



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2017 – 02:41 PM EDT

PDB ID : 5J43  
Title : CdiA-CT from uropathogenic Escherichia coli in complex with CysK  
Authors : Morse, R.P.; Goulding, C.W.; Johnson, P.M.  
Deposited on : 2016-03-31  
Resolution : 2.70 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.

---

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20029077  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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-20029077

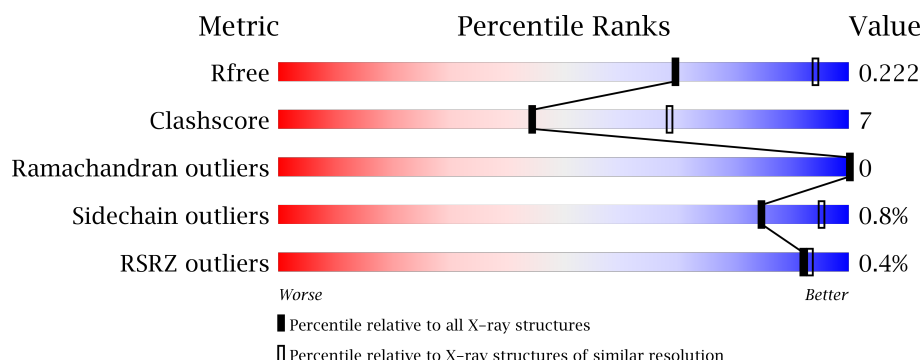
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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))
$R_{free}$	100719	2259 (2.70-2.70)
Clashscore	112137	2590 (2.70-2.70)
Ramachandran outliers	110173	2550 (2.70-2.70)
Sidechain outliers	110143	2550 (2.70-2.70)
RSRZ outliers	101464	2275 (2.70-2.70)

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	323	<div> <div>80%</div> <div>17%</div> <div>.</div> </div>
1	E	323	<div> <div>78%</div> <div>19%</div> <div>.</div> </div>
2	B	239	<div> <div>%</div> <div>36%</div> <div>6%</div> <div>58%</div> </div>
2	F	239	<div> <div>37%</div> <div>5%</div> <div>58%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6328 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 Cysteine synthase A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	E	314	Total	C	N	O	P	S	0	0	0
			2351	1483	405	456	1	6			
1	A	313	Total	C	N	O	P	S	0	0	0
			2346	1480	404	455	1	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	2	GLY	SER	conflict	UNP P0ABK6
A	2	GLY	SER	conflict	UNP P0ABK6

- Molecule 2 is a protein called tRNA nuclease CdiA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	101	Total	C	N	O	S	0	0	0
			765	469	141	152	3			
2	B	101	Total	C	N	O	S	0	0	0
			756	461	140	152	3			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	-11	MET	-	initiating methionine	UNP Q0T963
F	-10	ALA	-	expression tag	UNP Q0T963
F	-9	LYS	-	expression tag	UNP Q0T963
F	-8	SER	-	expression tag	UNP Q0T963
F	-7	HIS	-	expression tag	UNP Q0T963
F	-6	HIS	-	expression tag	UNP Q0T963
F	-5	HIS	-	expression tag	UNP Q0T963
F	-4	HIS	-	expression tag	UNP Q0T963
F	-3	HIS	-	expression tag	UNP Q0T963
F	-2	HIS	-	expression tag	UNP Q0T963

*Continued on next page...*

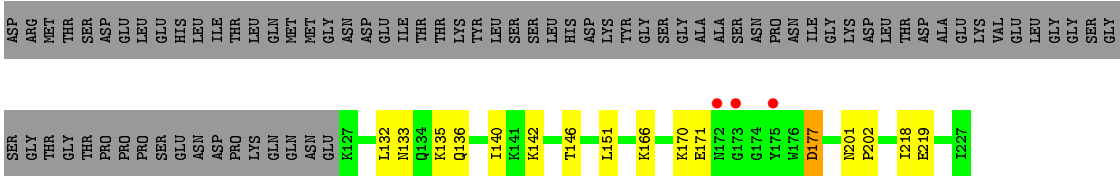
*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
F	-1	THR	-	expression tag	UNP Q0T963
F	0	SER	-	expression tag	UNP Q0T963
F	178	ALA	HIS	engineered mutation	UNP Q0T963
B	-11	MET	-	initiating methionine	UNP Q0T963
B	-10	ALA	-	expression tag	UNP Q0T963
B	-9	LYS	-	expression tag	UNP Q0T963
B	-8	SER	-	expression tag	UNP Q0T963
B	-7	HIS	-	expression tag	UNP Q0T963
B	-6	HIS	-	expression tag	UNP Q0T963
B	-5	HIS	-	expression tag	UNP Q0T963
B	-4	HIS	-	expression tag	UNP Q0T963
B	-3	HIS	-	expression tag	UNP Q0T963
B	-2	HIS	-	expression tag	UNP Q0T963
B	-1	THR	-	expression tag	UNP Q0T963
B	0	SER	-	expression tag	UNP Q0T963
B	178	ALA	HIS	engineered mutation	UNP Q0T963

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	E	40	Total O 40 40	0	0
3	F	9	Total O 9 9	0	0
3	A	48	Total O 48 48	0	0
3	B	13	Total O 13 13	0	0





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	64.01Å 64.01Å 365.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.92 – 2.70 44.92 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.0 (44.92-2.70) 99.3 (44.92-2.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.43 (at 2.69Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, $R_{free}$	0.200 , 0.224 0.198 , 0.222	Depositor DCC
$R_{free}$ test set	1980 reflections (4.99%)	DCC
Wilson B-factor (Å <sup>2</sup> )	47.5	Xtriage
Anisotropy	0.717	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 18.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.42$ , $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.499 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6328	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.99% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

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

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

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	0.46	0/2354	0.59	0/3185
1	E	0.47	0/2359	0.62	0/3192
2	B	0.45	0/762	0.56	0/1022
2	F	0.41	0/773	0.58	0/1038
All	All	0.46	0/6248	0.60	0/8437

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	2346	0	2412	36	0
1	E	2351	0	2414	45	0
2	B	756	0	749	11	0
2	F	765	0	757	7	0
3	A	48	0	0	2	0
3	B	13	0	0	0	0
3	E	40	0	0	0	0
3	F	9	0	0	0	0
All	All	6328	0	6332	94	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 7.

All (94) 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:E:48:ASN:HD22	1:E:152:ILE:HG22	1.56	0.70
2:F:167:PRO:HD2	2:F:169:PRO:HD3	1.74	0.69
1:E:65:LEU:HD22	1:E:141:LEU:HD21	1.76	0.67
1:A:65:LEU:HD22	1:A:141:LEU:HD21	1.78	0.66
1:E:152:ILE:HG12	1:E:153:HIS:N	2.11	0.65
1:E:256:ILE:HG12	1:E:272:ILE:HD12	1.80	0.64
1:E:260:ARG:HH22	1:E:313:ALA:HB3	1.63	0.62
2:B:177:ASP:OD1	2:B:177:ASP:N	2.24	0.61
1:A:228:GLN:HB3	3:A:429:HOH:O	2.03	0.59
1:A:114:THR:HG21	1:A:123:ALA:HA	1.83	0.59
2:F:151:LEU:HD21	2:F:218:ILE:HD11	1.84	0.58
2:B:142:LYS:O	2:B:146:THR:HG23	2.04	0.58
1:E:45:ILE:HG22	1:E:152:ILE:HD11	1.85	0.58
1:E:114:THR:HG21	1:E:123:ALA:HA	1.85	0.58
1:E:44:ARG:HB3	1:E:152:ILE:HG13	1.86	0.57
1:E:45:ILE:O	1:E:49:MET:HB2	2.04	0.57
1:E:164:ASP:HA	1:A:2:GLY:HA3	1.87	0.57
1:E:45:ILE:HB	1:E:49:MET:HE3	1.86	0.56
1:A:256:ILE:HG12	1:A:272:ILE:HD12	1.87	0.56
1:A:308:SER:OG	2:B:166:LYS:HE3	2.05	0.55
1:A:42:LLP:H2'1	1:A:72:ASN:ND2	2.21	0.54
1:E:38:SER:HB2	1:E:75:ILE:HG22	1.90	0.54
1:A:190:LYS:NZ	1:A:243:LEU:O	2.42	0.53
2:F:142:LYS:O	2:F:146:THR:HG23	2.09	0.53
1:A:45:ILE:HB	1:A:49:MET:HE3	1.91	0.53
1:E:162:TRP:NE1	1:E:193:LYS:HE2	2.24	0.53
1:E:263:MET:HG2	1:E:269:LEU:HA	1.90	0.53
1:E:155:LYS:O	1:E:159:PRO:HG2	2.09	0.52
1:E:2:GLY:HA3	1:A:164:ASP:HA	1.92	0.52
1:A:45:ILE:O	1:A:49:MET:HB2	2.09	0.52
1:E:48:ASN:HD22	1:E:152:ILE:CG2	2.21	0.52
1:E:152:ILE:O	1:E:156:THR:N	2.38	0.51
1:E:190:LYS:NZ	1:E:243:LEU:O	2.44	0.50
1:E:171:VAL:HB	1:E:295:ILE:HG12	1.94	0.50
1:A:307:LEU:HA	1:A:312:PHE:CE1	2.47	0.49
1:A:38:SER:HB2	1:A:75:ILE:HG22	1.95	0.49
1:A:263:MET:HG2	1:A:269:LEU:HA	1.95	0.48

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:257:SER:O	1:A:261:ARG:HG3	2.14	0.48
1:E:270:ALA:HB1	1:E:274:SER:HB2	1.95	0.48
1:E:49:MET:HG2	1:E:141:LEU:HD12	1.96	0.48
1:E:257:SER:O	1:E:261:ARG:HG3	2.14	0.48
1:E:157:THR:O	1:E:161:ILE:HG13	2.14	0.47
1:A:36:ASN:HB3	1:A:43:CYS:SG	2.56	0.46
1:E:30:ALA:HB1	1:E:299:LEU:HD22	1.98	0.46
1:A:222:PRO:HD2	2:B:219:GLU:OE2	2.15	0.46
1:E:209:PRO:O	1:E:213:GLN:HG3	2.15	0.46
1:E:153:HIS:CG	1:E:181:THR:HA	2.51	0.45
1:E:175:GLY:HA2	1:E:203:VAL:HB	1.98	0.45
1:E:42:LLP:H2'1	1:E:72:ASN:ND2	2.32	0.45
1:A:155:LYS:O	1:A:159:PRO:HG2	2.16	0.45
1:E:45:ILE:HG22	1:E:152:ILE:CD1	2.45	0.45
1:E:190:LYS:HA	1:E:195:LYS:O	2.17	0.45
1:A:33:GLU:OE2	1:A:44:ARG:NH1	2.50	0.45
2:B:136:GLN:O	2:B:140:ILE:HG13	2.16	0.45
1:A:171:VAL:HB	1:A:295:ILE:HG12	1.98	0.45
1:E:48:ASN:HA	1:E:51:TRP:HB2	1.99	0.45
1:A:146:ASN:HA	1:A:147:PRO:HD2	1.80	0.45
2:B:151:LEU:HD21	2:B:218:ILE:HD11	1.98	0.45
1:E:222:PRO:HD2	2:F:219:GLU:OE2	2.17	0.44
1:A:153:HIS:CG	1:A:181:THR:HA	2.52	0.44
1:E:272:ILE:HD13	1:E:311:LEU:HD13	2.00	0.44
1:A:42:LLP:OP3	1:A:42:LLP:H4'1	2.18	0.44
1:A:126:LYS:O	1:A:130:ILE:HG13	2.18	0.44
1:E:156:THR:OG1	1:E:157:THR:N	2.51	0.44
1:A:49:MET:HG2	1:A:141:LEU:HD12	2.00	0.43
1:A:79:TYR:CE2	1:A:80:VAL:HG23	2.54	0.43
2:F:147:ILE:O	2:F:151:LEU:HB2	2.18	0.43
1:A:156:THR:OG1	1:A:157:THR:N	2.52	0.42
1:E:44:ARG:HH22	1:E:160:GLU:CD	2.22	0.42
1:E:272:ILE:H	1:E:272:ILE:HG12	1.63	0.42
1:A:209:PRO:O	1:A:213:GLN:HG3	2.18	0.42
1:A:30:ALA:HB1	1:A:299:LEU:HD22	2.00	0.42
2:B:201:ASN:HA	2:B:202:PRO:HD3	1.88	0.42
1:E:63:VAL:HA	1:E:137:LYS:O	2.20	0.42
1:E:42:LLP:HE3	1:E:73:THR:OG1	2.19	0.42
2:B:170:LYS:HA	2:B:171:GLU:HA	1.76	0.42
1:E:146:ASN:HA	1:E:147:PRO:HD2	1.82	0.42
1:E:307:LEU:HA	1:E:312:PHE:CE1	2.55	0.42

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:133:ASN:OD1	2:F:136:GLN:HG3	2.20	0.41
1:E:288:GLU:HA	1:E:291:THR:OG1	2.19	0.41
1:A:256:ILE:HG12	1:A:272:ILE:CD1	2.49	0.41
1:A:15:THR:HG22	1:A:35:ARG:O	2.20	0.41
2:B:133:ASN:ND2	2:B:135:LYS:HB3	2.35	0.41
1:A:42:LLP:OP4	1:A:42:LLP:H4'1	2.21	0.41
1:A:162:TRP:NE1	1:A:193:LYS:HE2	2.36	0.41
1:A:35:ARG:HB3	3:A:430:HOH:O	2.21	0.41
1:E:120:MET:O	1:E:124:ILE:HG13	2.21	0.41
2:F:201:ASN:HA	2:F:202:PRO:HD3	1.92	0.41
1:E:256:ILE:HG12	1:E:272:ILE:CD1	2.48	0.40
1:E:60:LYS:HB2	1:E:61:PRO:HD2	2.03	0.40
1:A:228:GLN:HG3	1:A:228:GLN:H	1.72	0.40
1:A:60:LYS:HB2	1:A:61:PRO:HD2	2.03	0.40
2:B:151:LEU:HD23	2:B:151:LEU:HA	1.87	0.40
2:B:132:LEU:HD21	2:B:201:ASN:HB3	2.04	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	310/323 (96%)	300 (97%)	10 (3%)	0	100	100
1	E	311/323 (96%)	300 (96%)	11 (4%)	0	100	100
2	B	99/239 (41%)	98 (99%)	1 (1%)	0	100	100
2	F	99/239 (41%)	97 (98%)	2 (2%)	0	100	100
All	All	819/1124 (73%)	795 (97%)	24 (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	246/257 (96%)	245 (100%)	1 (0%)	93	98
1	E	246/257 (96%)	244 (99%)	2 (1%)	85	95
2	B	77/190 (40%)	76 (99%)	1 (1%)	73	91
2	F	78/190 (41%)	77 (99%)	1 (1%)	73	91
All	All	647/894 (72%)	642 (99%)	5 (1%)	85	95

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	92	MET
1	E	152	ILE
2	F	212	THR
1	A	92	MET
2	B	177	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	48	ASN
1	A	14	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

2 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	LLP	A	42	1	24,24,25	1.86	6 (25%)	28,32,34	2.17	4 (14%)
1	LLP	E	42	1	24,24,25	1.79	5 (20%)	28,32,34	2.29	4 (14%)

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	LLP	A	42	1	-	0/15/17/19	0/1/1/1
1	LLP	E	42	1	-	0/15/17/19	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	42	LLP	O3-C3	-5.58	1.24	1.37
1	E	42	LLP	O3-C3	-5.32	1.24	1.37
1	E	42	LLP	C4'-NZ	2.05	1.33	1.27
1	A	42	LLP	CE-NZ	2.13	1.51	1.46
1	E	42	LLP	C6-N1	2.26	1.39	1.34
1	A	42	LLP	C6-N1	2.33	1.39	1.34
1	A	42	LLP	CA-C	2.41	1.53	1.50
1	E	42	LLP	C2-N1	2.60	1.39	1.33
1	A	42	LLP	C2-N1	2.64	1.39	1.33
1	E	42	LLP	C4-C4'	3.10	1.52	1.46
1	A	42	LLP	C4-C4'	3.24	1.52	1.46

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	42	LLP	C5-C4-C4'	-5.68	112.86	121.36
1	A	42	LLP	C5-C4-C4'	-5.54	113.08	121.36
1	E	42	LLP	CE-NZ-C4'	-4.51	105.94	119.03
1	A	42	LLP	CE-NZ-C4'	-3.93	107.62	119.03
1	E	42	LLP	C3-C4-C4'	3.67	127.54	120.52
1	A	42	LLP	C3-C4-C4'	3.71	127.61	120.52
1	A	42	LLP	OP4-C5'-C5	6.72	122.83	109.32

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	42	LLP	OP4-C5'-C5	7.14	123.68	109.32

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	42	LLP	3	0
1	E	42	LLP	2	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 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	312/323 (96%)	-0.09	0 100 100	28, 38, 59, 74	0
1	E	313/323 (96%)	-0.09	0 100 100	28, 39, 60, 73	0
2	B	101/239 (42%)	0.08	3 (2%) 51 50	33, 45, 98, 128	0
2	F	101/239 (42%)	-0.03	0 100 100	30, 44, 96, 125	0
All	All	827/1124 (73%)	-0.06	3 (0%) 92 93	28, 40, 62, 128	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	172	ASN	4.1
2	B	173	GLY	3.4
2	B	175	TYR	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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
1	LLP	E	42	24/25	0.97	0.18	-	28,35,42,44	0
1	LLP	A	42	24/25	0.96	0.17	-	29,36,42,45	0

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.