



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 14, 2023 – 11:50 AM EDT

PDB ID : 4RZ4  
Title : Fructose-6-phosphate aldolase Q59E Y131F from E.coli  
Authors : Sandalova, T.; Stellmacher, L.; Leptihn, S.; Schneider, G.; Sprenger, G.A.; Samland, A.K.  
Deposited on : 2014-12-18  
Resolution : 1.75 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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35.1  
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) : 2.35.1

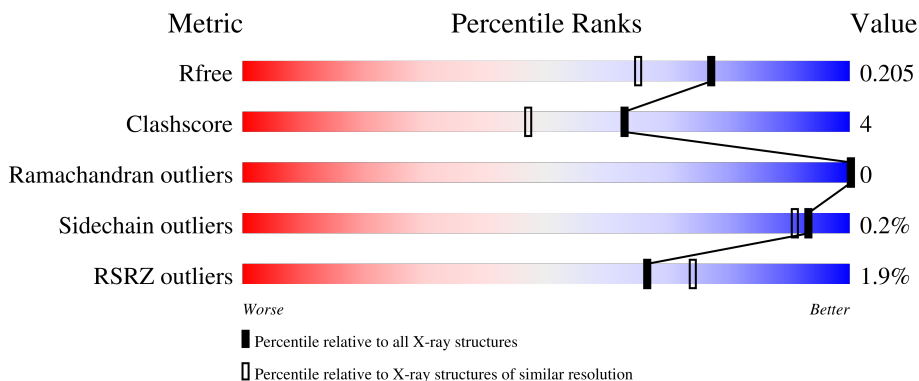
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.75 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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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	226	 90% 8% .
1	B	226	 3% 91% 6% .
1	C	226	 2% 91% 6% .
1	D	226	 92% 6% .
1	E	226	 2% 91% 7% .

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Mol	Chain	Length	Quality of chain
1	F	226	 % 90% 7%
1	G	226	 % 90% 7%
1	H	226	 4% 92% 6%
1	I	226	 % 89% 8%
1	J	226	 2% 89% 8%

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
4	CL	C	304	-	-	X	-
4	CL	D	304	-	-	X	-
4	CL	F	304	-	-	X	-
4	CL	G	304	-	-	X	-

## 2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 17605 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 Fructose-6-phosphate aldolase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	220	1617	1032	273	303	9	0	1	0
1	B	220	1614	1030	273	303	8	0	0	0
1	C	220	1617	1032	273	303	9	0	1	0
1	D	220	1617	1032	273	303	9	0	1	0
1	E	220	1617	1032	273	303	9	0	1	0
1	F	220	1617	1032	273	303	9	0	1	0
1	G	220	1617	1032	273	303	9	0	1	0
1	H	220	1617	1032	273	303	9	0	1	0
1	I	220	1617	1032	273	303	9	0	1	0
1	J	220	1617	1032	273	303	9	0	1	0

There are 90 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	MET	-	expression tag	UNP P78055
A	-4	HIS	-	expression tag	UNP P78055
A	-3	HIS	-	expression tag	UNP P78055
A	-2	HIS	-	expression tag	UNP P78055
A	-1	HIS	-	expression tag	UNP P78055
A	0	HIS	-	expression tag	UNP P78055
A	1	HIS	-	expression tag	UNP P78055
A	59	GLU	GLN	engineered mutation	UNP P78055
A	131	PHE	TYR	engineered mutation	UNP P78055

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-5	MET	-	expression tag	UNP P78055
B	-4	HIS	-	expression tag	UNP P78055
B	-3	HIS	-	expression tag	UNP P78055
B	-2	HIS	-	expression tag	UNP P78055
B	-1	HIS	-	expression tag	UNP P78055
B	0	HIS	-	expression tag	UNP P78055
B	1	HIS	-	expression tag	UNP P78055
B	59	GLU	GLN	engineered mutation	UNP P78055
B	131	PHE	TYR	engineered mutation	UNP P78055
C	-5	MET	-	expression tag	UNP P78055
C	-4	HIS	-	expression tag	UNP P78055
C	-3	HIS	-	expression tag	UNP P78055
C	-2	HIS	-	expression tag	UNP P78055
C	-1	HIS	-	expression tag	UNP P78055
C	0	HIS	-	expression tag	UNP P78055
C	1	HIS	-	expression tag	UNP P78055
C	59	GLU	GLN	engineered mutation	UNP P78055
C	131	PHE	TYR	engineered mutation	UNP P78055
D	-5	MET	-	expression tag	UNP P78055
D	-4	HIS	-	expression tag	UNP P78055
D	-3	HIS	-	expression tag	UNP P78055
D	-2	HIS	-	expression tag	UNP P78055
D	-1	HIS	-	expression tag	UNP P78055
D	0	HIS	-	expression tag	UNP P78055
D	1	HIS	-	expression tag	UNP P78055
D	59	GLU	GLN	engineered mutation	UNP P78055
D	131	PHE	TYR	engineered mutation	UNP P78055
E	-5	MET	-	expression tag	UNP P78055
E	-4	HIS	-	expression tag	UNP P78055
E	-3	HIS	-	expression tag	UNP P78055
E	-2	HIS	-	expression tag	UNP P78055
E	-1	HIS	-	expression tag	UNP P78055
E	0	HIS	-	expression tag	UNP P78055
E	1	HIS	-	expression tag	UNP P78055
E	59	GLU	GLN	engineered mutation	UNP P78055
E	131	PHE	TYR	engineered mutation	UNP P78055
F	-5	MET	-	expression tag	UNP P78055
F	-4	HIS	-	expression tag	UNP P78055
F	-3	HIS	-	expression tag	UNP P78055
F	-2	HIS	-	expression tag	UNP P78055
F	-1	HIS	-	expression tag	UNP P78055
F	0	HIS	-	expression tag	UNP P78055

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Chain	Residue	Modelled	Actual	Comment	Reference
F	1	HIS	-	expression tag	UNP P78055
F	59	GLU	GLN	engineered mutation	UNP P78055
F	131	PHE	TYR	engineered mutation	UNP P78055
G	-5	MET	-	expression tag	UNP P78055
G	-4	HIS	-	expression tag	UNP P78055
G	-3	HIS	-	expression tag	UNP P78055
G	-2	HIS	-	expression tag	UNP P78055
G	-1	HIS	-	expression tag	UNP P78055
G	0	HIS	-	expression tag	UNP P78055
G	1	HIS	-	expression tag	UNP P78055
G	59	GLU	GLN	engineered mutation	UNP P78055
G	131	PHE	TYR	engineered mutation	UNP P78055
H	-5	MET	-	expression tag	UNP P78055
H	-4	HIS	-	expression tag	UNP P78055
H	-3	HIS	-	expression tag	UNP P78055
H	-2	HIS	-	expression tag	UNP P78055
H	-1	HIS	-	expression tag	UNP P78055
H	0	HIS	-	expression tag	UNP P78055
H	1	HIS	-	expression tag	UNP P78055
H	59	GLU	GLN	engineered mutation	UNP P78055
H	131	PHE	TYR	engineered mutation	UNP P78055
I	-5	MET	-	expression tag	UNP P78055
I	-4	HIS	-	expression tag	UNP P78055
I	-3	HIS	-	expression tag	UNP P78055
I	-2	HIS	-	expression tag	UNP P78055
I	-1	HIS	-	expression tag	UNP P78055
I	0	HIS	-	expression tag	UNP P78055
I	1	HIS	-	expression tag	UNP P78055
I	59	GLU	GLN	engineered mutation	UNP P78055
I	131	PHE	TYR	engineered mutation	UNP P78055
J	-5	MET	-	expression tag	UNP P78055
J	-4	HIS	-	expression tag	UNP P78055
J	-3	HIS	-	expression tag	UNP P78055
J	-2	HIS	-	expression tag	UNP P78055
J	-1	HIS	-	expression tag	UNP P78055
J	0	HIS	-	expression tag	UNP P78055
J	1	HIS	-	expression tag	UNP P78055
J	59	GLU	GLN	engineered mutation	UNP P78055
J	131	PHE	TYR	engineered mutation	UNP P78055

- Molecule 2 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>4</sub>).



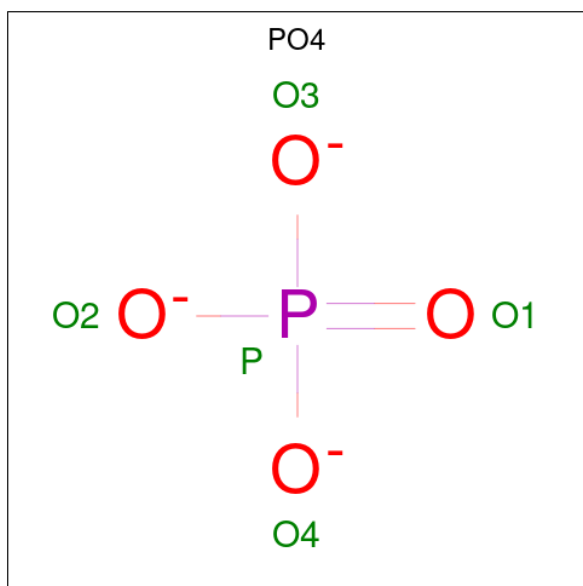
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	E	1	Total C O 4 2 2	0	0
2	E	1	Total C O 4 2 2	0	0
2	F	1	Total C O 4 2 2	0	0
2	F	1	Total C O 4 2 2	0	0
2	G	1	Total C O 4 2 2	0	0
2	G	1	Total C O 4 2 2	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	H	1	Total C O 4 2 2	0	0
2	H	1	Total C O 4 2 2	0	0
2	I	1	Total C O 4 2 2	0	0
2	I	1	Total C O 4 2 2	0	0
2	J	1	Total C O 4 2 2	0	0
2	J	1	Total C O 4 2 2	0	0

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O P 5 4 1	0	0
3	B	1	Total O P 5 4 1	0	0
3	C	1	Total O P 5 4 1	0	0
3	D	1	Total O P 5 4 1	0	0
3	E	1	Total O P 5 4 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	F	1	Total O P 5 4 1	0	0
3	G	1	Total O P 5 4 1	0	0
3	H	1	Total O P 5 4 1	0	0
3	I	1	Total O P 5 4 1	0	0
3	J	1	Total O P 5 4 1	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Cl 1 1	0	0
4	B	1	Total Cl 1 1	0	0
4	C	1	Total Cl 1 1	0	0
4	D	1	Total Cl 1 1	0	0
4	E	1	Total Cl 1 1	0	0
4	F	1	Total Cl 1 1	0	0
4	G	1	Total Cl 1 1	0	0
4	H	1	Total Cl 1 1	0	0
4	I	1	Total Cl 1 1	0	0
4	J	1	Total Cl 1 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	121	Total O 121 121	0	0
5	B	142	Total O 142 142	0	0

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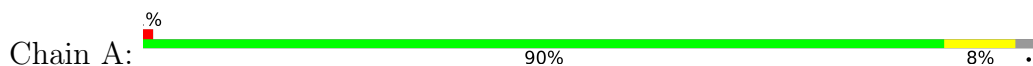
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	116	Total 116	O 116	0	0
5	D	124	Total 124	O 124	0	0
5	E	120	Total 120	O 120	0	0
5	F	134	Total 134	O 134	0	0
5	G	151	Total 151	O 151	0	0
5	H	121	Total 121	O 121	0	0
5	I	119	Total 119	O 119	0	0
5	J	150	Total 150	O 150	0	0

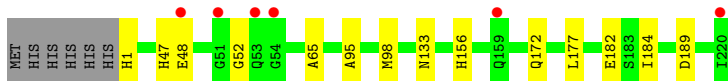
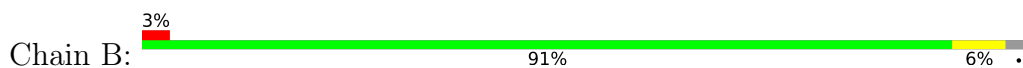
### 3 Residue-property plots [i](#)

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.

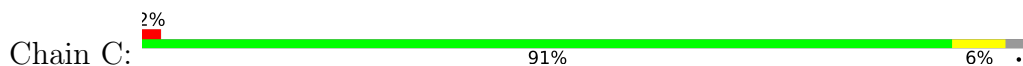
- Molecule 1: Fructose-6-phosphate aldolase 1



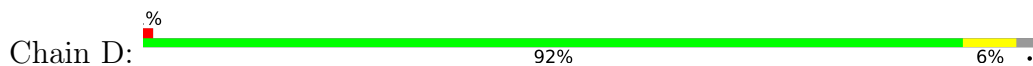
- Molecule 1: Fructose-6-phosphate aldolase 1



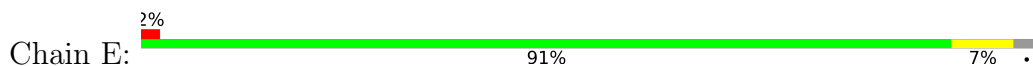
- Molecule 1: Fructose-6-phosphate aldolase 1



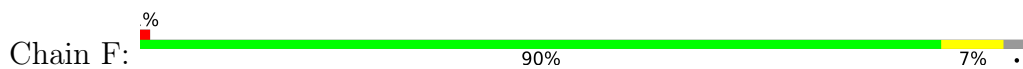
- Molecule 1: Fructose-6-phosphate aldolase 1



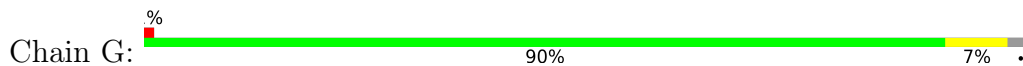
- Molecule 1: Fructose-6-phosphate aldolase 1



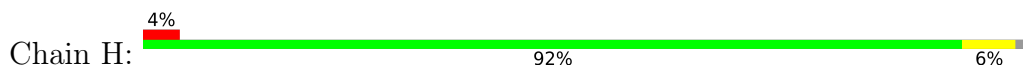
- Molecule 1: Fructose-6-phosphate aldolase 1



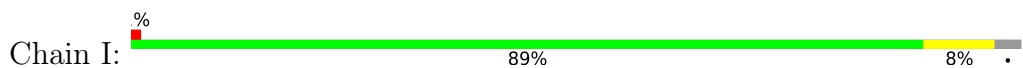
- Molecule 1: Fructose-6-phosphate aldolase 1



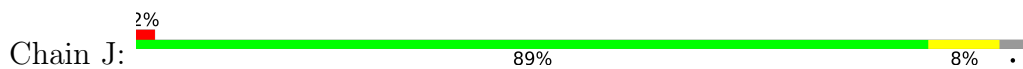
- Molecule 1: Fructose-6-phosphate aldolase 1



- Molecule 1: Fructose-6-phosphate aldolase 1



- Molecule 1: Fructose-6-phosphate aldolase 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.44Å 133.00Å 102.27Å 90.00° 92.37° 90.00°	Depositor
Resolution (Å)	88.44 – 1.75 40.51 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.5 (88.44-1.75) 99.6 (40.51-1.75)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.92 (at 1.75Å)	Xtrriage
Refinement program	REFMAC 5.8.0073	Depositor
R, $R_{free}$	0.179 , 0.196 0.189 , 0.205	Depositor DCC
$R_{free}$ test set	11921 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	16.4	Xtrriage
Anisotropy	0.032	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 41.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.020 for h,-k,-l	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	17605	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.33% 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: CL, PGE, PO4

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	0.49	0/1646	0.65	0/2238
1	B	0.47	0/1640	0.66	1/2230 (0.0%)
1	C	0.46	0/1646	0.64	0/2238
1	D	0.50	0/1646	0.64	0/2238
1	E	0.48	0/1646	0.67	1/2238 (0.0%)
1	F	0.50	0/1646	0.67	1/2238 (0.0%)
1	G	0.50	0/1646	0.68	2/2238 (0.1%)
1	H	0.50	0/1646	0.68	0/2238
1	I	0.51	0/1646	0.70	3/2238 (0.1%)
1	J	0.51	0/1646	0.71	2/2238 (0.1%)
All	All	0.49	0/16454	0.67	10/22372 (0.0%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	55	ARG	NE-CZ-NH2	-6.39	117.11	120.30
1	I	55	ARG	NE-CZ-NH1	6.10	123.35	120.30
1	G	195	ILE	CG1-CB-CG2	5.72	123.99	111.40
1	E	189	ASP	CB-CG-OD1	5.44	123.20	118.30
1	J	217	ARG	CA-CB-CG	5.36	125.20	113.40
1	B	189	ASP	CB-CG-OD1	5.34	123.10	118.30
1	G	189	ASP	CB-CG-OD1	5.20	122.98	118.30
1	F	189	ASP	CB-CG-OD1	5.15	122.94	118.30
1	J	189	ASP	CB-CG-OD1	5.14	122.93	118.30
1	I	55	ARG	CG-CD-NE	5.12	122.55	111.80

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	1617	0	1677	16	0
1	B	1614	0	1672	14	0
1	C	1617	0	1677	15	0
1	D	1617	0	1677	16	0
1	E	1617	0	1677	15	0
1	F	1617	0	1677	13	0
1	G	1617	0	1677	15	0
1	H	1617	0	1677	13	0
1	I	1617	0	1677	23	0
1	J	1617	0	1677	19	0
2	A	8	0	10	0	0
2	B	8	0	10	0	0
2	C	8	0	10	0	0
2	D	8	0	10	0	0
2	E	8	0	10	0	0
2	F	8	0	10	0	0
2	G	8	0	10	0	0
2	H	8	0	10	0	0
2	I	8	0	10	0	0
2	J	8	0	10	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	5	0	0	0	0
3	D	5	0	0	0	0
3	E	5	0	0	0	0
3	F	5	0	0	0	0
3	G	5	0	0	0	0
3	H	5	0	0	0	0
3	I	5	0	0	0	0
3	J	5	0	0	0	0
4	A	1	0	0	1	0
4	B	1	0	0	1	0
4	C	1	0	0	2	0
4	D	1	0	0	2	0
4	E	1	0	0	1	0
4	F	1	0	0	2	0
4	G	1	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	1	0	0	1	0
4	I	1	0	0	1	0
4	J	1	0	0	1	0
5	A	121	0	0	3	0
5	B	142	0	0	2	0
5	C	116	0	0	2	0
5	D	124	0	0	3	0
5	E	120	0	0	1	0
5	F	134	0	0	1	0
5	G	151	0	0	2	0
5	H	121	0	0	0	0
5	I	119	0	0	2	0
5	J	150	0	0	1	0
All	All	17605	0	16865	144	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 (144) 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:I:195:ILE:CD1	1:J:93:LEU:HD12	1.93	0.99
1:D:47:HIS:HD2	5:D:476:HOH:O	1.48	0.94
1:E:151:GLN:HG3	5:E:498:HOH:O	1.75	0.86
1:I:195:ILE:HD12	1:J:93:LEU:HD12	1.57	0.84
1:E:133:ASN:HD22	1:E:172:GLN:HE22	1.28	0.81
1:C:133:ASN:HD22	1:C:172:GLN:HE22	1.28	0.81
1:D:133:ASN:HD22	1:D:172:GLN:HE22	1.27	0.81
1:J:133:ASN:HD22	1:J:172:GLN:HE22	1.28	0.80
1:I:195:ILE:HD11	1:J:93:LEU:HD12	1.60	0.80
1:A:151:GLN:HG3	5:A:516:HOH:O	1.81	0.79
1:A:133:ASN:HD22	1:A:172:GLN:HE22	1.29	0.78
1:B:133:ASN:HD22	1:B:172:GLN:HE22	1.29	0.78
1:I:133:ASN:HD22	1:I:172:GLN:HE22	1.30	0.78
1:F:133:ASN:HD22	1:F:172:GLN:HE22	1.29	0.78
1:H:133:ASN:HD22	1:H:172:GLN:HE22	1.31	0.78
1:G:133:ASN:HD22	1:G:172:GLN:HE22	1.28	0.77
1:B:65:ALA:HB1	1:B:98:MET:CE	2.17	0.74
1:F:65:ALA:HB1	1:F:98:MET:CE	2.18	0.74
1:C:65:ALA:HB1	1:C:98:MET:CE	2.18	0.74
1:D:65:ALA:HB1	1:D:98:MET:CE	2.17	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:65:ALA:HB1	1:J:98:MET:CE	2.18	0.74
1:A:65:ALA:HB1	1:A:98:MET:CE	2.18	0.73
1:E:65:ALA:HB1	1:E:98:MET:CE	2.17	0.73
1:H:65:ALA:HB1	1:H:98:MET:CE	2.19	0.73
1:I:65:ALA:HB1	1:I:98:MET:CE	2.18	0.72
1:A:156:HIS:HD2	5:A:494:HOH:O	1.71	0.72
1:D:47:HIS:CD2	5:D:476:HOH:O	2.32	0.71
1:G:65:ALA:HB1	1:G:98:MET:CE	2.19	0.71
1:C:65:ALA:HB1	1:C:98:MET:HE2	1.72	0.71
1:I:65:ALA:HB1	1:I:98:MET:HE2	1.72	0.70
1:D:65:ALA:HB1	1:D:98:MET:HE2	1.72	0.69
1:B:156:HIS:HD2	5:B:516:HOH:O	1.75	0.69
1:F:156:HIS:HD2	5:F:512:HOH:O	1.75	0.69
1:A:65:ALA:HB1	1:A:98:MET:HE2	1.72	0.69
1:B:65:ALA:HB1	1:B:98:MET:HE2	1.74	0.69
1:D:156:HIS:HD2	5:D:521:HOH:O	1.75	0.68
1:H:65:ALA:HB1	1:H:98:MET:HE2	1.75	0.67
1:J:65:ALA:HB1	1:J:98:MET:HE2	1.76	0.67
5:G:541:HOH:O	1:H:156:HIS:HD2	1.76	0.67
1:I:195:ILE:HD12	1:J:93:LEU:CD1	2.25	0.66
1:E:65:ALA:HB1	1:E:98:MET:HE1	1.76	0.65
1:F:65:ALA:HB1	1:F:98:MET:HE2	1.78	0.64
1:G:65:ALA:HB1	1:G:98:MET:HE2	1.78	0.64
1:J:156:HIS:HD2	5:J:530:HOH:O	1.82	0.63
5:A:512:HOH:O	1:E:156:HIS:HD2	1.81	0.63
1:I:55:ARG:HH11	1:I:55:ARG:HG3	1.65	0.62
1:C:48:GLU:HG2	5:C:494:HOH:O	2.02	0.59
1:F:65:ALA:HB1	1:F:98:MET:HE1	1.84	0.59
1:E:65:ALA:HB1	1:E:98:MET:HE2	1.85	0.58
1:I:156:HIS:HD2	5:I:510:HOH:O	1.86	0.57
1:A:1:HIS:HE1	1:A:184:ILE:HG22	1.70	0.56
1:B:48:GLU:HG2	5:B:526:HOH:O	2.04	0.56
1:J:65:ALA:HB1	1:J:98:MET:HE1	1.86	0.56
1:C:156:HIS:HD2	5:C:512:HOH:O	1.89	0.55
1:B:65:ALA:HB1	1:B:98:MET:HE1	1.88	0.55
1:H:65:ALA:HB1	1:H:98:MET:HE1	1.89	0.55
1:G:156:HIS:HD2	5:G:539:HOH:O	1.89	0.55
1:G:65:ALA:HB1	1:G:98:MET:HE1	1.87	0.54
1:D:65:ALA:HB1	1:D:98:MET:HE1	1.90	0.54
1:A:156:HIS:HE1	1:B:177:LEU:O	1.91	0.53
1:I:195:ILE:O	1:I:195:ILE:HG12	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:1:HIS:HE1	1:J:184:ILE:HG22	1.74	0.53
1:D:1:HIS:HE1	1:D:184:ILE:HG22	1.74	0.53
1:I:1:HIS:HE1	1:I:184:ILE:HG22	1.74	0.53
1:C:1:HIS:HE1	1:C:184:ILE:HG22	1.74	0.52
1:B:1:HIS:HE1	1:B:184:ILE:HG22	1.73	0.52
1:E:1:HIS:HE1	1:E:184:ILE:HG22	1.73	0.52
1:I:65:ALA:HB1	1:I:98:MET:HE1	1.91	0.52
1:C:65:ALA:HB1	1:C:98:MET:HE1	1.90	0.52
1:H:1:HIS:HE1	1:H:184:ILE:HG22	1.75	0.51
1:F:1:HIS:HE1	1:F:184:ILE:HG22	1.75	0.51
1:C:156:HIS:HE1	1:D:177:LEU:O	1.94	0.50
1:F:156:HIS:HE1	1:J:177:LEU:O	1.95	0.50
1:G:1:HIS:HE1	1:G:184:ILE:HG22	1.76	0.50
1:B:156:HIS:HE1	1:C:177:LEU:O	1.95	0.50
1:H:177:LEU:O	1:I:156:HIS:HE1	1.94	0.50
1:G:195:ILE:HD13	1:G:195:ILE:O	2.12	0.49
1:I:195:ILE:CD1	1:J:93:LEU:CD1	2.77	0.48
1:D:156:HIS:HE1	1:E:177:LEU:O	1.96	0.48
1:I:55:ARG:NH2	5:I:473:HOH:O	2.41	0.48
1:E:65:ALA:CB	1:E:98:MET:HE1	2.43	0.48
1:A:65:ALA:HB1	1:A:98:MET:HE1	1.91	0.48
1:D:182:GLU:OE2	4:D:304:CL:CL	2.69	0.48
1:A:177:LEU:O	1:E:156:HIS:HE1	1.97	0.47
1:G:177:LEU:O	1:H:156:HIS:HE1	1.97	0.47
1:H:182:GLU:OE2	4:H:304:CL:CL	2.69	0.47
1:F:177:LEU:O	1:G:156:HIS:HE1	1.99	0.46
1:I:182:GLU:OE2	4:I:304:CL:CL	2.70	0.46
1:A:1:HIS:CE1	1:A:184:ILE:HG22	2.50	0.46
1:G:182:GLU:OE2	4:G:304:CL:CL	2.71	0.46
1:A:95:ALA:HA	1:A:98:MET:HE3	1.98	0.46
1:I:95:ALA:HA	1:I:98:MET:HE3	1.98	0.45
1:C:95:ALA:HA	1:C:98:MET:HE3	1.99	0.45
1:E:182:GLU:OE2	4:E:304:CL:CL	2.72	0.45
1:J:182:GLU:OE2	4:J:304:CL:CL	2.71	0.44
1:F:182:GLU:OE2	4:F:304:CL:CL	2.73	0.44
1:D:95:ALA:HA	1:D:98:MET:HE3	2.00	0.44
1:H:95:ALA:HA	1:H:98:MET:CE	2.48	0.44
1:E:1:HIS:CE1	1:E:184:ILE:HG22	2.52	0.44
1:I:95:ALA:HA	1:I:98:MET:CE	2.48	0.43
1:J:95:ALA:HA	1:J:98:MET:CE	2.48	0.43
1:B:47:HIS:CE1	1:B:52:GLY:HA2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:177:LEU:O	1:J:156:HIS:HE1	2.01	0.43
1:B:1:HIS:CE1	1:B:184:ILE:HG22	2.53	0.43
1:D:1:HIS:CE1	1:D:184:ILE:HG22	2.53	0.43
1:I:47:HIS:CE1	1:I:52:GLY:HA2	2.54	0.43
1:B:182:GLU:OE2	4:B:304:CL:CL	2.74	0.43
1:B:95:ALA:HA	1:B:98:MET:CE	2.48	0.43
1:C:1:HIS:CE1	1:C:184:ILE:HG22	2.53	0.43
1:A:95:ALA:HA	1:A:98:MET:CE	2.48	0.43
1:A:182:GLU:OE2	4:A:304:CL:CL	2.74	0.43
1:J:1:HIS:CE1	1:J:184:ILE:HG22	2.53	0.43
1:C:182:GLU:OE2	4:C:304:CL:CL	2.74	0.42
1:E:95:ALA:HA	1:E:98:MET:CE	2.49	0.42
1:H:95:ALA:HA	1:H:98:MET:HE3	2.00	0.42
1:J:79:ILE:HB	1:J:82:ILE:HB	2.01	0.42
1:C:95:ALA:HA	1:C:98:MET:CE	2.48	0.42
1:G:47:HIS:CE1	1:G:52:GLY:HA2	2.54	0.42
1:I:1:HIS:CE1	1:I:184:ILE:HG22	2.53	0.42
1:D:95:ALA:HA	1:D:98:MET:CE	2.49	0.42
1:J:47:HIS:CE1	1:J:52:GLY:HA2	2.53	0.42
1:C:47:HIS:CE1	1:C:52:GLY:HA2	2.55	0.42
1:E:47:HIS:CE1	1:E:52:GLY:HA2	2.55	0.42
1:F:150:HIS:ND1	4:F:304:CL:CL	2.88	0.42
1:F:47:HIS:CE1	1:F:52:GLY:HA2	2.55	0.42
1:G:1:HIS:CE1	1:G:184:ILE:HG22	2.54	0.42
1:H:47:HIS:CE1	1:H:52:GLY:HA2	2.55	0.42
1:G:95:ALA:HA	1:G:98:MET:CE	2.49	0.42
1:J:95:ALA:HA	1:J:98:MET:HE3	2.02	0.41
1:A:47:HIS:CE1	1:A:52:GLY:HA2	2.55	0.41
1:F:95:ALA:HA	1:F:98:MET:CE	2.50	0.41
1:C:150:HIS:ND1	4:C:304:CL:CL	2.90	0.41
1:B:95:ALA:HA	1:B:98:MET:HE3	2.02	0.41
1:I:79:ILE:HB	1:I:82:ILE:HB	2.02	0.41
1:D:47:HIS:CE1	1:D:52:GLY:HA2	2.56	0.41
1:F:24:GLY:HA2	1:F:50:MET:SD	2.60	0.41
1:I:24:GLY:HA2	1:I:50:MET:SD	2.61	0.41
1:A:24:GLY:HA2	1:A:50:MET:SD	2.61	0.40
1:H:55:ARG:HH11	1:H:55:ARG:HG2	1.84	0.40
1:D:150:HIS:ND1	4:D:304:CL:CL	2.88	0.40
1:G:79:ILE:HB	1:G:82:ILE:HB	2.03	0.40
1:G:150:HIS:ND1	4:G:304:CL:CL	2.87	0.40
1:A:79:ILE:HB	1:A:82:ILE:HB	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:107:LEU:C	1:E:107:LEU:HD23	2.42	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	219/226 (97%)	217 (99%)	2 (1%)	0	100	100
1	B	218/226 (96%)	216 (99%)	2 (1%)	0	100	100
1	C	219/226 (97%)	217 (99%)	2 (1%)	0	100	100
1	D	219/226 (97%)	217 (99%)	2 (1%)	0	100	100
1	E	219/226 (97%)	217 (99%)	2 (1%)	0	100	100
1	F	219/226 (97%)	217 (99%)	2 (1%)	0	100	100
1	G	219/226 (97%)	217 (99%)	2 (1%)	0	100	100
1	H	219/226 (97%)	217 (99%)	2 (1%)	0	100	100
1	I	219/226 (97%)	217 (99%)	2 (1%)	0	100	100
1	J	219/226 (97%)	217 (99%)	2 (1%)	0	100	100
All	All	2189/2260 (97%)	2169 (99%)	20 (1%)	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	168/173 (97%)	168 (100%)	0	100	100
1	B	167/173 (96%)	167 (100%)	0	100	100
1	C	168/173 (97%)	168 (100%)	0	100	100
1	D	168/173 (97%)	168 (100%)	0	100	100
1	E	168/173 (97%)	168 (100%)	0	100	100
1	F	168/173 (97%)	168 (100%)	0	100	100
1	G	168/173 (97%)	167 (99%)	1 (1%)	86	79
1	H	168/173 (97%)	168 (100%)	0	100	100
1	I	168/173 (97%)	167 (99%)	1 (1%)	86	79
1	J	168/173 (97%)	166 (99%)	2 (1%)	71	56
All	All	1679/1730 (97%)	1675 (100%)	4 (0%)	93	91

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	195	ILE
1	I	55	ARG
1	J	55	ARG
1	J	217	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (53) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1	HIS
1	A	47	HIS
1	A	133	ASN
1	A	144	GLN
1	A	156	HIS
1	A	212	GLN
1	B	1	HIS
1	B	133	ASN
1	B	144	GLN
1	B	156	HIS
1	B	212	GLN
1	C	1	HIS
1	C	47	HIS
1	C	133	ASN
1	C	156	HIS

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Mol	Chain	Res	Type
1	C	212	GLN
1	D	1	HIS
1	D	47	HIS
1	D	133	ASN
1	D	144	GLN
1	D	156	HIS
1	D	212	GLN
1	E	1	HIS
1	E	47	HIS
1	E	133	ASN
1	E	144	GLN
1	E	156	HIS
1	E	212	GLN
1	F	1	HIS
1	F	47	HIS
1	F	133	ASN
1	F	144	GLN
1	F	156	HIS
1	G	1	HIS
1	G	133	ASN
1	G	144	GLN
1	G	156	HIS
1	G	159	GLN
1	G	212	GLN
1	H	1	HIS
1	H	47	HIS
1	H	133	ASN
1	H	156	HIS
1	H	212	GLN
1	I	1	HIS
1	I	47	HIS
1	I	133	ASN
1	I	144	GLN
1	I	156	HIS
1	I	212	GLN
1	J	1	HIS
1	J	133	ASN
1	J	156	HIS

### 5.3.3 RNA ⓘ

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 40 ligands modelled in this entry, 10 are monoatomic - leaving 30 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	PGE	I	302	-	3,3,9	0.47	0	2,2,8	0.79	0
3	PO4	B	303	-	4,4,4	1.13	0	6,6,6	0.59	0
3	PO4	E	303	-	4,4,4	1.33	0	6,6,6	0.42	0
2	PGE	C	302	-	3,3,9	0.45	0	2,2,8	0.29	0
2	PGE	G	302	-	3,3,9	0.50	0	2,2,8	0.49	0
3	PO4	C	303	-	4,4,4	1.02	0	6,6,6	0.43	0
3	PO4	D	303	-	4,4,4	1.19	0	6,6,6	0.61	0
2	PGE	A	301	-	3,3,9	0.57	0	2,2,8	0.20	0
3	PO4	J	303	-	4,4,4	0.93	0	6,6,6	0.82	0
2	PGE	H	301	-	3,3,9	0.62	0	2,2,8	0.35	0
2	PGE	J	302	-	3,3,9	0.42	0	2,2,8	0.31	0
2	PGE	B	301	-	3,3,9	0.57	0	2,2,8	0.10	0
3	PO4	G	303	-	4,4,4	1.27	0	6,6,6	0.40	0
2	PGE	F	302	-	3,3,9	0.46	0	2,2,8	0.44	0
3	PO4	F	303	-	4,4,4	1.17	0	6,6,6	1.11	1 (16%)
2	PGE	E	301	-	3,3,9	0.63	0	2,2,8	0.03	0
2	PGE	D	301	-	3,3,9	0.53	0	2,2,8	0.06	0
2	PGE	G	301	-	3,3,9	0.65	0	2,2,8	0.49	0
2	PGE	I	301	-	3,3,9	0.69	0	2,2,8	0.20	0
2	PGE	C	301	-	3,3,9	0.56	0	2,2,8	0.05	0
2	PGE	H	302	-	3,3,9	0.51	0	2,2,8	0.46	0
2	PGE	F	301	-	3,3,9	0.64	0	2,2,8	0.18	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	PO4	H	303	-	4,4,4	1.16	0	6,6,6	0.33	0
2	PGE	E	302	-	3,3,9	0.50	0	2,2,8	0.48	0
2	PGE	J	301	-	3,3,9	0.80	0	2,2,8	0.41	0
2	PGE	D	302	-	3,3,9	0.41	0	2,2,8	0.22	0
2	PGE	B	302	-	3,3,9	0.42	0	2,2,8	0.42	0
3	PO4	I	303	-	4,4,4	1.22	0	6,6,6	0.19	0
3	PO4	A	303	-	4,4,4	1.25	0	6,6,6	0.73	0
2	PGE	A	302	-	3,3,9	0.41	0	2,2,8	0.13	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	PGE	I	302	-	-	0/1/1/7	-
2	PGE	C	302	-	-	0/1/1/7	-
2	PGE	G	302	-	-	0/1/1/7	-
2	PGE	A	301	-	-	0/1/1/7	-
2	PGE	H	301	-	-	0/1/1/7	-
2	PGE	J	302	-	-	0/1/1/7	-
2	PGE	B	301	-	-	0/1/1/7	-
2	PGE	F	302	-	-	0/1/1/7	-
2	PGE	E	301	-	-	0/1/1/7	-
2	PGE	D	301	-	-	0/1/1/7	-
2	PGE	G	301	-	-	0/1/1/7	-
2	PGE	I	301	-	-	0/1/1/7	-
2	PGE	C	301	-	-	0/1/1/7	-
2	PGE	H	302	-	-	0/1/1/7	-
2	PGE	F	301	-	-	0/1/1/7	-
2	PGE	E	302	-	-	0/1/1/7	-
2	PGE	J	301	-	-	0/1/1/7	-
2	PGE	D	302	-	-	0/1/1/7	-
2	PGE	B	302	-	-	0/1/1/7	-
2	PGE	A	302	-	-	0/1/1/7	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	303	PO4	O4-P-O3	2.19	115.00	107.97



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	220/226 (97%)	-0.05	3 (1%) 75 82	10, 15, 30, 47	0
1	B	220/226 (97%)	-0.04	6 (2%) 54 60	10, 15, 32, 50	0
1	C	220/226 (97%)	0.04	5 (2%) 60 67	11, 19, 34, 51	0
1	D	220/226 (97%)	0.09	3 (1%) 75 82	11, 18, 32, 51	0
1	E	220/226 (97%)	-0.13	4 (1%) 68 76	10, 15, 36, 50	0
1	F	220/226 (97%)	-0.16	2 (0%) 84 89	9, 13, 30, 46	0
1	G	220/226 (97%)	-0.15	2 (0%) 84 89	10, 14, 28, 47	0
1	H	220/226 (97%)	0.01	9 (4%) 37 44	10, 16, 33, 49	0
1	I	220/226 (97%)	0.04	3 (1%) 75 82	11, 17, 31, 44	0
1	J	220/226 (97%)	-0.09	4 (1%) 68 76	9, 14, 28, 46	0
All	All	2200/2260 (97%)	-0.04	41 (1%) 66 74	9, 16, 32, 51	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	51	GLY	4.4
1	I	220	ILE	4.4
1	J	54	GLY	4.2
1	D	220	ILE	4.0
1	H	54	GLY	3.9
1	E	220	ILE	3.4
1	H	220	ILE	3.4
1	A	159	GLN	3.3
1	E	53	GLN	3.2
1	A	220	ILE	3.2
1	B	53	GLN	3.2
1	D	216	GLY	3.1
1	B	220	ILE	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	54	GLY	3.0
1	B	54	GLY	3.0
1	H	11	VAL	3.0
1	G	220	ILE	2.9
1	H	216	GLY	2.9
1	C	53	GLN	2.8
1	E	48	GLU	2.8
1	H	48	GLU	2.8
1	B	48	GLU	2.8
1	A	53	GLN	2.8
1	J	220	ILE	2.6
1	I	195	ILE	2.6
1	C	220	ILE	2.6
1	B	51	GLY	2.5
1	B	159	GLN	2.4
1	G	53	GLN	2.4
1	C	74	LYS	2.3
1	C	48	GLU	2.3
1	H	15	ALA	2.3
1	E	215	PHE	2.2
1	I	51	GLY	2.2
1	H	215	PHE	2.2
1	D	218	THR	2.2
1	H	212	GLN	2.1
1	J	53	GLN	2.1
1	H	218	THR	2.1
1	F	159	GLN	2.1
1	F	220	ILE	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	PGE	I	301	4/10	0.92	0.09	22,25,25,25	0
2	PGE	D	301	4/10	0.93	0.10	26,27,27,28	0
2	PGE	F	301	4/10	0.93	0.10	20,20,21,21	0
2	PGE	C	301	4/10	0.93	0.10	26,27,27,28	0
2	PGE	B	301	4/10	0.94	0.10	22,23,23,23	0
2	PGE	I	302	4/10	0.94	0.10	22,22,23,23	0
2	PGE	H	301	4/10	0.95	0.09	21,22,22,23	0
2	PGE	C	302	4/10	0.95	0.09	22,22,22,22	0
2	PGE	G	301	4/10	0.95	0.07	18,19,19,20	0
3	PO4	F	303	5/5	0.95	0.09	34,37,39,39	0
2	PGE	D	302	4/10	0.96	0.09	21,21,21,22	0
2	PGE	H	302	4/10	0.96	0.07	19,19,19,20	0
2	PGE	E	301	4/10	0.96	0.08	19,20,20,21	0
2	PGE	B	302	4/10	0.96	0.08	17,17,18,19	0
2	PGE	J	301	4/10	0.96	0.09	18,19,20,20	0
2	PGE	A	302	4/10	0.96	0.09	18,18,19,19	0
2	PGE	E	302	4/10	0.97	0.09	17,17,17,17	0
2	PGE	G	302	4/10	0.97	0.07	18,19,20,20	0
2	PGE	A	301	4/10	0.97	0.08	21,23,23,25	0
3	PO4	E	303	5/5	0.97	0.09	31,37,37,38	0
2	PGE	F	302	4/10	0.97	0.11	17,17,17,18	0
3	PO4	G	303	5/5	0.97	0.09	33,38,39,39	0
3	PO4	H	303	5/5	0.97	0.09	37,38,43,43	0
3	PO4	I	303	5/5	0.97	0.08	34,38,39,39	0
3	PO4	J	303	5/5	0.97	0.08	30,33,33,35	0
4	CL	D	304	1/1	0.97	0.18	30,30,30,30	0
4	CL	I	304	1/1	0.97	0.14	28,28,28,28	0
3	PO4	C	303	5/5	0.98	0.10	37,38,40,42	0
3	PO4	D	303	5/5	0.98	0.09	36,40,42,43	0
2	PGE	J	302	4/10	0.98	0.08	16,16,17,18	0
4	CL	B	304	1/1	0.98	0.15	30,30,30,30	0
3	PO4	A	303	5/5	0.98	0.08	33,36,38,39	0
4	CL	E	304	1/1	0.98	0.14	28,28,28,28	0
4	CL	H	304	1/1	0.98	0.18	27,27,27,27	0
3	PO4	B	303	5/5	0.98	0.09	33,38,40,43	0
4	CL	A	304	1/1	0.99	0.14	29,29,29,29	0
4	CL	C	304	1/1	0.99	0.11	28,28,28,28	0
4	CL	J	304	1/1	0.99	0.15	26,26,26,26	0
4	CL	G	304	1/1	1.00	0.14	28,28,28,28	0
4	CL	F	304	1/1	1.00	0.14	24,24,24,24	0

## 6.5 Other polymers

There are no such residues in this entry.