASP:OD2	1:D:35:SER:HB3	2.13	0.48
1:D:92[B]:SER:HB3	1:D:93:GLN:H	1.52	0.48
1:D:220:ILE:HA	1:D:304:PHE:CE1	2.49	0.47
1:C:29:THR:HG23	1:C:119:VAL:HG12	1.96	0.47
1:B:-5:ILE:HD12	1:B:144:ASN:HD21	1.79	0.47
1:A:271:SER:HA	1:A:274:TYR:CE2	2.50	0.46
1:B:237:CYS:HB3	1:B:247:LEU:HB3	1.98	0.46
1:C:18:THR:HG22	1:C:91:ILE:HG12	1.97	0.46
1:B:146:ILE:HD13	1:B:167:PHE:HB3	1.98	0.45
1:B:15:ALA:HA	1:B:29:THR:O	2.17	0.45
1:A:42:SER:HB3	1:A:103:LEU:HG	1.99	0.45
1:D:176:ASP:HB3	1:D:324:ARG:HB2	1.98	0.45
1:A:191:LYS:HD2	1:A:211:GLN:HE21	1.82	0.44
1:B:313:PHE:N	1:B:313:PHE:CD1	2.86	0.44
1:D:47:GLY:HA2	1:D:51:VAL:CG2	2.48	0.44
1:C:53:LEU:HD12	1:C:118:ASP:HA	2.00	0.44
1:B:221:ILE:HG12	1:B:301:ILE:HA	1.99	0.44
1:D:271:SER:HA	1:D:274:TYR:CE2	2.53	0.44
1:B:92[A]:SER:OG	1:B:93:GLN:NE2	2.51	0.43
1:D:155:HIS:HA	1:D:163:GLY:HA3	2.01	0.43
1:D:148:ALA:O	1:D:150:VAL:N	2.45	0.42
1:D:218:LYS:HD3	1:D:222:VAL:HG23	2.00	0.42
1:B:214:ILE:HD13	1:B:301:ILE:HB	2.01	0.42
1:B:243:ARG:HA	4:B:511:HOH:O	2.19	0.42
1:D:51(A):CYS:HA	1:D:51(B):PRO:HD2	1.92	0.42
1:D:30:VAL:HG23	1:D:117:ALA:CB	2.49	0.42
1:A:165:ILE:O	1:A:165:ILE:HG23	2.18	0.42
1:C:25:GLN:HB3	1:C:27:PHE:CE1	2.54	0.42
1:D:52:ASN:HD22	1:D:52:ASN:HA	1.68	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:269:ILE:HD11	1:D:305:PHE:CD1	2.51	0.41
1:A:33:SER:HA	1:A:123:ILE:HG13	2.02	0.41
1:B:110:GLN:O	1:B:114:ILE:HG12	2.20	0.41
1:D:149:PRO:O	1:D:315:TRP:HB3	2.21	0.41
1:A:72:GLN:HG2	1:A:81:VAL:HG22	2.03	0.41
1:D:287:GLN:HA	1:D:288:PRO:HD3	1.92	0.41
1:D:205:ALA:HA	1:D:206:PRO:HD3	1.90	0.40
1:D:124:ALA:HA	1:D:190:TRP:CD1	2.57	0.40
1:D:247:LEU:HD21	1:D:275:ILE:HD12	2.03	0.40
1:D:47:GLY:HA2	1:D:51:VAL:HG22	2.03	0.40
1:B:238:VAL:HG22	1:D:93:GLN:HG2	2.04	0.40
1:A:18:THR:HG22	1:A:91:ILE:HG12	2.02	0.40
1:D:176:ASP:HB2	1:D:326:VAL:HG23	2.03	0.40
1:B:195:ASP:HB2	1:B:261:VAL:HG12	2.04	0.40
1:D:241:LYS:O	1:D:241:LYS:HG3	2.21	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	331/330 (100%)	310 (94%)	20 (6%)	1 (0%)	46	79
1	B	331/330 (100%)	315 (95%)	14 (4%)	2 (1%)	30	67
1	C	331/330 (100%)	312 (94%)	16 (5%)	3 (1%)	21	57
1	D	331/330 (100%)	312 (94%)	18 (5%)	1 (0%)	46	79
All	All	1324/1320 (100%)	1249 (94%)	68 (5%)	7 (0%)	34	71

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	-6	SER
1	C	76	ASP
1	C	242(B)	THR
1	C	225	LYS
1	B	76	ASP
1	B	126	PRO
1	D	128	PRO

### 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	283/280 (101%)	264 (93%)	19 (7%)	20	50
1	B	283/280 (101%)	275 (97%)	8 (3%)	51	84
1	C	283/280 (101%)	270 (95%)	13 (5%)	33	69
1	D	283/280 (101%)	266 (94%)	17 (6%)	24	57
All	All	1132/1120 (101%)	1075 (95%)	57 (5%)	29	65

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-6	SER
1	A	-5	ILE
1	A	71	VAL
1	A	72	GLN
1	A	74	LYS
1	A	76	ASP
1	A	97	SER
1	A	110	GLN
1	A	135	THR
1	A	137	LEU
1	A	155	HIS
1	A	225	LYS
1	A	242(C)	ARG
1	A	243	ARG
1	A	246	LYS

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Mol	Chain	Res	Type
1	A	247	LEU
1	A	270	SER
1	A	297	ASP
1	A	326	VAL
1	B	-5	ILE
1	B	96	THR
1	B	179	PHE
1	B	242(C)	ARG
1	B	243	ARG
1	B	247	LEU
1	B	270	SER
1	B	287	GLN
1	C	51	VAL
1	C	57	GLU
1	C	134	LYS
1	C	137	LEU
1	C	155	HIS
1	C	244	ILE
1	C	246	LYS
1	C	248	ASP
1	C	269	ILE
1	C	270	SER
1	C	292	SER
1	C	297	ASP
1	C	316	GLU
1	D	-5	ILE
1	D	-2	LEU
1	D	0	LYS
1	D	17	ILE
1	D	35	SER
1	D	52	ASN
1	D	57	GLU
1	D	72	GLN
1	D	77	THR
1	D	95	THR
1	D	96	THR
1	D	150	VAL
1	D	240	GLU
1	D	241	LYS
1	D	244	ILE
1	D	246	LYS
1	D	247	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	ASN
1	A	52	ASN
1	A	98	GLN
1	A	144	ASN
1	A	211	GLN
1	B	26	ASN
1	B	52	ASN
1	B	93	GLN
1	B	98	GLN
1	B	211	GLN
1	B	287	GLN
1	C	26	ASN
1	C	52	ASN
1	C	93	GLN
1	C	98	GLN
1	C	159	GLN
1	C	211	GLN
1	C	272	GLN
1	C	278	ASN
1	D	52	ASN
1	D	110	GLN
1	D	159	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 ⓘ

8 carbohydrates are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected



value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	NAG	A	501	1,2	14,14,15	0.80	0	15,19,21	1.00	1 (6%)
2	NAG	A	502	2	14,14,15	0.45	0	15,19,21	1.03	0
2	NAG	B	401	1,2	14,14,15	0.57	0	15,19,21	1.42	1 (6%)
2	NAG	B	402	2	14,14,15	0.66	0	15,19,21	1.14	1 (6%)
2	NAG	C	401	1,2	14,14,15	0.52	0	15,19,21	1.25	2 (13%)
2	NAG	C	402	2	14,14,15	0.59	0	15,19,21	0.99	1 (6%)
2	NAG	D	401	1,2	14,14,15	0.57	0	15,19,21	1.17	1 (6%)
2	NAG	D	402	2	14,14,15	0.55	0	15,19,21	1.00	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
2	NAG	A	501	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	502	2	-	0/6/23/26	0/1/1/1
2	NAG	B	401	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	402	2	-	0/6/23/26	0/1/1/1
2	NAG	C	401	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	402	2	-	0/6/23/26	0/1/1/1
2	NAG	D	401	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	402	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NAG	C4-C3-C2	2.06	114.53	111.34
2	C	401	NAG	C4-C3-C2	2.40	115.07	111.34
2	C	402	NAG	C4-C3-C2	2.58	115.34	111.34
2	B	402	NAG	C4-C3-C2	3.17	116.27	111.34
2	D	401	NAG	C1-O5-C5	3.38	117.11	112.14
2	C	401	NAG	C1-O5-C5	3.64	117.49	112.14
2	B	401	NAG	C1-O5-C5	4.56	118.84	112.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.6 Ligand geometry

Of 4 ligands modelled in this entry, 4 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

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	330/330 (100%)	-0.21	5 (1%) 76 74	21, 39, 60, 109	0
1	B	330/330 (100%)	-0.28	5 (1%) 76 74	24, 42, 63, 112	0
1	C	330/330 (100%)	-0.11	4 (1%) 81 78	32, 51, 75, 130	0
1	D	330/330 (100%)	0.05	11 (3%) 50 42	31, 63, 94, 185	0
All	All	1320/1320 (100%)	-0.14	25 (1%) 70 66	21, 47, 82, 185	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	-8	GLY	9.8
1	D	-8	GLY	9.2
1	A	-8	GLY	7.8
1	C	-7	ALA	7.6
1	B	-7	ALA	6.8
1	B	-8	GLY	6.2
1	A	-7	ALA	6.1
1	D	242(A)	THR	5.1
1	C	-6	SER	4.8
1	D	-7	ALA	4.4
1	C	327	GLU	4.1
1	A	-6	SER	4.0
1	B	-6	SER	3.7
1	D	180	THR	3.7
1	D	-6	SER	3.0
1	B	-5	ILE	2.7
1	D	179	PHE	2.4
1	B	327	GLU	2.4
1	D	168	GLY	2.3
1	D	150	VAL	2.3
1	D	178	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	181	TYR	2.2
1	D	-5	ILE	2.2
1	A	172	TRP	2.1
1	A	-5	ILE	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](#)

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
2	NAG	D	401	14/15	0.94	0.30	2.95	57,64,67,73	0
2	NAG	C	401	14/15	0.94	0.25	0.80	54,61,65,71	0
2	NAG	B	401	14/15	0.93	0.14	0.43	51,57,64,70	0
2	NAG	A	501	14/15	0.95	0.13	-1.50	45,50,55,61	0
2	NAG	B	402	14/15	0.83	0.22	-	74,77,78,78	0
2	NAG	D	402	14/15	0.87	0.31	-	77,80,82,83	0
2	NAG	A	502	14/15	0.91	0.23	-	67,71,72,73	0
2	NAG	C	402	14/15	0.87	0.35	-	77,81,83,85	0

## 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
3	ZN	B	403	1/1	1.00	0.14	-0.17	35,35,35,35	0
3	ZN	A	503	1/1	1.00	0.15	-0.35	31,31,31,31	0
3	ZN	D	403	1/1	0.98	0.12	-1.05	51,51,51,51	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	ZN	C	403	1/1	1.00	0.09	-2.56	45,45,45,45	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.