



Full wwPDB X-ray Structure Validation Report

May 16, 2020 – 08:02 am BST

PDB ID : 6WHL
Title : The crystal structure of a beta-lactamase from Legionella pneumophila str. Paris
Authors : Tan, K.; Wu, R.; Endres, M.; Joachimiak, A.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2020-04-08
Resolution : 2.30 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the  symbol.

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 : 2.11
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) : 2.11

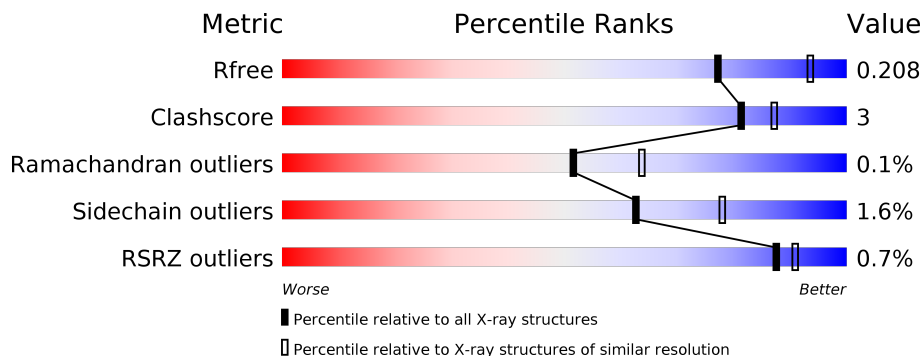
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 ≥ 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	249	
1	B	249	
1	C	249	
1	D	249	
1	E	249	
1	F	249	

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Mol	Chain	Length	Quality of chain
1	G	249	 89% 10% .
1	H	249	 % 96% ..
1	I	249	 92% 7% .
1	J	249	 % 91% 6% ..
1	K	249	 2% 88% 10% .
1	L	249	 % 90% 8% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 24649 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 Beta-lactamase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	246	2011	1291	342	371	3	4	0	0	0
1	B	246	2012	1292	341	372	3	4	0	0	0
1	C	246	2006	1288	340	371	3	4	0	0	0
1	D	246	1997	1285	340	365	3	4	0	0	0
1	E	246	2015	1294	343	371	3	4	0	0	0
1	F	246	1997	1285	338	367	3	4	0	0	0
1	G	246	1999	1287	340	365	3	4	0	0	0
1	H	246	2007	1290	341	369	3	4	0	0	0
1	I	246	2011	1293	341	370	3	4	0	0	0
1	J	246	1996	1283	339	367	3	4	0	0	0
1	K	246	2004	1287	340	370	3	4	0	0	0
1	L	246	1993	1282	338	366	3	4	0	0	0

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	18	SER	-	expression tag	UNP A0A378KDC5
A	19	ASN	-	expression tag	UNP A0A378KDC5
A	20	ALA	-	expression tag	UNP A0A378KDC5
A	22	ASP	GLY	conflict	UNP A0A378KDC5
B	18	SER	-	expression tag	UNP A0A378KDC5

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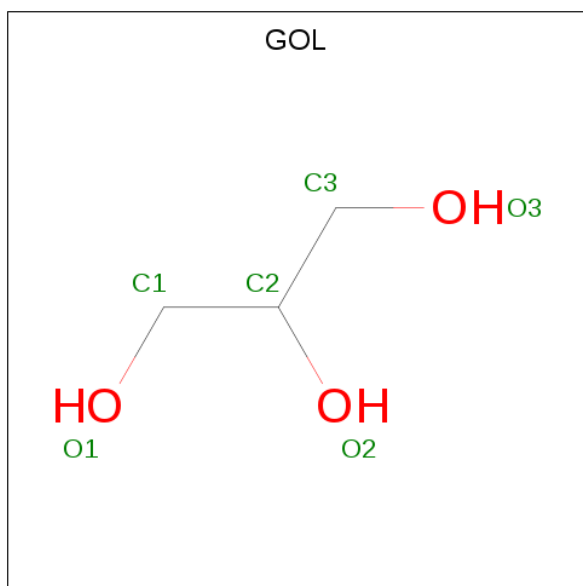
Chain	Residue	Modelled	Actual	Comment	Reference
B	19	ASN	-	expression tag	UNP A0A378KDC5
B	20	ALA	-	expression tag	UNP A0A378KDC5
B	22	ASP	GLY	conflict	UNP A0A378KDC5
C	18	SER	-	expression tag	UNP A0A378KDC5
C	19	ASN	-	expression tag	UNP A0A378KDC5
C	20	ALA	-	expression tag	UNP A0A378KDC5
C	22	ASP	GLY	conflict	UNP A0A378KDC5
D	18	SER	-	expression tag	UNP A0A378KDC5
D	19	ASN	-	expression tag	UNP A0A378KDC5
D	20	ALA	-	expression tag	UNP A0A378KDC5
D	22	ASP	GLY	conflict	UNP A0A378KDC5
E	18	SER	-	expression tag	UNP A0A378KDC5
E	19	ASN	-	expression tag	UNP A0A378KDC5
E	20	ALA	-	expression tag	UNP A0A378KDC5
E	22	ASP	GLY	conflict	UNP A0A378KDC5
F	18	SER	-	expression tag	UNP A0A378KDC5
F	19	ASN	-	expression tag	UNP A0A378KDC5
F	20	ALA	-	expression tag	UNP A0A378KDC5
F	22	ASP	GLY	conflict	UNP A0A378KDC5
G	18	SER	-	expression tag	UNP A0A378KDC5
G	19	ASN	-	expression tag	UNP A0A378KDC5
G	20	ALA	-	expression tag	UNP A0A378KDC5
G	22	ASP	GLY	conflict	UNP A0A378KDC5
H	18	SER	-	expression tag	UNP A0A378KDC5
H	19	ASN	-	expression tag	UNP A0A378KDC5
H	20	ALA	-	expression tag	UNP A0A378KDC5
H	22	ASP	GLY	conflict	UNP A0A378KDC5
I	18	SER	-	expression tag	UNP A0A378KDC5
I	19	ASN	-	expression tag	UNP A0A378KDC5
I	20	ALA	-	expression tag	UNP A0A378KDC5
I	22	ASP	GLY	conflict	UNP A0A378KDC5
J	18	SER	-	expression tag	UNP A0A378KDC5
J	19	ASN	-	expression tag	UNP A0A378KDC5
J	20	ALA	-	expression tag	UNP A0A378KDC5
J	22	ASP	GLY	conflict	UNP A0A378KDC5
K	18	SER	-	expression tag	UNP A0A378KDC5
K	19	ASN	-	expression tag	UNP A0A378KDC5
K	20	ALA	-	expression tag	UNP A0A378KDC5
K	22	ASP	GLY	conflict	UNP A0A378KDC5
L	18	SER	-	expression tag	UNP A0A378KDC5
L	19	ASN	-	expression tag	UNP A0A378KDC5
L	20	ALA	-	expression tag	UNP A0A378KDC5

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Chain	Residue	Modelled	Actual	Comment	Reference
L	22	ASP	GLY	conflict	UNP A0A378KDC5

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	E	1	Total	C	O	0	0
			6	3	3		
2	F	1	Total	C	O	0	0
			6	3	3		
2	G	1	Total	C	O	0	0
			6	3	3		
2	H	1	Total	C	O	0	0
			6	3	3		
2	I	1	Total	C	O	0	0
			6	3	3		
2	J	1	Total	C	O	0	0
			6	3	3		
2	K	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	L	1	6	3	3	0	0


- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	49	Total 49	O 49	0	0
3	B	30	Total 30	O 30	0	0
3	C	23	Total 23	O 23	0	0
3	D	54	Total 54	O 54	0	0
3	E	52	Total 52	O 52	0	0
3	F	19	Total 19	O 19	0	0
3	G	81	Total 81	O 81	0	0
3	H	61	Total 61	O 61	0	0
3	I	68	Total 68	O 68	0	0
3	J	54	Total 54	O 54	0	0
3	K	28	Total 28	O 28	0	0
3	L	10	Total 10	O 10	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 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: Beta-lactamase

Chain A: 

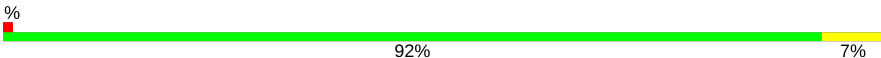


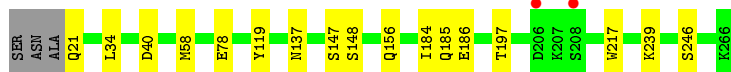
- Molecule 1: Beta-lactamase

Chain B: 




- Molecule 1: Beta-lactamase

Chain C: 




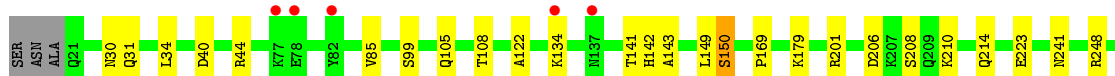
- Molecule 1: Beta-lactamase

Chain D: 

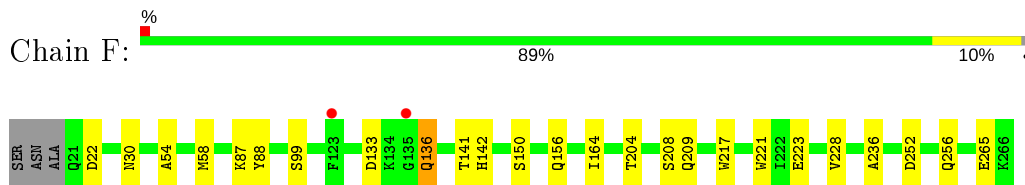


- Molecule 1: Beta-lactamase

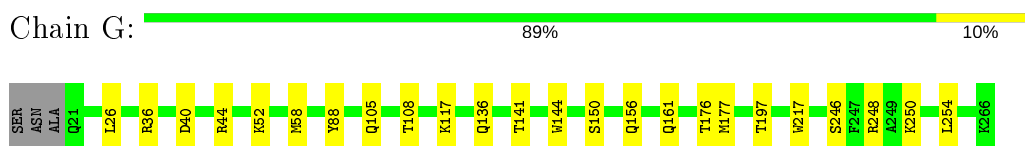
Chain E: 



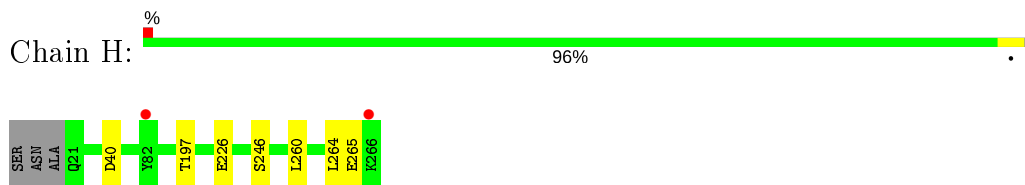
• Molecule 1: Beta-lactamase



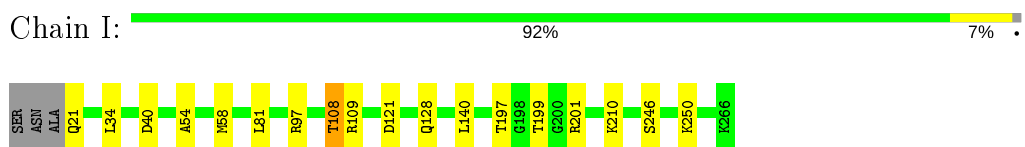
• Molecule 1: Beta-lactamase



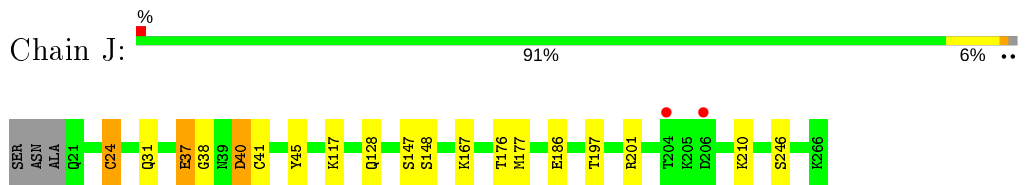
• Molecule 1: Beta-lactamase



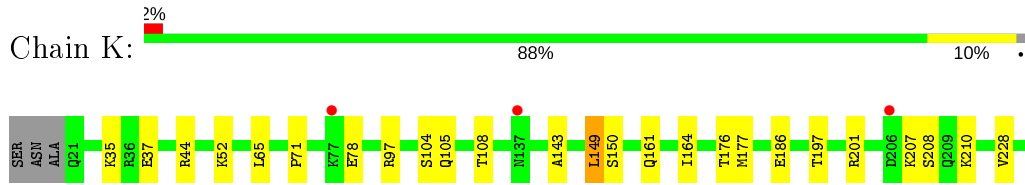
• Molecule 1: Beta-lactamase



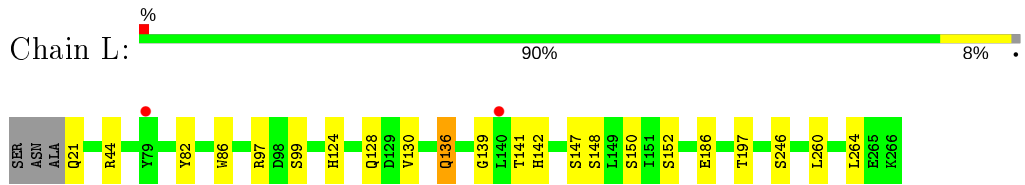
• Molecule 1: Beta-lactamase



• Molecule 1: Beta-lactamase



• Molecule 1: Beta-lactamase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	80.22Å 178.20Å 143.75Å 90.00° 97.04° 90.00°	Depositor
Resolution (Å)	42.52 – 2.30 47.83 – 2.29	Depositor EDS
% Data completeness (in resolution range)	87.9 (42.52-2.30) 83.1 (47.83-2.29)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.58 (at 2.29Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.163 , 0.208 0.163 , 0.208	Depositor DCC
R_{free} test set	7619 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	27.3	Xtrriage
Anisotropy	0.058	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 41.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	24649	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.53% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: GOL, KCX

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	0.38	0/2047	0.56	0/2761
1	B	0.39	0/2048	0.56	0/2762
1	C	0.37	0/2042	0.54	0/2756
1	D	0.41	0/2033	0.56	0/2743
1	E	0.41	0/2051	0.55	0/2765
1	F	0.34	0/2033	0.52	0/2743
1	G	0.41	0/2035	0.57	0/2745
1	H	0.41	0/2043	0.57	0/2756
1	I	0.43	0/2047	0.58	0/2760
1	J	0.42	1/2032 (0.0%)	0.57	0/2743
1	K	0.36	0/2040	0.52	0/2753
1	L	0.33	0/2029	0.49	0/2740
All	All	0.39	1/24480 (0.0%)	0.55	0/33027

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	167	LYS	C-N	-5.00	1.22	1.34

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	2011	0	1966	13	0
1	B	2012	0	1966	8	0
1	C	2006	0	1948	9	0
1	D	1997	0	1948	12	0
1	E	2015	0	1975	16	0
1	F	1997	0	1946	12	0
1	G	1999	0	1952	16	0
1	H	2007	0	1959	3	0
1	I	2011	0	1966	11	0
1	J	1996	0	1941	11	0
1	K	2004	0	1949	14	0
1	L	1993	0	1931	11	0
2	A	6	0	8	0	0
2	B	6	0	8	0	0
2	C	12	0	16	0	0
2	E	6	0	8	0	0
2	F	6	0	8	1	0
2	G	6	0	8	1	0
2	H	6	0	8	1	0
2	I	6	0	8	1	0
2	J	6	0	8	0	0
2	K	6	0	8	0	0
2	L	6	0	8	0	0
3	A	49	0	0	0	0
3	B	30	0	0	0	0
3	C	23	0	0	0	0
3	D	54	0	0	1	0
3	E	52	0	0	0	0
3	F	19	0	0	0	0
3	G	81	0	0	4	0
3	H	61	0	0	0	0
3	I	68	0	0	0	0
3	J	54	0	0	1	0
3	K	28	0	0	0	0
3	L	10	0	0	0	0
All	All	24649	0	23543	123	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 3.

All (123) 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:40:ASP:OD2	1:D:45:TYR:OH	2.03	0.73
1:D:147:SER:OG	1:D:148:SER:N	2.23	0.71
1:H:197:THR:HB	1:H:246:SER:HB2	1.80	0.63
1:B:197:THR:HB	1:B:246:SER:HB2	1.80	0.63
1:F:223:GLU:HG2	1:F:228:VAL:HG22	1.79	0.62
1:C:197:THR:HB	1:C:246:SER:HB2	1.81	0.61
1:G:136:GLN:NE2	3:G:401:HOH:O	2.32	0.61
1:J:197:THR:HB	1:J:246:SER:HB2	1.82	0.61
1:F:141:THR:HG23	1:F:142:HIS:ND1	2.17	0.60
1:A:133:ASP:OD1	1:A:147:SER:OG	2.20	0.59
1:B:36:ARG:HD3	1:C:34:LEU:O	2.04	0.58
1:I:108:THR:HG23	1:I:140:LEU:HG	1.85	0.58
1:E:85:VAL:HG12	1:F:265:GLU:HG3	1.85	0.58
1:D:58:MSE:HG2	1:D:119:TYR:CG	2.39	0.58
1:I:197:THR:HB	1:J:246:SER:HB2	1.87	0.57
1:E:105:GLN:HA	1:E:108:THR:HG22	1.86	0.56
1:C:184:ILE:HG22	1:C:185:GLN:HG2	1.87	0.56
1:F:87:LYS:HE2	1:F:88:TYR:CE2	2.41	0.56
1:D:197:THR:HB	1:D:246:SER:HB2	1.88	0.56
1:D:161:GLN:HG2	1:D:230:THR:HG21	1.89	0.54
1:B:141:THR:HG23	1:B:142:HIS:ND1	2.23	0.54
1:E:141:THR:HG23	1:E:142:HIS:ND1	2.22	0.54
1:A:121:ASP:OD1	1:A:128:GLN:NE2	2.42	0.53
1:E:201:ARG:HH21	1:E:210:LYS:HD2	1.74	0.53
1:G:36:ARG:HD3	1:I:34:LEU:O	2.08	0.53
1:G:176:THR:HG22	1:G:177:MSE:HE2	1.89	0.53
1:A:186:GLU:OE2	1:B:97:ARG:NH1	2.35	0.53
1:K:97:ARG:NH1	1:L:186:GLU:OE2	2.33	0.53
1:B:24:CYS:HG	1:B:41:CYS:HG	1.57	0.52
1:L:197:THR:HB	1:L:246:SER:HB2	1.91	0.52
1:G:88:TYR:OH	1:H:265:GLU:OE1	2.16	0.52
1:J:176:THR:HG22	1:J:177:MSE:HE2	1.92	0.51
1:J:24:CYS:HG	1:J:41:CYS:HG	1.57	0.51
1:A:204:THR:O	1:A:208:SER:HA	2.10	0.51
1:J:117:LYS:NZ	1:J:128:GLN:O	2.43	0.51
1:J:31:GLN:HB2	3:J:438:HOH:O	2.09	0.51
1:I:97:ARG:NH1	1:J:186:GLU:OE2	2.37	0.51
1:G:250:LYS:HD3	2:G:301:GOL:H32	1.93	0.50
1:J:38:GLY:O	1:K:35:LYS:NZ	2.44	0.50
1:K:44:ARG:HB3	1:K:150:SER:OG	2.12	0.50
1:G:197:THR:HB	1:G:246:SER:HB2	1.93	0.49
1:K:143:ALA:O	1:K:149:LEU:HB2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:141:THR:HG23	1:D:142:HIS:ND1	2.27	0.49
1:J:147:SER:OG	1:J:148:SER:N	2.45	0.49
1:K:201:ARG:HH21	1:K:210:LYS:HD2	1.79	0.48
1:E:179:LYS:NZ	1:E:223:GLU:OE1	2.46	0.48
1:K:164:ILE:HD12	1:K:228:VAL:HG11	1.95	0.48
1:J:201:ARG:HH21	1:J:210:LYS:HD2	1.78	0.47
1:L:141:THR:HG23	1:L:142:HIS:ND1	2.28	0.47
1:E:143:ALA:O	1:E:149:LEU:HB2	2.14	0.47
1:I:121:ASP:OD1	1:I:128:GLN:NE2	2.43	0.47
1:K:197:THR:HB	1:K:246:SER:HB2	1.97	0.47
1:L:44:ARG:HB3	1:L:150:SER:OG	2.15	0.47
1:G:52:KCX:OQ2	1:G:144:TRP:NE1	2.41	0.47
1:F:22:ASP:HB2	1:F:236:ALA:HB3	1.96	0.47
1:A:256:GLN:H	1:A:256:GLN:HG2	1.60	0.45
1:A:197:THR:HB	1:A:246:SER:HB2	1.97	0.45
1:K:105:GLN:HA	1:K:108:THR:HG22	1.97	0.45
1:G:26:LEU:CD2	1:G:36:ARG:HG3	2.47	0.45
1:E:214:GLN:NE2	1:E:241:ASN:O	2.50	0.44
1:L:82:TYR:HB3	1:L:86:TRP:CZ3	2.52	0.44
1:G:26:LEU:HD21	1:G:36:ARG:HG3	1.99	0.44
1:H:260:LEU:O	1:H:264:LEU:HG	2.18	0.44
1:B:26:LEU:O	1:B:231:PHE:HA	2.18	0.44
1:C:186:GLU:OE2	1:D:97:ARG:NH1	2.45	0.44
1:K:186:GLU:OE2	1:L:97:ARG:NH1	2.45	0.44
1:D:75:TYR:CZ	1:D:81:LEU:HD13	2.53	0.43
1:E:134:LYS:HG3	1:E:208:SER:HA	1.99	0.43
1:C:156:GLN:HG3	1:C:217:TRP:CZ3	2.52	0.43
1:I:250:LYS:NZ	2:I:301:GOL:H11	2.33	0.43
1:G:117:LYS:HD2	3:G:464:HOH:O	2.17	0.43
1:K:78:GLU:HG3	1:K:78:GLU:H	1.64	0.43
1:E:254:LEU:HD13	2:F:301:GOL:H2	2.00	0.43
1:E:44:ARG:HB3	1:E:150:SER:HB3	2.00	0.43
1:I:54:ALA:O	1:I:58:MSE:HG3	2.19	0.43
1:D:36:ARG:HD3	1:E:34:LEU:O	2.18	0.43
1:F:204:THR:O	1:F:208:SER:HA	2.19	0.43
1:G:156:GLN:HG3	1:G:217:TRP:CZ3	2.53	0.43
1:A:251:ASN:ND2	1:B:251:ASN:HB3	2.34	0.42
1:G:248:ARG:HD2	3:G:418:HOH:O	2.18	0.42
1:C:58:MSE:HG2	1:C:119:TYR:CG	2.53	0.42
1:D:164:ILE:HD11	1:D:221:TRP:CD2	2.54	0.42
1:G:117:LYS:HE2	3:G:471:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:GLN:HG3	1:A:217:TRP:CZ3	2.54	0.42
1:I:108:THR:HG21	1:I:140:LEU:O	2.20	0.42
1:C:147:SER:OG	1:C:148:SER:N	2.51	0.42
1:E:206:ASP:N	1:E:206:ASP:OD1	2.49	0.42
1:A:266:LYS:HB3	1:A:266:LYS:HE3	1.79	0.42
1:I:201:ARG:HH21	1:I:210:LYS:HD2	1.84	0.42
1:C:239:LYS:HE2	1:C:239:LYS:HB3	1.79	0.42
1:E:30:ASN:OD1	1:E:31:GLN:HG2	2.19	0.42
1:B:157:ILE:O	1:B:161:GLN:HG3	2.20	0.42
1:A:44:ARG:HB3	1:A:150:SER:OG	2.20	0.42
1:F:133:ASP:HB2	1:F:136:GLN:HG3	2.02	0.42
1:L:147:SER:OG	1:L:148:SER:N	2.52	0.42
1:F:252:ASP:O	1:F:256:GLN:HG3	2.20	0.41
1:G:105:GLN:HA	1:G:108:THR:HG22	2.01	0.41
1:A:147:SER:HB3	1:A:148:SER:H	1.47	0.41
1:D:214:GLN:HG3	3:D:321:HOH:O	2.20	0.41
1:F:221:TRP:CD1	1:F:223:GLU:HG3	2.55	0.41
1:I:81:LEU:HA	1:I:81:LEU:HD12	1.95	0.41
1:A:164:ILE:HD11	1:A:221:TRP:CD2	2.55	0.41
1:J:40:ASP:OD2	1:J:45:TYR:OH	2.16	0.41
1:E:134:LYS:HB3	1:E:134:LYS:HE2	1.94	0.41
1:F:156:GLN:HG3	1:F:217:TRP:CZ3	2.56	0.41
1:J:37:GLU:HG2	1:K:37:GLU:HG2	2.03	0.41
1:L:124:HIS:HA	1:L:128:GLN:NE2	2.36	0.41
1:L:136:GLN:HG2	1:L:136:GLN:H	1.60	0.41
1:F:54:ALA:O	1:F:58:MSE:HG3	2.21	0.41
1:L:130:VAL:O	1:L:139:GLY:HA3	2.21	0.41
1:D:143:ALA:O	1:D:149:LEU:HB2	2.20	0.41
1:E:122:ALA:HB1	1:E:169:PRO:HG3	2.03	0.41
1:G:254:LEU:HB3	2:H:301:GOL:H2	2.02	0.41
1:E:248:ARG:HD2	1:E:248:ARG:HH21	1.76	0.41
1:F:164:ILE:HD11	1:F:221:TRP:CD2	2.56	0.41
1:I:109:ARG:HA	1:I:140:LEU:HD21	2.03	0.41
1:G:44:ARG:HB3	1:G:150:SER:OG	2.21	0.41
1:K:52:KCX:HE2	1:K:104:SER:OG	2.20	0.40
1:L:260:LEU:O	1:L:264:LEU:HG	2.21	0.40
1:K:65:LEU:HA	1:K:71:PRO:HD3	2.03	0.40
1:A:203:LEU:HA	1:A:209:GLN:O	2.21	0.40
1:C:21:GLN:OE1	1:C:21:GLN:N	2.55	0.40
1:K:176:THR:HG22	1:K:177:MSE:HE2	2.03	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	243/249 (98%)	232 (96%)	11 (4%)	0	100	100
1	B	243/249 (98%)	234 (96%)	9 (4%)	0	100	100
1	C	243/249 (98%)	234 (96%)	9 (4%)	0	100	100
1	D	243/249 (98%)	229 (94%)	14 (6%)	0	100	100
1	E	243/249 (98%)	231 (95%)	12 (5%)	0	100	100
1	F	243/249 (98%)	232 (96%)	10 (4%)	1 (0%)	34	42
1	G	243/249 (98%)	237 (98%)	6 (2%)	0	100	100
1	H	243/249 (98%)	235 (97%)	8 (3%)	0	100	100
1	I	243/249 (98%)	230 (95%)	13 (5%)	0	100	100
1	J	243/249 (98%)	235 (97%)	8 (3%)	0	100	100
1	K	243/249 (98%)	233 (96%)	9 (4%)	1 (0%)	34	42
1	L	243/249 (98%)	227 (93%)	16 (7%)	0	100	100
All	All	2916/2988 (98%)	2789 (96%)	125 (4%)	2 (0%)	51	64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	K	207	LYS
1	F	30	ASN

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	215/216 (100%)	211 (98%)	4 (2%)	57	73
1	B	215/216 (100%)	211 (98%)	4 (2%)	57	73
1	C	213/216 (99%)	210 (99%)	3 (1%)	67	81
1	D	211/216 (98%)	207 (98%)	4 (2%)	57	73
1	E	216/216 (100%)	213 (99%)	3 (1%)	67	81
1	F	211/216 (98%)	207 (98%)	4 (2%)	57	73
1	G	211/216 (98%)	207 (98%)	4 (2%)	57	73
1	H	213/216 (99%)	211 (99%)	2 (1%)	78	89
1	I	214/216 (99%)	210 (98%)	4 (2%)	57	73
1	J	211/216 (98%)	208 (99%)	3 (1%)	67	81
1	K	213/216 (99%)	210 (99%)	3 (1%)	67	81
1	L	209/216 (97%)	205 (98%)	4 (2%)	57	73
All	All	2552/2592 (98%)	2510 (98%)	42 (2%)	62	78

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	ASP
1	A	136	GLN
1	A	238	SER
1	A	256	GLN
1	B	24	CYS
1	B	40	ASP
1	B	78	GLU
1	B	99	SER
1	C	40	ASP
1	C	78	GLU
1	C	137	ASN
1	D	40	ASP
1	D	99	SER
1	D	147	SER
1	D	150	SER
1	E	40	ASP
1	E	99	SER
1	E	150	SER
1	F	99	SER
1	F	136	GLN
1	F	150	SER
1	F	209	GLN

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Mol	Chain	Res	Type
1	G	40	ASP
1	G	58	MSE
1	G	141	THR
1	G	161	GLN
1	H	40	ASP
1	H	226	GLU
1	I	21	GLN
1	I	40	ASP
1	I	108	THR
1	I	199	THR
1	J	24	CYS
1	J	37	GLU
1	J	40	ASP
1	K	149	LEU
1	K	161	GLN
1	K	208	SER
1	L	21	GLN
1	L	99	SER
1	L	136	GLN
1	L	152	SER

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

Mol	Chain	Res	Type
1	C	21	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

12 non-standard protein/DNA/RNA residues 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
1	KCX	K	52	1	7,11,12	0.94	1 (14%)	4,12,14	0.38	0
1	KCX	A	52	1	7,11,12	0.73	0	4,12,14	1.32	1 (25%)
1	KCX	E	52	1	7,11,12	0.75	0	4,12,14	0.51	0
1	KCX	C	52	1	7,11,12	0.93	0	4,12,14	0.45	0
1	KCX	G	52	1	7,11,12	0.91	0	4,12,14	1.49	1 (25%)
1	KCX	J	52	1	7,11,12	0.77	0	4,12,14	1.34	1 (25%)
1	KCX	H	52	1	7,11,12	0.98	0	4,12,14	1.77	1 (25%)
1	KCX	L	52	1	7,11,12	0.92	0	4,12,14	0.91	0
1	KCX	B	52	1	7,11,12	0.69	0	4,12,14	1.63	1 (25%)
1	KCX	F	52	1	7,11,12	0.83	0	4,12,14	0.61	0
1	KCX	I	52	1	7,11,12	0.68	0	4,12,14	2.34	1 (25%)
1	KCX	D	52	1	7,11,12	1.20	1 (14%)	4,12,14	0.60	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
1	KCX	K	52	1	-	0/7/10/12	-
1	KCX	A	52	1	-	3/7/10/12	-
1	KCX	E	52	1	-	0/7/10/12	-
1	KCX	C	52	1	-	0/7/10/12	-
1	KCX	G	52	1	-	2/7/10/12	-
1	KCX	J	52	1	-	2/7/10/12	-
1	KCX	H	52	1	-	3/7/10/12	-
1	KCX	L	52	1	-	0/7/10/12	-
1	KCX	B	52	1	-	3/7/10/12	-
1	KCX	F	52	1	-	0/7/10/12	-
1	KCX	I	52	1	-	2/7/10/12	-
1	KCX	D	52	1	-	0/7/10/12	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	52	KCX	CE-NZ	2.47	1.51	1.45
1	K	52	KCX	CE-NZ	2.01	1.50	1.45

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	I	52	KCX	CE-NZ-CX	4.25	130.15	122.95
1	H	52	KCX	CE-NZ-CX	3.04	128.09	122.95
1	B	52	KCX	CD-CE-NZ	-2.64	104.40	111.49
1	J	52	KCX	CD-CE-NZ	-2.48	104.83	111.49
1	G	52	KCX	CD-CE-NZ	-2.37	105.13	111.49
1	A	52	KCX	CD-CE-NZ	-2.23	105.50	111.49

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	G	52	KCX	CG-CD-CE-NZ
1	H	52	KCX	CG-CD-CE-NZ
1	H	52	KCX	CE-CD-CG-CB
1	B	52	KCX	CG-CD-CE-NZ
1	A	52	KCX	CG-CD-CE-NZ
1	B	52	KCX	CE-CD-CG-CB
1	I	52	KCX	CG-CD-CE-NZ
1	J	52	KCX	CG-CD-CE-NZ
1	I	52	KCX	CE-CD-CG-CB
1	J	52	KCX	CE-CD-CG-CB
1	A	52	KCX	CE-CD-CG-CB
1	A	52	KCX	CD-CE-NZ-CX
1	G	52	KCX	CD-CE-NZ-CX
1	H	52	KCX	CD-CE-NZ-CX
1	B	52	KCX	CD-CE-NZ-CX

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	K	52	KCX	1	0
1	G	52	KCX	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

12 ligands 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	GOL	H	301	-	5,5,5	0.86	0	5,5,5	1.04	0
2	GOL	B	301	-	5,5,5	0.76	0	5,5,5	1.01	0
2	GOL	A	301	-	5,5,5	0.95	0	5,5,5	0.85	0
2	GOL	C	301	-	5,5,5	0.94	0	5,5,5	1.00	0
2	GOL	F	301	-	5,5,5	0.91	0	5,5,5	0.86	0
2	GOL	E	301	-	5,5,5	1.08	0	5,5,5	0.70	0
2	GOL	C	302	-	5,5,5	0.93	0	5,5,5	0.95	0
2	GOL	G	301	-	5,5,5	0.97	0	5,5,5	0.93	0
2	GOL	J	301	-	5,5,5	0.88	0	5,5,5	1.04	0
2	GOL	I	301	-	5,5,5	0.91	0	5,5,5	0.91	0
2	GOL	L	301	-	5,5,5	0.97	0	5,5,5	0.92	0
2	GOL	K	301	-	5,5,5	1.00	0	5,5,5	0.87	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	GOL	H	301	-	-	0/4/4/4	-
2	GOL	B	301	-	-	2/4/4/4	-
2	GOL	A	301	-	-	1/4/4/4	-
2	GOL	C	301	-	-	0/4/4/4	-
2	GOL	F	301	-	-	2/4/4/4	-
2	GOL	E	301	-	-	2/4/4/4	-
2	GOL	C	302	-	-	2/4/4/4	-
2	GOL	G	301	-	-	2/4/4/4	-
2	GOL	J	301	-	-	0/4/4/4	-
2	GOL	I	301	-	-	2/4/4/4	-
2	GOL	L	301	-	-	2/4/4/4	-
2	GOL	K	301	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	301	GOL	C1-C2-C3-O3
2	F	301	GOL	C1-C2-C3-O3
2	E	301	GOL	O1-C1-C2-C3
2	G	301	GOL	O1-C1-C2-O2
2	I	301	GOL	C1-C2-C3-O3
2	L	301	GOL	O1-C1-C2-C3
2	C	302	GOL	O1-C1-C2-C3
2	G	301	GOL	O1-C1-C2-C3
2	F	301	GOL	O2-C2-C3-O3
2	E	301	GOL	O1-C1-C2-O2
2	L	301	GOL	O1-C1-C2-O2
2	B	301	GOL	O2-C2-C3-O3
2	C	302	GOL	O1-C1-C2-O2
2	I	301	GOL	O2-C2-C3-O3
2	A	301	GOL	O1-C1-C2-C3

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	301	GOL	1	0
2	F	301	GOL	1	0
2	G	301	GOL	1	0
2	I	301	GOL	1	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 i

6.1 Protein, DNA and RNA chains i

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, 95th 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 ²)	Q<0.9
1	A	241/249 (96%)	-0.36	1 (0%) 92 95	19, 33, 67, 99	0
1	B	241/249 (96%)	-0.52	0 100 100	17, 33, 63, 90	0
1	C	241/249 (96%)	-0.36	2 (0%) 86 89	21, 36, 70, 88	0
1	D	241/249 (96%)	-0.52	0 100 100	18, 32, 61, 108	0
1	E	241/249 (96%)	-0.33	5 (2%) 63 70	18, 35, 67, 90	0
1	F	241/249 (96%)	-0.09	2 (0%) 86 89	28, 52, 79, 102	0
1	G	241/249 (96%)	-0.57	0 100 100	14, 28, 56, 87	0
1	H	241/249 (96%)	-0.52	2 (0%) 86 89	15, 30, 64, 87	0
1	I	241/249 (96%)	-0.55	0 100 100	13, 27, 54, 82	0
1	J	241/249 (96%)	-0.35	2 (0%) 86 89	15, 31, 64, 122	0
1	K	241/249 (96%)	-0.17	4 (1%) 70 76	23, 40, 78, 113	0
1	L	241/249 (96%)	0.02	2 (0%) 86 89	32, 58, 84, 102	0
All	All	2892/2988 (96%)	-0.36	20 (0%) 87 91	13, 36, 73, 122	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	206	ASP	6.3
1	E	137	ASN	3.6
1	E	134	LYS	3.2
1	L	140	LEU	3.2
1	K	137	ASN	3.1
1	E	77	LYS	2.9
1	J	204	THR	2.7
1	F	123	PHE	2.7
1	E	82	TYR	2.7
1	L	79	TYR	2.4
1	K	77	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	K	206	ASP	2.4
1	E	78	GLU	2.3
1	F	135	GLY	2.3
1	A	137	ASN	2.2
1	C	208	SER	2.1
1	H	266	LYS	2.1
1	C	206	ASP	2.1
1	H	82	TYR	2.0
1	K	266	LYS	2.0

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

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, 95th 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(Å ²)	Q<0.9
1	KCX	K	52	12/13	0.96	0.11	25,29,48,48	0
1	KCX	A	52	12/13	0.96	0.12	18,23,37,37	0
1	KCX	E	52	12/13	0.97	0.12	27,29,37,40	0
1	KCX	H	52	12/13	0.97	0.10	17,19,31,33	0
1	KCX	L	52	12/13	0.97	0.12	41,46,50,54	0
1	KCX	F	52	12/13	0.97	0.11	36,41,48,51	0
1	KCX	J	52	12/13	0.98	0.11	14,23,43,44	0
1	KCX	B	52	12/13	0.98	0.10	18,22,35,39	0
1	KCX	G	52	12/13	0.98	0.10	14,17,35,37	0
1	KCX	I	52	12/13	0.98	0.11	14,18,28,30	0
1	KCX	D	52	12/13	0.98	0.09	19,23,32,32	0
1	KCX	C	52	12/13	0.99	0.10	22,26,34,35	0

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. The B-factors column lists the minimum, median, 95th 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(\AA^2)	Q<0.9
2	GOL	E	301	6/6	0.78	0.20	69,71,73,74	0
2	GOL	C	301	6/6	0.82	0.21	72,75,78,80	0
2	GOL	B	301	6/6	0.85	0.17	54,59,64,71	0
2	GOL	A	301	6/6	0.86	0.14	56,60,62,64	0
2	GOL	K	301	6/6	0.86	0.24	62,65,66,66	0
2	GOL	L	301	6/6	0.87	0.15	63,65,65,66	0
2	GOL	H	301	6/6	0.89	0.13	63,67,69,70	0
2	GOL	G	301	6/6	0.90	0.16	62,65,66,66	0
2	GOL	J	301	6/6	0.91	0.16	57,58,62,62	0
2	GOL	C	302	6/6	0.92	0.14	62,69,70,71	0
2	GOL	F	301	6/6	0.94	0.14	58,59,61,64	0
2	GOL	I	301	6/6	0.95	0.12	51,55,61,63	0

6.5 Other polymers [i](#)

There are no such residues in this entry.