



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 17, 2024 – 11:38 AM EDT

PDB ID : 3OE3  
Title : Crystal structure of PliC-St, periplasmic lysozyme inhibitor of C-type lysozyme from *Salmonella typhimurium*  
Authors : Leysen, S.; Van Herreweghe, J.M.; Callewaert, L.; Heirbaut, M.; Buntinx, P.; Michiels, C.W.; Strelkov, S.V.  
Deposited on : 2010-08-12  
Resolution : 1.51 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 1.20.1  
EDS : 2.37.1  
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) : 2.37.1

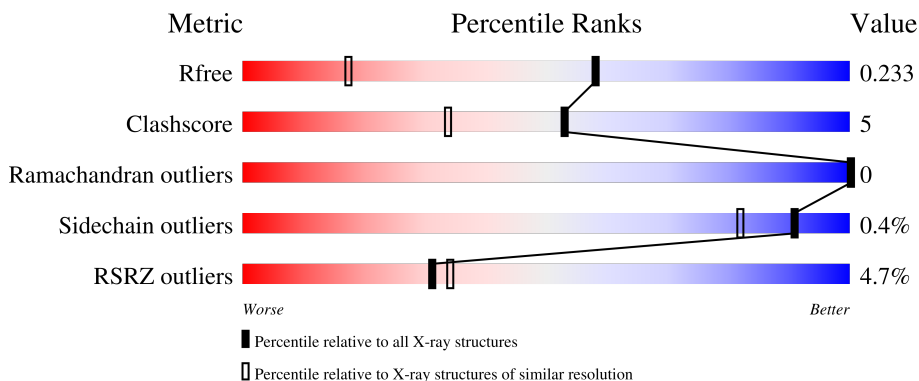
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.51 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.



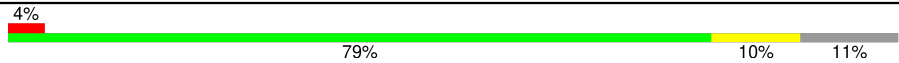
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	4009 (1.54-1.50)
Clashscore	141614	4249 (1.54-1.50)
Ramachandran outliers	138981	4148 (1.54-1.50)
Sidechain outliers	138945	4146 (1.54-1.50)
RSRZ outliers	127900	3943 (1.54-1.50)

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 of 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	98	 4% 73% 12% • 13%
1	B	98	 5% 80% 6% 14%
1	C	98	 2% 80% 9% • 10%
1	D	98	 2% 80% 10% 10%
1	E	98	 6% 67% 16% 16%

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Mol	Chain	Length	Quality of chain
1	F	98	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into four segments: a small red segment at the beginning labeled '4%', a large green segment labeled '79%', a yellow segment labeled '10%', and a grey segment at the end labeled '11%'.</p>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4434 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 Putative periplasmic protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	85	671	420	107	137	2	5	0	1	0
1	B	84	660	414	105	134	2	5	0	0	0
1	C	88	683	426	110	140	2	5	0	0	0
1	D	88	683	426	110	140	2	5	0	0	0
1	E	82	647	406	103	132	2	4	0	0	0
1	F	87	674	421	108	138	2	5	0	0	0

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	115	LEU	-	EXPRESSION TAG	UNP Q8ZPY8
A	116	GLU	-	EXPRESSION TAG	UNP Q8ZPY8
A	117	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
A	118	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
A	119	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
A	120	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
A	121	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
A	122	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
B	115	LEU	-	EXPRESSION TAG	UNP Q8ZPY8
B	116	GLU	-	EXPRESSION TAG	UNP Q8ZPY8
B	117	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
B	118	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
B	119	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
B	120	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
B	121	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
B	122	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
C	115	LEU	-	EXPRESSION TAG	UNP Q8ZPY8

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Chain	Residue	Modelled	Actual	Comment	Reference
C	116	GLU	-	EXPRESSION TAG	UNP Q8ZPY8
C	117	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
C	118	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
C	119	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
C	120	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
C	121	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
C	122	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
D	115	LEU	-	EXPRESSION TAG	UNP Q8ZPY8
D	116	GLU	-	EXPRESSION TAG	UNP Q8ZPY8
D	117	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
D	118	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
D	119	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
D	120	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
D	121	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
D	122	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
E	115	LEU	-	EXPRESSION TAG	UNP Q8ZPY8
E	116	GLU	-	EXPRESSION TAG	UNP Q8ZPY8
E	117	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
E	118	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
E	119	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
E	120	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
E	121	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
E	122	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
F	115	LEU	-	EXPRESSION TAG	UNP Q8ZPY8
F	116	GLU	-	EXPRESSION TAG	UNP Q8ZPY8
F	117	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
F	118	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
F	119	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
F	120	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
F	121	HIS	-	EXPRESSION TAG	UNP Q8ZPY8
F	122	HIS	-	EXPRESSION TAG	UNP Q8ZPY8

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Na 1 1	0	0
2	C	2	Total Na 2 2	0	0
2	D	1	Total Na 1 1	0	0
2	E	1	Total Na 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	F	1	Total	Na	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	64	Total	O	0	0
			64	64		
3	B	45	Total	O	0	0
			45	45		
3	C	101	Total	O	0	0
			101	101		
3	D	84	Total	O	0	0
			84	84		
3	E	50	Total	O	0	0
			50	50		
3	F	66	Total	O	0	0
			66	66		

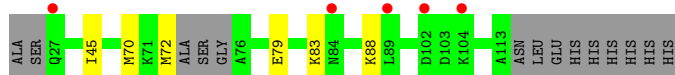
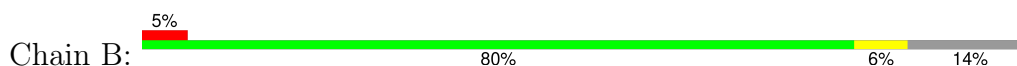
### 3 Residue-property plots [i](#)

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.

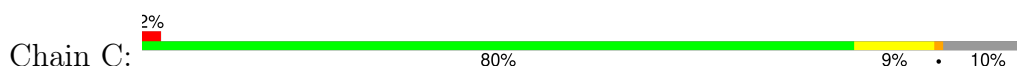
- Molecule 1: Putative periplasmic protein



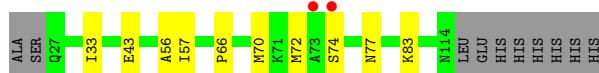
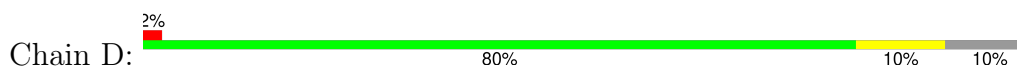
- Molecule 1: Putative periplasmic protein



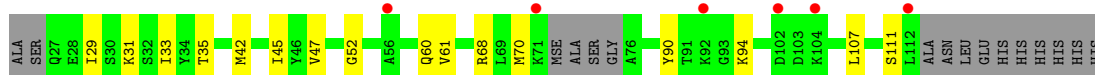
- Molecule 1: Putative periplasmic protein




- Molecule 1: Putative periplasmic protein



- Molecule 1: Putative periplasmic protein



- Molecule 1: Putative periplasmic protein

Chain F: 





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	192.38Å 34.69Å 74.91Å 90.00° 98.41° 90.00°	Depositor
Resolution (Å)	19.77 – 1.51 19.76 – 1.51	Depositor EDS
% Data completeness (in resolution range)	93.7 (19.77-1.51) 93.7 (19.76-1.51)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.98 (at 1.51Å)	Xtrriage
Refinement program	PHENIX 1.5_2	Depositor
R, $R_{free}$	0.208 , 0.242 0.204 , 0.233	Depositor DCC
$R_{free}$ test set	2087 reflections (2.87%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	18.4	Xtrriage
Anisotropy	0.280	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 63.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4434	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.27 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.3532e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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 [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.31	0/673	0.50	0/899
1	B	0.40	0/662	0.53	0/884
1	C	0.36	0/686	0.51	0/918
1	D	0.35	0/686	0.55	0/918
1	E	0.32	0/650	0.49	0/870
1	F	0.34	0/677	0.52	0/906
All	All	0.35	0/4034	0.52	0/5395

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	671	0	661	9	0
1	B	660	0	652	5	0
1	C	683	0	672	8	0
1	D	683	0	672	7	1
1	E	647	0	638	11	0
1	F	674	0	664	7	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	2	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	64	0	0	2	0
3	B	45	0	0	0	0
3	C	101	0	0	2	0
3	D	84	0	0	1	0
3	E	50	0	0	2	0
3	F	66	0	0	2	0
All	All	4434	0	3959	43	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 5.

All (43) 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:A:27:GLN:NE2	3:A:489:HOH:O	2.10	0.81
1:E:31:LYS:HE2	1:E:33:ILE:HD11	1.71	0.73
1:B:79:GLU:HG2	1:B:88:LYS:HG2	1.75	0.69
1:E:52:GLY:HA2	3:E:324:HOH:O	1.94	0.66
1:E:68:ARG:HD3	1:E:70:MSE:SE	2.46	0.64
1:A:92:LYS:O	1:A:95[B]:THR:HG22	2.02	0.59
1:E:60:GLN:O	1:E:61:VAL:HG22	2.05	0.57
1:C:74:SER:HB3	1:C:75:GLY:HA2	1.87	0.56
1:E:61:VAL:HG23	1:E:61:VAL:O	2.05	0.56
1:A:79:GLU:HG2	1:A:88:LYS:HG2	1.86	0.56
1:F:71:LYS:HD3	3:F:179:HOH:O	2.07	0.54
1:E:35:THR:OG1	3:E:481:HOH:O	1.82	0.54
1:E:29:ILE:HG22	1:E:47:VAL:HG22	1.90	0.54
1:C:70:MSE:SE	1:C:72:MSE:HE3	2.60	0.51
1:D:33:ILE:CD1	1:D:43:GLU:HG2	2.41	0.51
1:D:57:ILE:HD12	1:D:57:ILE:N	2.27	0.50
1:B:45:ILE:HD12	1:D:66:PRO:HG3	1.94	0.48
1:F:33:ILE:CD1	1:F:43:GLU:HG2	2.42	0.48
1:C:71:LYS:HE2	3:C:122:HOH:O	2.14	0.47
1:A:94:LYS:HD3	3:A:192:HOH:O	2.14	0.46
1:A:77:ASN:HD22	1:A:90:TYR:HA	1.82	0.45
1:D:56:ALA:C	1:D:57:ILE:HD12	2.36	0.45
1:D:70:MSE:HB3	1:D:72:MSE:HE3	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:42:MSE:HB2	1:E:107:LEU:HD13	1.99	0.45
1:C:70:MSE:HB3	1:C:72:MSE:HE3	2.00	0.43
1:D:74:SER:O	1:D:77:ASN:ND2	2.51	0.43
1:F:42:MSE:HB2	1:F:107:LEU:HD13	2.00	0.43
1:A:57:ILE:N	1:A:57:ILE:HD12	2.34	0.43
1:C:28:GLU:OE2	1:C:50:GLU:HG3	2.17	0.43
1:D:83:LYS:HG2	3:D:317:HOH:O	2.18	0.43
1:E:45:ILE:HD12	1:F:66:PRO:HG3	2.01	0.42
1:A:45:ILE:HD12	1:C:66:PRO:HG3	2.01	0.42
1:A:96:ALA:O	1:A:97:GLU:HG2	2.19	0.42
1:F:109:ASN:CG	1:F:109:ASN:O	2.58	0.41
1:B:83:LYS:HG3	1:E:90:TYR:CZ	2.54	0.41
1:C:42:MSE:HB2	1:C:107:LEU:HD13	2.03	0.41
1:A:51:ALA:HB3	3:C:126:HOH:O	2.20	0.41
1:B:72:MSE:HE3	1:B:72:MSE:HB2	2.00	0.41
1:C:74:SER:CB	1:C:75:GLY:HA2	2.48	0.41
1:E:94:LYS:O	1:E:111:SER:HA	2.21	0.41
1:F:35:THR:HG23	1:F:113:ALA:HB2	2.03	0.41
1:F:104:LYS:HE3	3:F:141:HOH:O	2.21	0.41
1:B:70:MSE:HG3	1:B:79:GLU:CD	2.42	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:74:SER:OG	1:D:74:SER:OG[2_656]	1.72	0.48

## 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	82/98 (84%)	81 (99%)	1 (1%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	80/98 (82%)	78 (98%)	2 (2%)	0	100	100
1	C	86/98 (88%)	81 (94%)	5 (6%)	0	100	100
1	D	86/98 (88%)	82 (95%)	4 (5%)	0	100	100
1	E	78/98 (80%)	77 (99%)	1 (1%)	0	100	100
1	F	85/98 (87%)	84 (99%)	1 (1%)	0	100	100
All	All	497/588 (84%)	483 (97%)	14 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	75/80 (94%)	74 (99%)	1 (1%)	69	43
1	B	74/80 (92%)	74 (100%)	0	100	100
1	C	76/80 (95%)	75 (99%)	1 (1%)	69	43
1	D	76/80 (95%)	76 (100%)	0	100	100
1	E	73/80 (91%)	73 (100%)	0	100	100
1	F	75/80 (94%)	75 (100%)	0	100	100
All	All	449/480 (94%)	447 (100%)	2 (0%)	91	82

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	C	71	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

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

Of 6 ligands modelled in this entry, 6 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 [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	80/98 (81%)	0.44	4 (5%) 28 31	14, 28, 48, 53	0
1	B	79/98 (80%)	0.31	5 (6%) 20 21	12, 26, 45, 50	0
1	C	83/98 (84%)	0.24	2 (2%) 59 63	13, 21, 38, 96	0
1	D	83/98 (84%)	0.30	2 (2%) 59 63	13, 21, 38, 91	0
1	E	78/98 (79%)	0.50	6 (7%) 13 14	17, 29, 51, 80	0
1	F	82/98 (83%)	0.29	4 (4%) 29 32	15, 24, 45, 67	0
All	All	485/588 (82%)	0.34	23 (4%) 31 34	12, 25, 46, 96	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	73	ALA	15.1
1	C	73	ALA	7.2
1	C	74	SER	6.2
1	E	102	ASP	3.7
1	B	27	GLN	3.2
1	A	38	ASP	3.1
1	F	83	LYS	2.9
1	E	71	LYS	2.9
1	D	74	SER	2.9
1	B	102	ASP	2.8
1	F	84	ASN	2.6
1	A	75	GLY	2.5
1	E	92	LYS	2.5
1	E	112	LEU	2.5
1	B	104	LYS	2.4
1	A	102	ASP	2.3
1	F	74	SER	2.3
1	E	56	ALA	2.2
1	B	89	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	E	104	LYS	2.1
1	F	85	TYR	2.1
1	A	61	VAL	2.0
1	B	84	ASN	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(Å <sup>2</sup> )	Q<0.9
2	NA	E	3	1/1	0.79	0.07	65,65,65,65	0
2	NA	D	2	1/1	0.90	0.07	30,30,30,30	0
2	NA	C	4	1/1	0.92	0.21	38,38,38,38	0
2	NA	C	5	1/1	0.92	0.24	49,49,49,49	0
2	NA	A	6	1/1	0.96	0.07	47,47,47,47	0
2	NA	F	1	1/1	0.96	0.10	28,28,28,28	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.