



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:39 AM GMT

PDB ID : 3MOH
Title : The structure of rat cytosolic PEPCK mutant A467G in complex with phosphoglycolate and GDP
Authors : Johnson, T.A.; Holyoak, T.
Deposited on : 2010-04-22
Resolution : 2.10 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

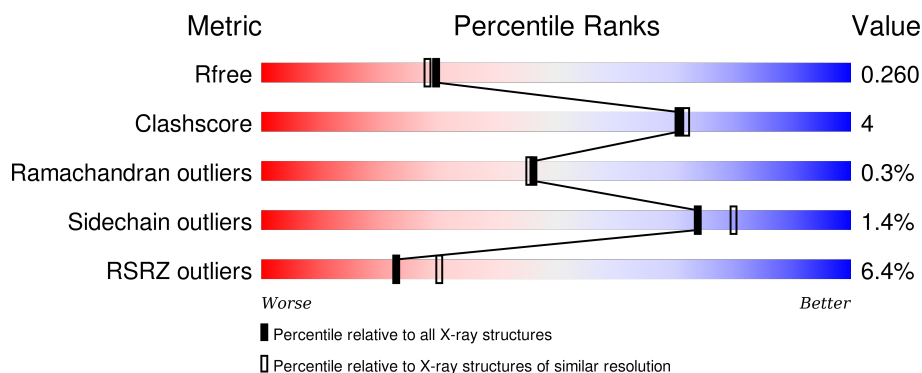
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	91344	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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>6%</div> <div>87%</div> <div>9%</div> <div>••</div> </div>
1	B	624	<div> <div>6%</div> <div>88%</div> <div>10%</div> <div>••</div> </div>

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
7	1PE	A	1300	-	-	-	X

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 10430 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.

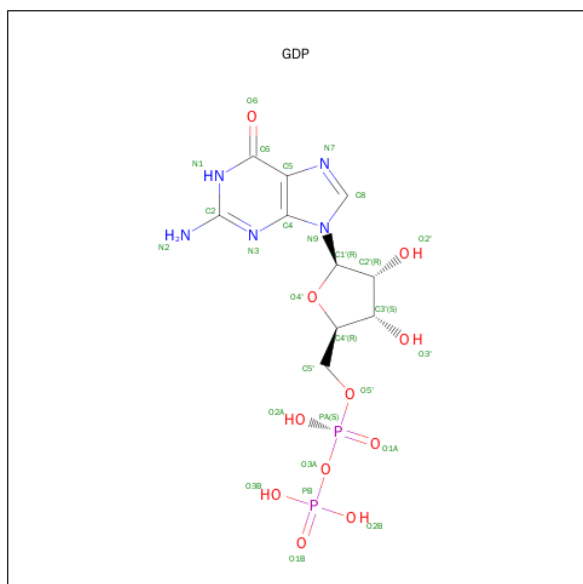
- Molecule 1 is a protein called Phosphoenolpyruvate carboxykinase, cytosolic [GTP].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	604	Total	C	N	O	S	0	10	0
			4811	3079	820	881	31			
1	B	613	Total	C	N	O	S	0	6	0
			4852	3104	828	887	33			

There are 6 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
A	467	GLY	ALA	ENGINEERED	UNP P07379
B	-1	GLY	-	EXPRESSION TAG	UNP P07379
B	0	SER	-	EXPRESSION TAG	UNP P07379
B	467	GLY	ALA	ENGINEERED	UNP P07379

- Molecule 2 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).

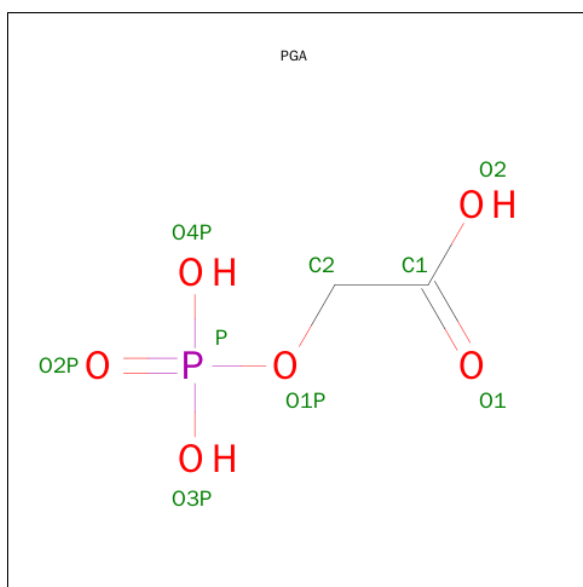


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	B	1	Total	C	N	O	P	0	0
			28	10	5	11	2		

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

- Molecule 4 is SUGAR (2-PHOSPHOGLYCOLIC ACID) (three-letter code: PGA) (formula: C₂H₅O₆P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	P	0	0
			9	2	6	1		
4	B	1	Total	C	O	P	0	0
			9	2	6	1		

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

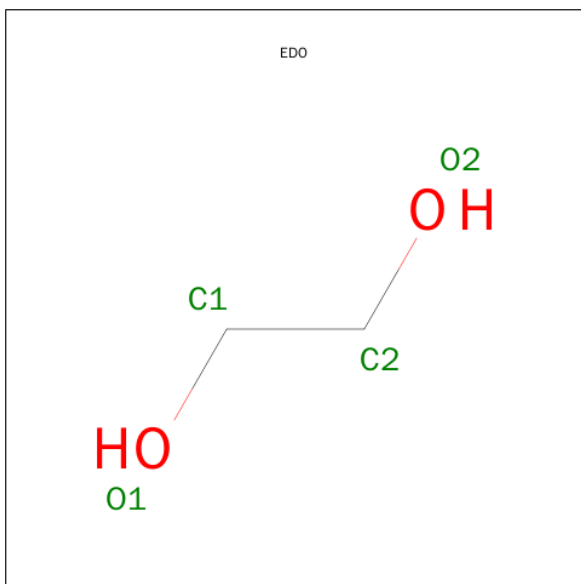
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Na	0	0
			1	1		

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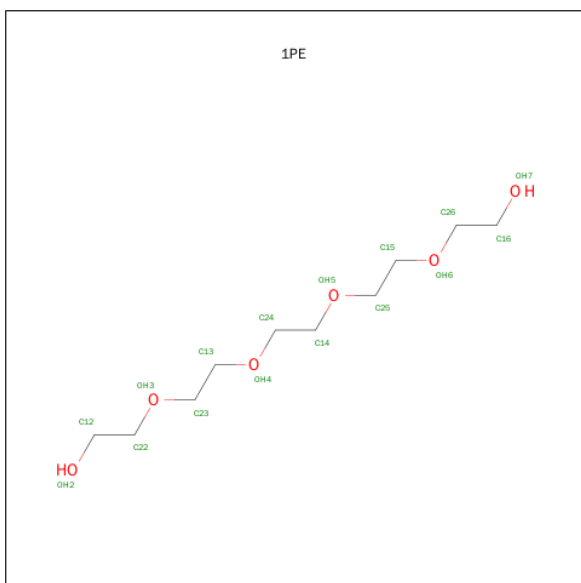
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Na	0	0
			1	1		

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			5	3	2		
7	B	1	Total	C	O	0	0
			10	6	4		

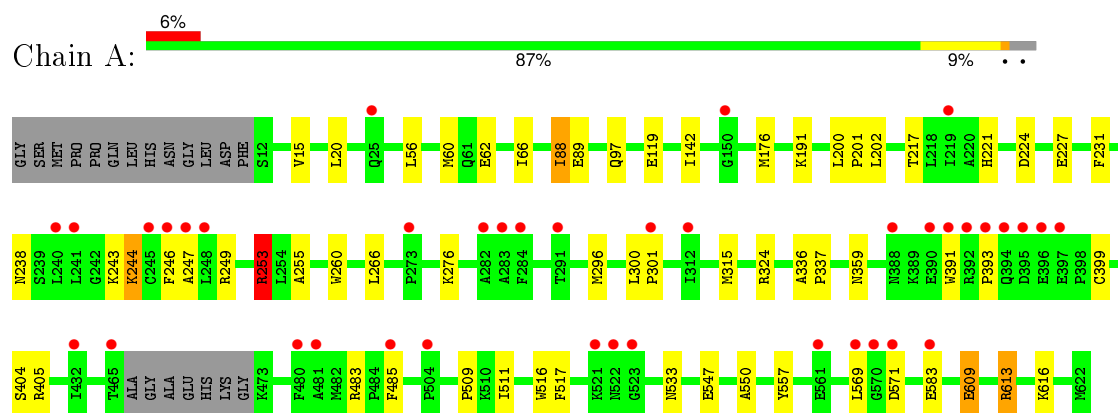
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	370	Total	O	0	0
			370	370		
8	B	298	Total	O	0	0
			298	298		

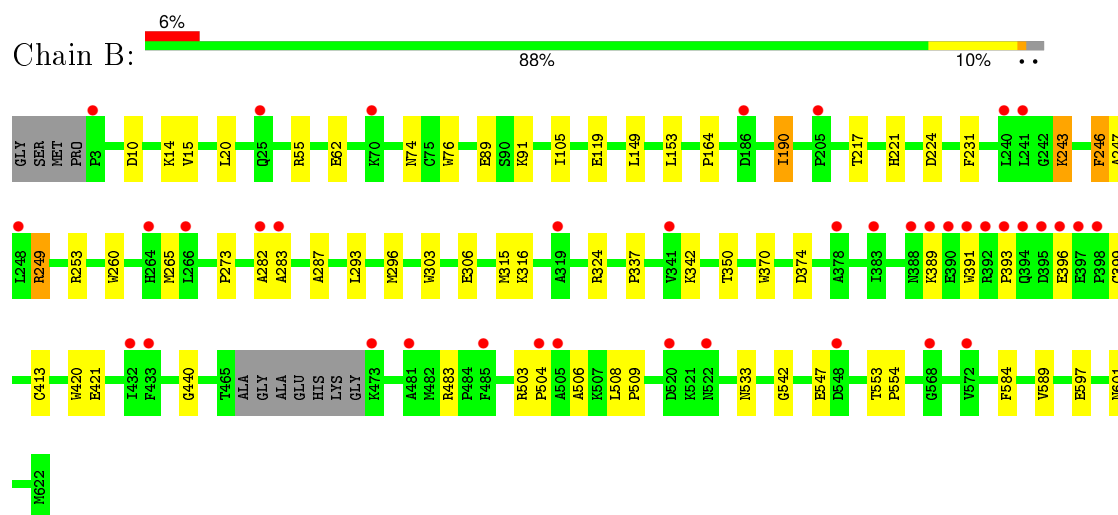
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ($\text{RSRZ} > 2$). 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, α , β , γ	62.08 Å 119.52 Å 87.02 Å 90.00° 107.15° 90.00°	Depositor
Resolution (Å)	35.93 – 2.10 35.93 – 2.10	Depositor EDS
% Data completeness (in resolution range)	96.8 (35.93-2.10) 96.8 (35.93-2.10)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.10 Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.190 , 0.247 0.205 , 0.260	Depositor DCC
R_{free} test set	3477 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	27.4	Xtriage
Anisotropy	0.056	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 33.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	2 of 68845 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10430	wwPDB-VP
Average B, all atoms (Å ²)	12.0	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 75.98 % 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 1.1384e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: GDP, PGA, NA, MN, EDO, 1PE

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	0.70	0/4965	0.71	3/6717 (0.0%)
1	B	0.68	0/4997	0.71	3/6765 (0.0%)
All	All	0.69	0/9962	0.71	6/13482 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	249	ARG	NE-CZ-NH2	-9.04	115.78	120.30
1	B	249	ARG	NE-CZ-NH1	6.73	123.66	120.30
1	A	613	ARG	NE-CZ-NH1	6.58	123.59	120.30
1	A	613	ARG	NE-CZ-NH2	-5.84	117.38	120.30
1	A	253	ARG	NE-CZ-NH2	-5.21	117.69	120.30
1	B	10	ASP	CB-CG-OD1	5.03	122.83	118.30

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	4811	0	4787	42	0
1	B	4852	0	4812	40	0
2	A	28	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	28	0	12	3	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	9	0	2	1	0
4	B	9	0	2	1	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	4	0	6	0	0
7	A	5	0	5	2	0
7	B	10	0	13	1	0
8	A	370	0	0	5	0
8	B	298	0	0	8	0
All	All	10430	0	9651	85	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 (85) 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:296:MET:HE2	1:A:533:ASN:HB2	1.59	0.83
1:A:119[B]:GLU:HG3	8:A:916:HOH:O	1.86	0.74
2:B:1000:GDP:O1B	8:B:746:HOH:O	2.11	0.69
1:B:89:GLU:OE2	1:B:483:ARG:NH1	2.28	0.66
1:B:224:ASP:OD2	8:B:626:HOH:O	2.14	0.65
1:B:190:ILE:HG21	1:B:246:PHE:CE2	2.36	0.60
1:B:287:ALA:HA	2:B:1000:GDP:H5'	1.83	0.60
1:A:509:PRO:O	7:A:1300:1PE:H122	2.02	0.60
1:A:511:ILE:H	7:A:1300:1PE:C23	2.14	0.59
1:B:91:LYS:HD3	1:B:217:THR:O	2.03	0.58
1:A:62[B]:GLU:OE1	1:A:62[B]:GLU:HA	2.03	0.58
1:A:217:THR:HA	1:A:231:PHE:O	2.04	0.58
1:B:14:LYS:NZ	7:B:1300:1PE:H122	2.20	0.57
1:A:315:MET:HA	1:A:324:ARG:O	2.05	0.57
1:B:62[B]:GLU:HG2	8:B:796:HOH:O	2.05	0.56
1:A:89:GLU:OE1	1:A:483:ARG:NH2	2.28	0.56
1:B:224:ASP:OD1	1:B:253:ARG:NH2	2.39	0.56
1:A:243:LYS:HG2	1:A:244:LYS:HD2	1.89	0.55
1:B:55:ARG:NH2	8:B:665:HOH:O	2.39	0.55
1:A:97[B]:GLN:OE1	1:A:119[B]:GLU:OE1	2.25	0.55
1:A:224:ASP:OD1	1:A:253:ARG:NH2	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:420:TRP:CE2	1:B:421:GLU:HG3	2.42	0.54
1:A:62[B]:GLU:HG2	8:A:664:HOH:O	2.08	0.54
1:A:244:LYS:HE2	1:A:485:PHE:CZ	2.44	0.53
1:A:296:MET:CE	1:A:533:ASN:HB2	2.35	0.53
1:B:503:ARG:HG2	1:B:506:ALA:HB2	1.91	0.52
1:A:255:ALA:HB1	1:A:260:TRP:O	2.10	0.52
1:A:244:LYS:HE2	1:A:485:PHE:CE1	2.44	0.52
1:B:217:THR:HA	1:B:231:PHE:O	2.10	0.51
1:A:609:GLU:OE1	1:A:613:ARG:HD2	2.10	0.51
1:B:243:LYS:O	1:B:247:ALA:HB3	2.11	0.51
1:B:153:LEU:HD11	1:B:316:LYS:HD3	1.93	0.51
1:B:391:TRP:CD2	1:B:399:CYS:HB3	2.46	0.51
1:A:391:TRP:CD2	1:A:399:CYS:HB3	2.47	0.50
1:B:164:PRO:HD2	8:B:705:HOH:O	2.12	0.49
1:B:293:LEU:HD12	1:B:296:MET:HE2	1.94	0.49
1:A:391:TRP:CE2	1:A:399:CYS:HB3	2.47	0.49
1:B:105:ILE:O	1:B:601[B]:ASN:ND2	2.46	0.47
1:A:276:LYS:HE3	8:B:658:HOH:O	2.14	0.47
2:B:1000:GDP:PB	8:B:746:HOH:O	2.73	0.47
1:B:15:VAL:HG21	1:B:20:LEU:HG	1.97	0.47
1:A:266:LEU:C	1:A:266:LEU:HD12	2.35	0.47
1:B:350:THR:OG1	1:B:413:CYS:HA	2.15	0.47
1:A:202:LEU:HD22	8:A:682:HOH:O	2.15	0.47
1:B:89:GLU:CD	1:B:483:ARG:HH12	2.19	0.46
1:A:88:ILE:H	1:A:238:ASN:HD21	1.63	0.46
1:B:553:THR:HB	1:B:554:PRO:HD2	1.98	0.46
1:A:15:VAL:HG21	1:A:20:LEU:HG	1.98	0.46
1:B:265:MET:HA	1:B:282:ALA:O	2.16	0.45
1:A:244:LYS:HG3	1:A:485:PHE:CE1	2.51	0.45
1:B:342:LYS:HE2	1:B:374:ASP:OD2	2.16	0.45
1:B:391:TRP:CE2	1:B:399:CYS:HB3	2.52	0.44
1:B:508:LEU:HA	1:B:509:PRO:HD2	1.88	0.44
1:A:336:ALA:HB3	1:A:337:PRO:HD3	1.98	0.44
1:A:547[A]:GLU:OE1	8:A:890:HOH:O	2.21	0.44
1:B:273:PRO:HD3	1:B:303:TRP:CE2	2.52	0.44
1:A:191:LYS:HB2	1:A:227:GLU:HG2	2.00	0.44
4:B:900:PGA:P	8:B:845:HOH:O	2.76	0.44
1:B:74:ASN:ND2	1:B:76:TRP:HE1	2.16	0.44
1:A:516:TRP:HB2	1:A:517:PHE:CE2	2.53	0.43
1:B:389:LYS:HA	1:B:389:LYS:HD3	1.84	0.43
1:B:296:MET:HE1	1:B:533:ASN:HB2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:TRP:HB3	1:B:316:LYS:HD2	1.99	0.43
1:A:200:LEU:HA	1:A:201:PRO:C	2.38	0.43
1:B:260:TRP:HB2	1:B:315:MET:O	2.19	0.43
1:A:89:GLU:OE2	1:A:483:ARG:NH1	2.40	0.42
1:B:503:ARG:HA	1:B:504:PRO:HD3	1.78	0.42
1:A:221:HIS:CE1	1:A:249:ARG:HD2	2.55	0.42
1:A:89:GLU:HB2	8:A:862:HOH:O	2.18	0.42
1:A:300:LEU:HA	1:A:301:PRO:HD3	1.95	0.42
1:A:616:LYS:HE3	1:A:616:LYS:HB2	1.79	0.42
1:B:315:MET:HA	1:B:324:ARG:O	2.20	0.42
1:A:550:ALA:HB1	1:A:557:TYR:HB3	2.02	0.42
1:B:221:HIS:CE1	1:B:249:ARG:HE	2.37	0.42
1:B:597:GLU:O	1:B:601[B]:ASN:HB3	2.20	0.41
1:A:359:ASN:O	1:A:404:SER:HB3	2.21	0.41
1:A:56:LEU:HD12	1:A:56:LEU:HA	1.89	0.41
1:B:265:MET:HG2	1:B:283:ALA:HB2	2.03	0.41
1:A:405:ARG:NH1	4:A:900:PGA:O2P	2.40	0.41
1:B:542:GLY:O	1:B:547:GLU:HB2	2.21	0.41
1:A:142:ILE:HG23	1:A:176:MET:CG	2.51	0.40
1:B:440:GLY:HA3	1:B:584:PHE:CZ	2.57	0.40
1:B:337:PRO:HA	1:B:370:TRP:CD1	2.56	0.40
1:A:243:LYS:O	1:A:247:ALA:HB3	2.22	0.40
1:A:60:MET:HB3	1:A:66:ILE:HG12	2.02	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	610/624 (98%)	590 (97%)	19 (3%)	1 (0%)	52	53
1	B	615/624 (99%)	587 (95%)	25 (4%)	3 (0%)	34	30

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1225/1248 (98%)	1177 (96%)	44 (4%)	4 (0%)	46	45

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	149	LEU
1	B	243	LYS
1	B	393	PRO
1	A	393	PRO

5.3.2 Protein sidechains ⓘ

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	516/520 (99%)	508 (98%)	8 (2%)	70	76
1	B	520/520 (100%)	514 (99%)	6 (1%)	78	84
All	All	1036/1040 (100%)	1022 (99%)	14 (1%)	74	80

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	88	ILE
1	A	244	LYS
1	A	246	PHE
1	A	253	ARG
1	A	569	LEU
1	A	571	ASP
1	A	583	GLU
1	A	609	GLU
1	B	119	GLU
1	B	190	ILE
1	B	246	PHE
1	B	306	GLU
1	B	396	GLU
1	B	589	VAL

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

Mol	Chain	Res	Type
1	A	238	ASN
1	B	74	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 13 ligands modelled in this entry, 6 are monoatomic - leaving 7 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	GDP	A	1000	3	23,30,30	1.18	2 (8%)	30,47,47	1.73	9 (30%)
6	EDO	A	1200	-	3,3,3	0.91	0	2,2,2	0.47	0
7	1PE	A	1300	-	4,4,15	0.62	0	3,3,14	0.63	0
4	PGA	A	900	3	5,8,8	0.69	0	6,11,11	1.22	1 (16%)
2	GDP	B	1000	3	23,30,30	1.41	3 (13%)	30,47,47	1.80	7 (23%)
7	1PE	B	1300	-	9,9,15	0.69	0	8,8,14	0.58	0
4	PGA	B	900	3	5,8,8	0.74	0	6,11,11	1.13	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	GDP	A	1000	3	-	0/12/32/32	0/3/3/3
6	EDO	A	1200	-	-	0/1/1/1	0/0/0/0
7	1PE	A	1300	-	-	0/2/2/13	0/0/0/0
4	PGA	A	900	3	-	0/4/6/6	0/0/0/0
2	GDP	B	1000	3	-	0/12/32/32	0/3/3/3
7	1PE	B	1300	-	-	0/7/7/13	0/0/0/0
4	PGA	B	900	3	-	0/4/6/6	0/0/0/0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1000	GDP	O4'-C1'	2.24	1.44	1.41
2	A	1000	GDP	C5-C4	2.48	1.46	1.40
2	B	1000	GDP	C6-C5	3.00	1.47	1.41
2	A	1000	GDP	C6-C5	3.45	1.48	1.41
2	B	1000	GDP	C5-C4	3.84	1.49	1.40

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1000	GDP	C2'-C1'-N9	-4.99	106.67	114.29
2	B	1000	GDP	C5-C6-N1	-3.93	118.21	123.59
2	A	1000	GDP	C5-C6-N1	-3.84	118.34	123.59
2	A	1000	GDP	C6-C5-C4	-2.93	117.40	120.90
2	A	1000	GDP	N3-C2-N1	-2.92	122.99	127.44
2	A	1000	GDP	C4-C5-N7	-2.58	107.11	109.48
2	B	1000	GDP	N3-C2-N1	-2.52	123.61	127.44
2	B	1000	GDP	O3'-C3'-C2'	-2.51	103.67	111.83
2	A	1000	GDP	C1'-N9-C4	-2.45	123.25	126.94
2	B	1000	GDP	PA-O3A-PB	-2.37	124.71	132.67
2	B	1000	GDP	C6-C5-C4	-2.32	118.13	120.90
2	A	1000	GDP	PA-O3A-PB	-2.19	125.32	132.67
4	A	900	PGA	O4P-P-O3P	2.15	115.58	107.38
2	A	1000	GDP	O2A-PA-O3A	2.27	115.38	105.09
2	A	1000	GDP	N2-C2-N1	2.35	121.09	117.20
2	A	1000	GDP	C6-N1-C2	3.71	121.08	115.94
2	B	1000	GDP	C6-N1-C2	3.93	121.39	115.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	1300	1PE	2	0
4	A	900	PGA	1	0
2	B	1000	GDP	3	0
7	B	1300	1PE	1	0
4	B	900	PGA	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	604/624 (96%)	0.35	39 (6%) 22 29	6, 12, 16, 21	0
1	B	613/624 (98%)	0.39	39 (6%) 23 30	7, 13, 17, 21	0
All	All	1217/1248 (97%)	0.37	78 (6%) 23 30	6, 12, 17, 21	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	394	GLN	8.5
1	B	394	GLN	6.5
1	A	396	GLU	5.6
1	A	504	PRO	5.3
1	B	392	ARG	5.1
1	A	395	ASP	4.9
1	A	571	ASP	4.5
1	B	393	PRO	4.5
1	B	504	PRO	4.3
1	A	392	ARG	4.2
1	A	248	LEU	3.9
1	B	396	GLU	3.9
1	B	572	VAL	3.7
1	B	391	TRP	3.6
1	A	397	GLU	3.6
1	B	341	VAL	3.6
1	B	473	LYS	3.4
1	B	3	PRO	3.4
1	B	240	LEU	3.3
1	B	205	PRO	3.3
1	A	245	CYS	3.2
1	A	481	ALA	3.2
1	B	505	ALA	3.2
1	B	388	ASN	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	240	LEU	3.2
1	B	481	ALA	3.1
1	A	241	LEU	3.1
1	A	312	ILE	3.0
1	A	521	LYS	3.0
1	A	273	PRO	3.0
1	A	465	THR	2.9
1	A	391	TRP	2.9
1	B	520	ASP	2.8
1	B	395	ASP	2.8
1	B	186	ASP	2.7
1	A	570	GLY	2.7
1	B	390	GLU	2.7
1	A	283	ALA	2.6
1	A	246	PHE	2.6
1	A	569	LEU	2.6
1	B	383	ILE	2.6
1	B	248	LEU	2.6
1	A	247	ALA	2.6
1	B	522	ASN	2.6
1	B	397	GLU	2.5
1	A	523	GLY	2.5
1	B	241	LEU	2.5
1	B	398	PRO	2.5
1	B	264	HIS	2.5
1	B	266	LEU	2.5
1	B	319	ALA	2.4
1	A	25	GLN	2.4
1	A	522	ASN	2.4
1	A	284	PHE	2.4
1	B	433	PHE	2.4
1	A	388	ASN	2.3
1	B	70	LYS	2.3
1	A	480	PHE	2.3
1	A	485	PHE	2.3
1	A	583	GLU	2.3
1	B	283	ALA	2.3
1	A	432	ILE	2.3
1	B	25	GLN	2.3
1	B	282	ALA	2.3
1	B	378	ALA	2.2
1	B	485	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	548	ASP	2.1
1	A	150	GLY	2.1
1	A	219	ILE	2.1
1	B	432	ILE	2.1
1	A	390	GLU	2.1
1	A	282	ALA	2.1
1	A	291	THR	2.0
1	A	393	PRO	2.0
1	A	561	GLU	2.0
1	B	389	LYS	2.0
1	A	301	PRO	2.0
1	B	568	GLY	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 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. 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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
7	1PE	A	1300	5/16	0.87	0.24	5.21	17,20,24,25	0
7	1PE	B	1300	10/16	0.84	0.18	1.07	23,28,29,30	0
2	GDP	A	1000	28/28	0.97	0.09	-1.01	14,17,19,23	0
5	NA	B	1100	1/1	0.97	0.09	-1.11	24,24,24,24	0
2	GDP	B	1000	28/28	0.96	0.09	-1.28	15,18,21,22	0
4	PGA	B	900	9/9	0.96	0.10	-1.84	19,20,21,21	0
4	PGA	A	900	9/9	0.97	0.10	-2.28	15,16,18,18	0
5	NA	A	1100	1/1	0.95	0.06	-4.38	22,22,22,22	0
3	MN	A	700	1/1	0.99	0.05	-4.46	18,18,18,18	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	MN	B	701	1/1	0.99	0.03	-4.71	20,20,20,20	0
3	MN	A	701	1/1	0.99	0.04	-6.21	19,19,19,19	0
3	MN	B	700	1/1	0.99	0.03	-7.08	19,19,19,19	0
6	EDO	A	1200	4/4	0.84	0.15	-	25,27,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.