



# Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 10:31 am BST

PDB ID : 2RK8  
Title : The Structure of rat cytosolic PEPCK in complex with phosphonoformate  
Authors : Sullivan, S.M.; Stiffin, R.M.; Carlson, G.M.; Holyoak, T.  
Deposited on : 2007-10-16  
Resolution : 2.00 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.

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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 : 2.11  
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

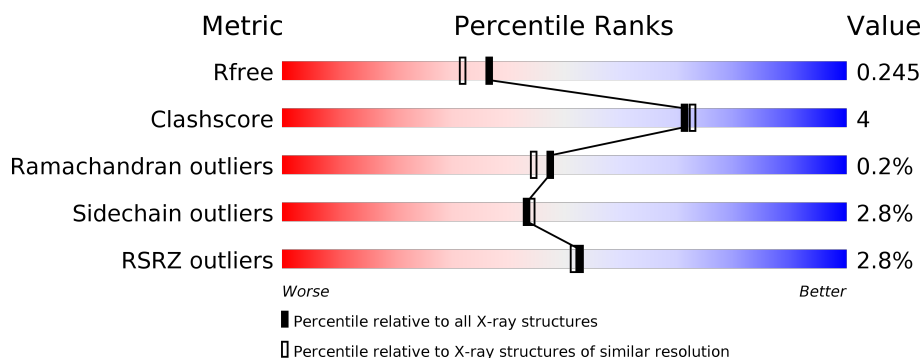
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 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	624	<div> <div>2%</div> <div> <div></div> <div>90%</div> <div>9%</div> </div> <div></div> </div>
1	B	624	<div> <div>3%</div> <div> <div></div> <div>85%</div> <div>12%</div> </div> <div></div> </div>

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 10831 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 Phosphoenolpyruvate carboxykinase, cytosolic [GTP].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	622	Total	C	N	O	S	0	23	0
			5030	3228	859	910	33			
1	B	611	Total	C	N	O	S	0	21	0
			4931	3165	837	896	33			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP P07379
A	0	SER	-	EXPRESSION TAG	UNP P07379
B	-1	GLY	-	EXPRESSION TAG	UNP P07379
B	0	SER	-	EXPRESSION TAG	UNP P07379

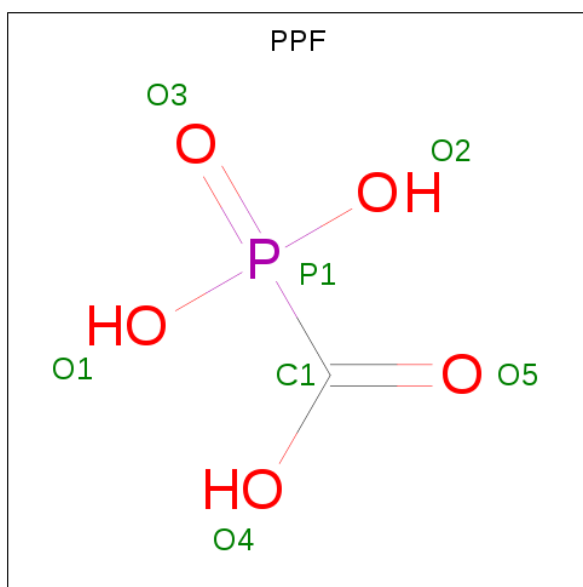
- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mn	0	0
			1	1		
2	A	1	Total	Mn	0	0
			1	1		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

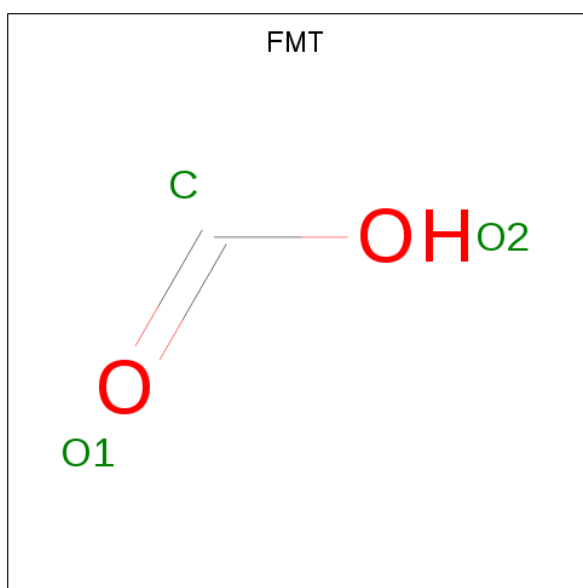
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Na	0	0
			1	1		
3	A	1	Total	Na	0	0
			1	1		

- Molecule 4 is PHOSPHONOFORMIC ACID (three-letter code: PPF) (formula: CH<sub>3</sub>O<sub>5</sub>P).



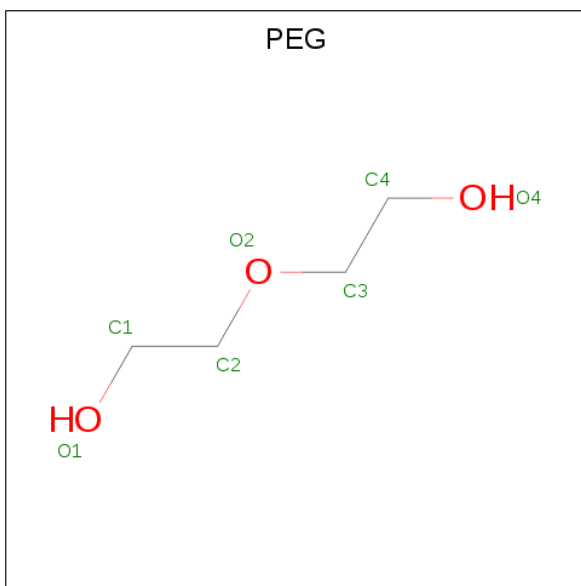
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	P	0	0
			7	1	5	1		
4	B	1	Total	C	O	P	0	0
			7	1	5	1		

- Molecule 5 is FORMIC ACID (three-letter code: FMT) (formula:  $\text{CH}_2\text{O}_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			3	1	2		
5	B	1	Total	C	O	0	1
			3	1	2		

- Molecule 6 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula:  $C_4H_{10}O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	4	3		

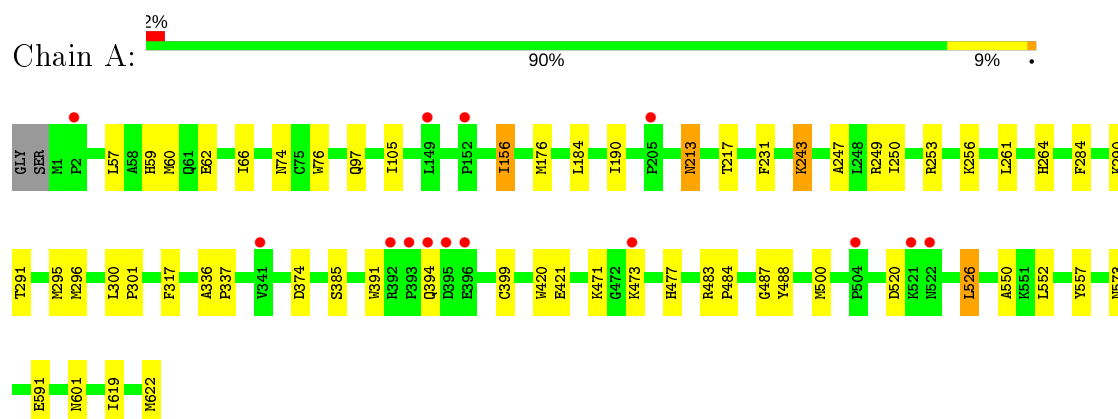
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	438	Total	O	0	0
			438	438		
7	B	401	Total	O	0	0
			401	401		

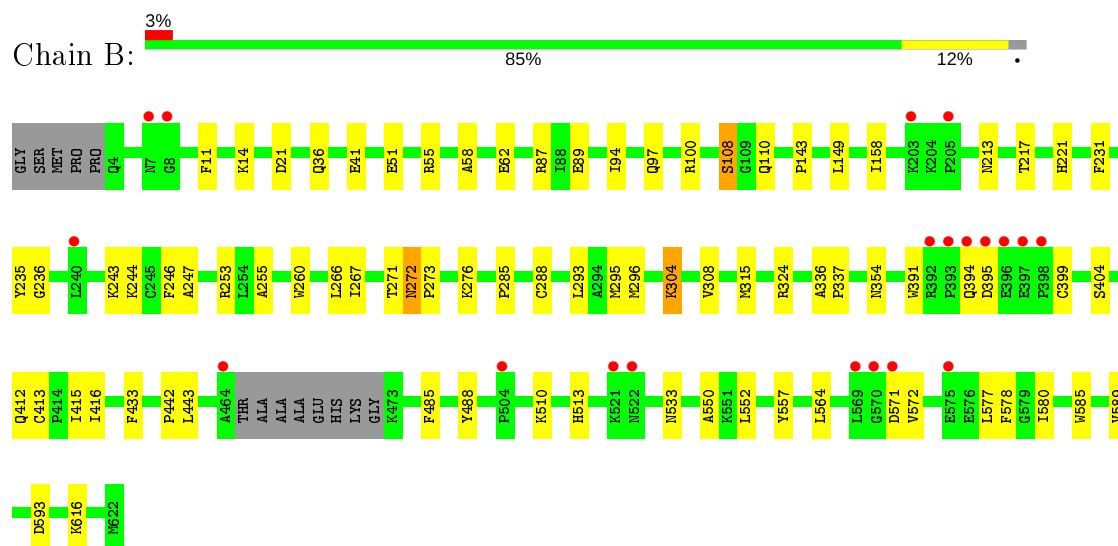
### 3 Residue-property plots [i](#)

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: Phosphoenolpyruvate carboxykinase, cytosolic [GTP]



- Molecule 1: Phosphoenolpyruvate carboxykinase, cytosolic [GTP]



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	62.25Å 119.47Å 86.87Å 90.00° 107.06° 90.00°	Depositor
Resolution (Å)	30.06 – 2.00 30.06 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.2 (30.06-2.00) 98.1 (30.06-2.00)	Depositor EDS
$R_{merge}$	0.71	Depositor
$R_{sym}$	0.52	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.43 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.195 , 0.249 0.195 , 0.245	Depositor DCC
$R_{free}$ test set	4000 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	24.2	Xtriage
Anisotropy	0.099	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 40.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10831	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 75.59 % 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 1.2277e-06. 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: PEG, NA, FMT, MN, PPF

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	0.41	0/5233	0.53	0/7080
1	B	0.43	4/5121 (0.1%)	0.55	1/6927 (0.0%)
All	All	0.42	4/10354 (0.0%)	0.54	1/14007 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	11	PHE	CG-CD1	6.66	1.48	1.38
1	B	11	PHE	CG-CD2	5.70	1.47	1.38
1	B	11	PHE	CE2-CZ	5.68	1.48	1.37
1	B	11	PHE	CE1-CZ	5.41	1.47	1.37

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	108	SER	N-CA-C	5.41	125.61	111.00

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	5030	0	5057	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	4931	0	4938	40	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	7	0	0	0	0
4	B	7	0	0	0	0
5	A	3	0	1	0	0
5	B	3	0	1	0	0
6	A	7	0	10	0	0
7	A	438	0	0	0	1
7	B	401	0	0	2	0
All	All	10831	0	10007	70	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 4.

All (70) 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:B:354:ASN:HD22	1:B:412:GLN:NE2	1.87	0.71
1:B:293:LEU:HD12	1:B:296:MET:HE3	1.70	0.71
1:B:354:ASN:HD22	1:B:412:GLN:HE22	1.45	0.62
1:A:156[A]:ILE:HD11	1:A:184:LEU:HD22	1.81	0.62
1:A:253[B]:ARG:HG2	1:A:488:TYR:HB3	1.83	0.60
1:B:51[B]:GLU:OE2	1:B:55[B]:ARG:NH1	2.34	0.60
1:B:354:ASN:HB2	1:B:412:GLN:HE22	1.69	0.58
1:B:443:LEU:HB2	1:B:577:LEU:HD22	1.86	0.57
1:A:420:TRP:CE2	1:A:421:GLU:HG3	2.40	0.56
1:B:253:ARG:HG2	1:B:488:TYR:HB3	1.88	0.55
1:B:550:ALA:HB1	1:B:557:TYR:HB3	1.89	0.54
1:A:619:ILE:O	1:A:622:MET:HG2	2.08	0.53
1:B:36:GLN:HG2	7:B:4300:HOH:O	2.08	0.53
1:A:336:ALA:HB3	1:A:337:PRO:HD3	1.92	0.52
1:B:413:CYS:HB3	1:B:416:ILE:HB	1.93	0.50
1:B:255:ALA:HB1	1:B:260:TRP:O	2.11	0.50
1:B:272:ASN:C	1:B:272:ASN:HD22	2.14	0.50
1:A:217:THR:HA	1:A:231:PHE:O	2.12	0.49
1:A:264:HIS:CD2	1:A:290[A]:LYS:NZ	2.81	0.49
1:B:296:MET:HE2	1:B:533:ASN:HB2	1.93	0.49
1:A:156[A]:ILE:HG12	1:A:184:LEU:HD13	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:PHE:HB2	1:A:290[A]:LYS:HG3	1.96	0.48
1:A:105:ILE:O	1:A:601[A]:ASN:ND2	2.46	0.48
1:B:336:ALA:HB3	1:B:337:PRO:HD3	1.97	0.47
1:A:317:PHE:HE2	1:A:500[B]:MET:CE	2.28	0.47
1:A:57:LEU:HD13	1:A:76:TRP:NE1	2.29	0.46
1:B:295:MET:HA	1:B:415:ILE:HD11	1.98	0.46
1:B:585:TRP:O	1:B:589:VAL:HG22	2.16	0.46
1:B:285:PRO:O	1:B:288[B]:CYS:HB2	2.15	0.46
1:A:243:LYS:O	1:A:247:ALA:HB3	2.16	0.45
1:B:14[B]:LYS:HD3	1:B:41:GLU:O	2.17	0.45
1:B:94:ILE:HG12	1:B:221:HIS:HB2	1.98	0.45
1:A:249:ARG:O	1:A:487:GLY:HA3	2.18	0.44
1:A:550:ALA:HB1	1:A:557:TYR:HB3	1.99	0.44
1:B:315:MET:HA	1:B:324:ARG:O	2.17	0.44
1:B:244:LYS:HD2	1:B:485:PHE:CE1	2.53	0.44
1:A:59:HIS:O	1:A:62[B]:GLU:HG2	2.18	0.44
1:A:74:ASN:ND2	1:A:76:TRP:HE1	2.15	0.44
1:B:272:ASN:HD21	1:B:276:LYS:H	1.66	0.44
1:A:261:LEU:HD11	1:A:500[B]:MET:SD	2.59	0.43
1:B:217:THR:HA	1:B:231:PHE:O	2.18	0.43
1:B:442:PRO:HB3	1:B:578:PHE:HA	2.00	0.43
1:A:213:ASN:HD22	1:A:213:ASN:C	2.22	0.43
1:B:593:ASP:OD2	1:B:616[A]:LYS:NZ	2.51	0.43
1:B:89:GLU:HG3	7:B:4004:HOH:O	2.18	0.43
1:B:58:ALA:O	1:B:62[B]:GLU:HG3	2.19	0.43
1:A:520:ASP:HB3	1:A:526:LEU:HD11	2.00	0.43
1:A:66:ILE:HD12	1:A:76:TRP:CE3	2.54	0.42
1:B:143:PRO:HA	1:B:158:ILE:HD13	2.00	0.42
1:B:266:LEU:HD21	1:B:293:LEU:HD23	2.00	0.42
1:B:337:PRO:HD3	1:B:404:SER:OG	2.20	0.42
1:A:291:THR:O	1:A:295:MET:HG2	2.19	0.42
1:B:433:PHE:O	1:B:513:HIS:HA	2.19	0.42
1:A:477:HIS:HE2	1:A:591:GLU:CD	2.23	0.42
1:A:60:MET:HB3	1:A:66:ILE:HG12	2.01	0.42
1:B:87:ARG:N	1:B:235[B]:TYR:CE2	2.88	0.42
1:B:296:MET:CE	1:B:533:ASN:HB2	2.50	0.41
1:B:243:LYS:O	1:B:247:ALA:HB3	2.20	0.41
1:B:272:ASN:ND2	1:B:276:LYS:H	2.18	0.41
1:A:300:LEU:HA	1:A:301:PRO:HD3	1.91	0.41
1:A:385:SER:HB3	1:A:391:TRP:HB2	2.03	0.41
1:B:271:THR:HB	1:B:304:LYS:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:391:TRP:CD2	1:A:399:CYS:HB3	2.56	0.40
1:B:272:ASN:HB2	1:B:273:PRO:CD	2.51	0.40
1:A:483:ARG:HB3	1:A:484:PRO:HD3	2.03	0.40
1:B:267:ILE:HB	1:B:308:VAL:HB	2.04	0.40
1:B:391:TRP:CD2	1:B:399:CYS:HB3	2.57	0.40
1:A:190:ILE:HD13	1:A:250:ILE:HG21	2.03	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:B:108:SER:O	7:A:6194:HOH:O[1_454]	2.07	0.13

## 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	644/624 (103%)	625 (97%)	18 (3%)	1 (0%)	47	44
1	B	628/624 (101%)	610 (97%)	17 (3%)	1 (0%)	47	44
All	All	1272/1248 (102%)	1235 (97%)	35 (3%)	2 (0%)	47	44

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	236	GLY
1	A	243	LYS

### 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	543/520 (104%)	529 (97%)	14 (3%)	46	48
1	B	533/520 (102%)	514 (96%)	19 (4%)	35	34
All	All	1076/1040 (104%)	1043 (97%)	33 (3%)	43	40

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	97	GLN
1	A	156[A]	ILE
1	A	156[B]	ILE
1	A	176	MET
1	A	213	ASN
1	A	256	LYS
1	A	374	ASP
1	A	394	GLN
1	A	471[A]	LYS
1	A	471[B]	LYS
1	A	473	LYS
1	A	526	LEU
1	A	552	LEU
1	A	573	ASN
1	B	21	ASP
1	B	97	GLN
1	B	100[A]	ARG
1	B	100[B]	ARG
1	B	110	GLN
1	B	149	LEU
1	B	213	ASN
1	B	246	PHE
1	B	272	ASN
1	B	304	LYS
1	B	394	GLN
1	B	395	ASP
1	B	510	LYS
1	B	552	LEU
1	B	564	LEU
1	B	571	ASP
1	B	572	VAL

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Mol	Chain	Res	Type
1	B	580[A]	ILE
1	B	580[B]	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	ASN
1	A	97	GLN
1	A	213	ASN
1	A	238	ASN
1	A	515	ASN
1	A	573	ASN
1	B	34	ASN
1	B	74	ASN
1	B	99	GLN
1	B	208	ASN
1	B	213	ASN
1	B	272	ASN
1	B	297	ASN
1	B	388	ASN
1	B	394	GLN
1	B	412	GLN
1	B	502	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](#)

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

### 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

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

Of 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 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	$\# Z  > 2$
5	FMT	B	3406[A]	-	0,2,2	0.00	-	0,1,1	0.00	-
4	PPF	B	3969	2	3,6,6	1.32	0	5,9,9	0.88	0
5	FMT	A	3406	-	0,2,2	0.00	-	0,1,1	0.00	-
6	PEG	A	5759	-	6,6,6	0.43	0	5,5,5	0.31	0
4	PPF	A	3969	2	3,6,6	1.52	0	5,9,9	1.14	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
4	PPF	B	3969	2	-	0/0/6/6	-
6	PEG	A	5759	-	-	3/4/4/4	-
4	PPF	A	3969	2	-	0/0/6/6	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	5759	PEG	O2-C3-C4-O4
6	A	5759	PEG	O1-C1-C2-O2
6	A	5759	PEG	C4-C3-O2-C2

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	622/624 (99%)	0.02	14 (2%) 60 59	7, 17, 28, 38	0
1	B	611/624 (97%)	0.02	20 (3%) 46 45	8, 18, 30, 43	0
All	All	1233/1248 (98%)	0.02	34 (2%) 53 51	7, 17, 29, 43	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	394	GLN	6.0
1	B	394	GLN	5.7
1	A	393	PRO	5.1
1	B	396	GLU	4.4
1	A	504	PRO	4.3
1	B	504	PRO	4.3
1	B	570	GLY	4.2
1	B	392	ARG	4.2
1	A	392	ARG	4.1
1	A	152	PRO	4.1
1	B	521	LYS	3.8
1	B	7	ASN	3.8
1	B	8	GLY	3.7
1	B	522	ASN	3.4
1	B	571	ASP	3.3
1	A	396	GLU	3.1
1	B	397	GLU	3.1
1	B	205	PRO	3.1
1	A	395	ASP	3.0
1	B	393	PRO	2.8
1	A	473	LYS	2.5
1	A	521	LYS	2.4
1	A	149	LEU	2.4
1	A	341	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	395	ASP	2.3
1	B	203	LYS	2.3
1	B	398	PRO	2.3
1	A	205	PRO	2.3
1	B	569	LEU	2.3
1	B	464	ALA	2.3
1	A	2	PRO	2.3
1	B	575	GLU	2.3
1	B	240	LEU	2.2
1	A	522	ASN	2.1

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

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

## 6.3 Carbohydrates [i](#)

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, 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
3	NA	A	624	1/1	0.86	0.08	43,43,43,43	0
6	PEG	A	5759	7/7	0.88	0.13	54,54,54,54	0
5	FMT	B	3406[A]	3/3	0.90	0.19	16,16,17,17	3
5	FMT	A	3406	3/3	0.93	0.12	37,37,38,38	0
3	NA	B	624	1/1	0.96	0.07	35,35,35,35	0
4	PPF	A	3969	7/7	0.97	0.08	12,16,17,17	0
4	PPF	B	3969	7/7	0.98	0.08	23,24,25,26	0
2	MN	B	700	1/1	0.99	0.06	18,18,18,18	0
2	MN	A	700	1/1	1.00	0.03	17,17,17,17	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.