



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 11:35 PM GMT

PDB ID : 2POC
Title : The crystal structure of isomerase domain of glucosamine-6-phosphatesynthase from *Candida albicans*
Authors : Raczynska, J.; Olchow, J.; Milewski, S.; Rypniewski, W.
Deposited on : 2007-04-26
Resolution : 1.80 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

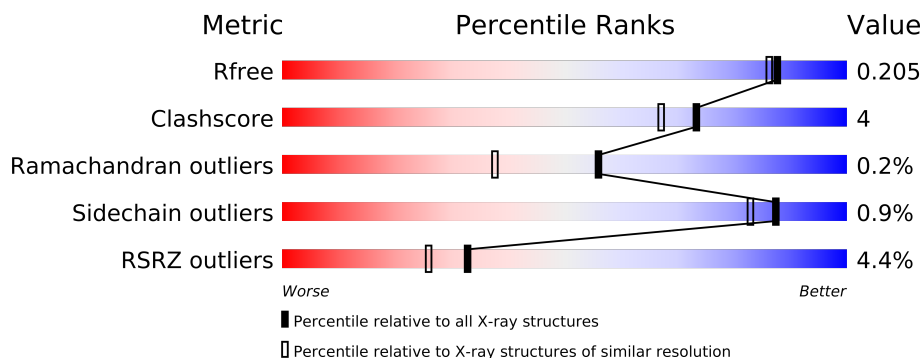
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.80 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}	66092	3513 (1.80-1.80)
Clashscore	79885	4461 (1.80-1.80)
Ramachandran outliers	78287	4404 (1.80-1.80)
Sidechain outliers	78261	4403 (1.80-1.80)
RSRZ outliers	66119	3515 (1.80-1.80)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	367	
1	B	367	
1	C	367	
1	D	367	

2 Entry composition i

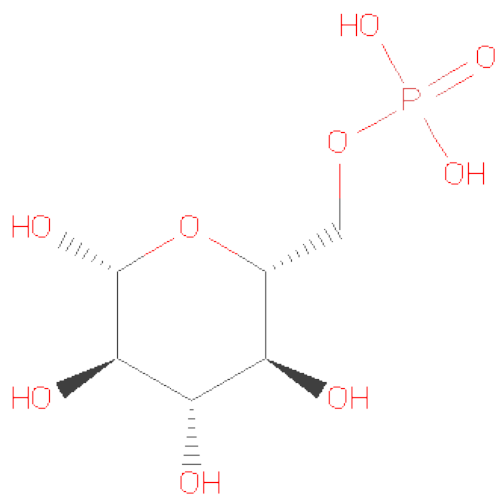
There are 6 unique types of molecules in this entry. The entry contains 11928 atoms, of which 0 are hydrogen and 0 are deuterium.

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 isomerase domain of glutamine-fructose-6-phosphatetransaminase (isomerizing).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	339	Total	C	N	O	S	0	4	0
			2647	1671	462	497	17			
1	B	352	Total	C	N	O	S	0	6	0
			2753	1740	473	522	18			
1	C	352	Total	C	N	O	S	0	2	0
			2739	1729	473	519	18			
1	D	340	Total	C	N	O	S	0	5	0
			2665	1681	461	506	17			

- Molecule 2 is SUGAR (BETA-D-GLUCOSE-6-PHOSPHATE) (three-letter code: BG6) (formula: C₆H₁₃O₉P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			16	6	9	1		

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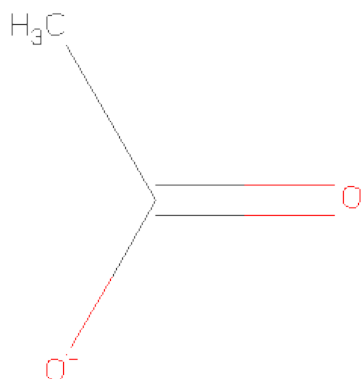
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	O	P	0	0
			16	6	9	1		
2	C	1	Total	C	O	P	0	0
			16	6	9	1		
2	D	1	Total	C	O	P	0	0
			16	6	9	1		

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

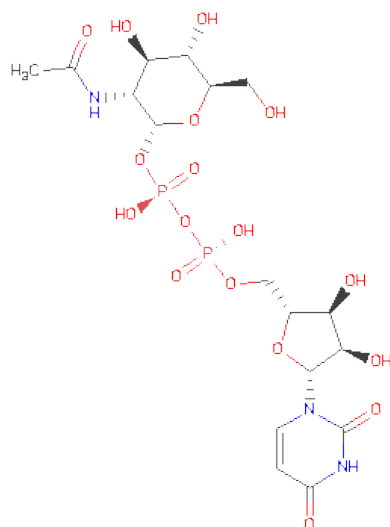
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Na	0	0
			1	1		
3	A	1	Total	Na	0	0
			1	1		
3	D	1	Total	Na	0	0
			1	1		
3	C	1	Total	Na	0	0
			1	1		

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



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

- Molecule 5 is URIDINE-DIPHOSPHATE-N-ACETYLGLUCOSAMINE (three-letter code: UD1) (formula: $C_{17}H_{27}N_3O_{17}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			39	17	3	17	2		
5	B	1	Total	C	N	O	P	0	0
			39	17	3	17	2		
5	C	1	Total	C	N	O	P	0	0
			39	17	3	17	2		
5	D	1	Total	C	N	O	P	0	0
			39	17	3	17	2		

- Molecule 6 is water.

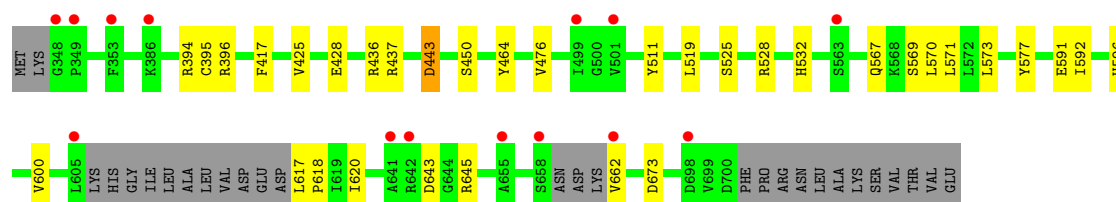
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	199	Total	O	0	0
			199	199		
6	B	234	Total	O	0	0
			234	234		
6	C	231	Total	O	0	0
			231	231		
6	D	228	Total	O	0	0
			228	228		

3 Residue-property plots

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 ($RSRZ > 2$). 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.

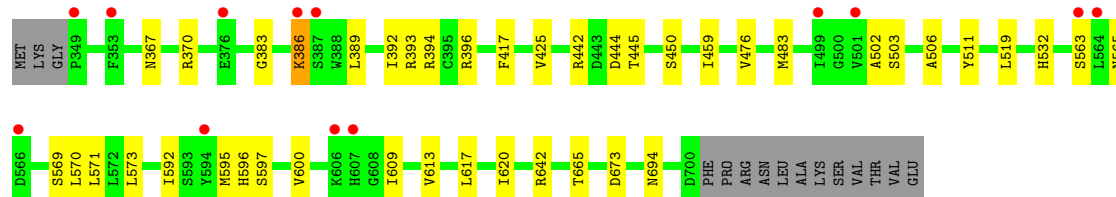
- Molecule 1: isomerase domain of glutamine-fructose-6-phosphatetransaminase (isomerizing)

Chain A: 



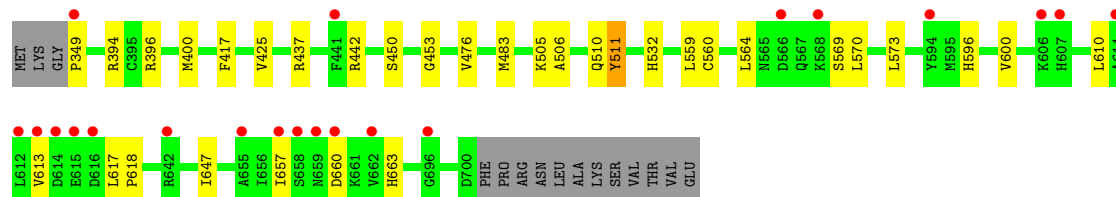
- Molecule 1: isomerase domain of glutamine-fructose-6-phosphatetransaminase (isomerizing)

Chain B: 



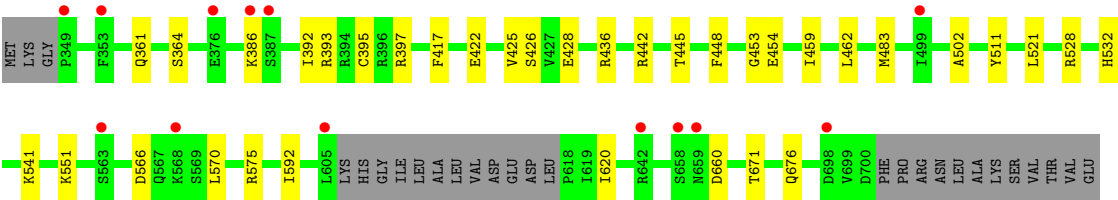
- Molecule 1: isomerase domain of glutamine-fructose-6-phosphatetransaminase (isomerizing)

Chain C: 



- Molecule 1: isomerase domain of glutamine-fructose-6-phosphatetransaminase (isomerizing)

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	65.98Å 117.83Å 99.71Å 90.00° 91.60° 90.00°	Depositor
Resolution (Å)	19.94 – 1.80 19.93 – 1.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.94-1.80) 99.9 (19.93-1.80)	Depositor EDS
R_{merge}	0.00	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.02 (at 1.80Å)	Xtriage
Refinement program	REFMAC 5.2	Depositor
R, R_{free}	0.175 , 0.206 0.174 , 0.205	Depositor DCC
R_{free} test set	2819 reflections (2.05%)	DCC
Wilson B-factor (Å ²)	22.7	Xtriage
Anisotropy	0.040	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 51.9	EDS
Estimated twinning fraction	0.035 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 140518 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11928	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.98% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA, UD1, BG6, ACT

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	0.82	1/2704 (0.0%)	0.82	2/3652 (0.1%)
1	B	0.75	0/2819	0.79	2/3810 (0.1%)
1	C	0.77	0/2787	0.79	2/3766 (0.1%)
1	D	0.82	0/2721	0.83	2/3674 (0.1%)
All	All	0.79	1/11031 (0.0%)	0.81	8/14902 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	577	TYR	CD1-CE1	5.09	1.47	1.39

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	673	ASP	CB-CG-OD2	-5.84	113.04	118.30
1	D	575	ARG	NE-CZ-NH2	-5.83	117.39	120.30
1	B	673	ASP	CB-CG-OD2	-5.73	113.14	118.30
1	C	570	LEU	CA-CB-CG	5.68	128.35	115.30
1	A	443	ASP	CB-CG-OD1	5.65	123.39	118.30
1	B	673	ASP	CB-CG-OD1	5.48	123.23	118.30
1	D	575	ARG	NE-CZ-NH1	5.32	122.96	120.30
1	C	511	TYR	CB-CG-CD2	-5.15	117.91	121.00

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2647	0	2689	24	0
1	B	2753	0	2807	32	0
1	C	2739	0	2794	22	0
1	D	2665	0	2718	20	0
2	A	16	0	11	0	0
2	B	16	0	11	0	0
2	C	16	0	11	0	0
2	D	16	0	11	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	B	4	0	3	0	0
4	C	4	0	3	0	0
5	A	39	0	25	1	0
5	B	39	0	25	0	0
5	C	39	0	25	0	0
5	D	39	0	25	1	0
6	A	199	0	0	2	0
6	B	234	0	0	4	0
6	C	231	0	0	10	0
6	D	228	0	0	4	0
All	All	11928	0	11158	94	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 4.

All (94) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:459[A]:ILE:HD13	1:B:483[A]:MET:SD	1.98	1.02
1:C:400:MET:SD	6:C:5190:HOH:O	2.41	0.78
1:A:596[B]:HIS:NE2	1:B:609:ILE:HD12	1.98	0.78
1:B:459[A]:ILE:CD1	1:B:483[A]:MET:SD	2.74	0.76
1:B:571:LEU:HG	1:B:617:LEU:HD21	1.68	0.76

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:428:GLU:OE1	1:A:436[A]:ARG:NH2	2.21	0.73
1:D:428:GLU:OE1	1:D:436:ARG:NH2	2.22	0.72
1:D:426:SER:HB3	6:D:5220:HOH:O	1.90	0.70
1:A:571:LEU:HD11	1:A:617:LEU:HD21	1.75	0.68
1:C:394:ARG:HB2	6:C:5210:HOH:O	1.94	0.68
1:C:349:PRO:N	6:C:5153:HOH:O	2.27	0.67
1:D:502:ALA:HB2	1:D:592:ILE:HD11	1.76	0.67
1:A:395:CYS:HB3	1:C:442[B]:ARG:NH2	2.09	0.66
1:C:505:LYS:HB3	6:C:5234:HOH:O	1.96	0.65
1:D:502:ALA:CB	1:D:592:ILE:HD11	2.28	0.64
1:B:367:ASN:OD1	1:B:370:ARG:NH2	2.30	0.64
1:A:643:ASP:HB2	6:A:5193:HOH:O	1.99	0.62
1:D:393:ARG:HD2	1:D:521:LEU:O	2.00	0.62
1:A:443:ASP:CG	6:A:5200:HOH:O	2.38	0.61
1:D:453:GLY:O	1:D:483:MET:HG3	2.01	0.61
1:B:389:LEU:O	1:B:393:ARG:HG3	2.02	0.60
1:A:596[B]:HIS:HE2	1:B:609:ILE:HD12	1.67	0.59
1:B:613:VAL:O	1:B:642:ARG:NH2	2.34	0.58
1:B:459[A]:ILE:HD11	1:B:483[A]:MET:HG2	1.87	0.57
1:C:394:ARG:HB3	6:C:5151:HOH:O	2.04	0.57
1:A:569:SER:O	1:A:618:PRO:HD2	2.04	0.57
1:A:591:GLU:HG3	1:A:592:ILE:HG12	1.85	0.56
1:B:396:ARG:HD3	6:B:5071:HOH:O	2.04	0.56
1:B:502:ALA:CB	1:B:592:ILE:HD11	2.35	0.56
1:D:551:LYS:HE3	6:D:5161:HOH:O	2.06	0.55
1:B:565:ASN:OD1	1:B:694:ASN:HB3	2.08	0.54
1:A:571:LEU:CD1	1:A:617:LEU:HD21	2.37	0.54
1:A:567:GLN:HG3	1:A:618:PRO:HG3	1.90	0.54
1:C:505:LYS:CB	6:C:5234:HOH:O	2.55	0.53
1:B:502:ALA:HB2	1:B:592:ILE:HD11	1.92	0.52
1:B:442[B]:ARG:NH2	1:D:395:CYS:HB3	2.24	0.52
1:C:657:ILE:HD12	1:C:663:HIS:CG	2.43	0.52
1:B:573:LEU:HD23	1:B:600:VAL:HB	1.91	0.52
1:B:503:SER:OG	1:B:506:ALA:HB3	2.10	0.51
1:C:417:PHE:HB3	1:C:425:VAL:HG21	1.92	0.51
1:A:570:LEU:HD23	1:A:570:LEU:C	2.32	0.50
1:B:571:LEU:HD21	1:B:609:ILE:HD11	1.95	0.48
1:A:519:LEU:O	1:A:532[A]:HIS:HE1	1.96	0.47
1:B:450:SER:O	1:B:476:VAL:HA	2.15	0.47
1:A:570:LEU:HD21	1:A:620:ILE:HD12	1.96	0.47
1:C:506:ALA:O	1:C:510:GLN:HG3	2.14	0.47
1:D:417:PHE:HB3	1:D:425:VAL:HG21	1.96	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:541:LYS:HE3	6:D:5127:HOH:O	2.14	0.47
1:B:442[B]:ARG:NH2	1:D:445:THR:OG1	2.37	0.47
1:B:665[A]:THR:HG21	6:B:5163:HOH:O	2.14	0.47
1:C:617:LEU:HD12	1:C:618:PRO:HD2	1.97	0.47
1:D:570:LEU:HD21	1:D:620:ILE:HD12	1.97	0.47
1:A:567:GLN:HG3	1:A:618:PRO:CG	2.45	0.46
1:A:643:ASP:OD2	1:A:645:ARG:NH1	2.49	0.46
1:A:450:SER:O	1:A:476:VAL:HA	2.16	0.45
5:A:5002:UD1:H5'2	5:A:5002:UD1:H8'2	1.98	0.45
1:C:613:VAL:HB	1:C:617:LEU:HD23	1.99	0.45
1:B:642:ARG:HG3	6:B:5154:HOH:O	2.16	0.45
1:A:525:SER:OG	1:A:528:ARG:HD3	2.17	0.44
1:B:444:ASP:O	1:D:442[B]:ARG:NH2	2.50	0.44
1:A:417:PHE:HB3	1:A:425:VAL:HG21	1.99	0.44
1:B:417:PHE:HB3	1:B:425:VAL:HG21	1.99	0.44
1:C:560:CYS:HA	1:C:564:LEU:HB2	1.98	0.44
1:C:573:LEU:HD23	1:C:600:VAL:HB	1.99	0.44
1:B:519:LEU:O	1:B:532:HIS:HE1	2.01	0.44
1:C:400:MET:HE1	6:C:5190:HOH:O	2.17	0.43
1:D:454:GLU:HA	1:D:459:ILE:HD11	2.00	0.43
1:C:559:LEU:HD21	1:C:647:ILE:HD13	2.01	0.43
1:B:392:ILE:HG23	1:B:445:THR:HG21	2.00	0.43
1:B:665[A]:THR:HG23	6:B:5161:HOH:O	2.18	0.43
1:B:570:LEU:HD21	1:B:620:ILE:HD12	2.00	0.43
1:A:573:LEU:HD23	1:A:600:VAL:HB	2.01	0.43
1:A:394:ARG:HG2	1:A:394:ARG:O	2.18	0.43
1:D:361:GLN:HA	1:D:364:SER:OG	2.19	0.43
1:C:453:GLY:O	1:C:483[B]:MET:HG2	2.18	0.42
1:A:645:ARG:HE	1:A:662:VAL:HG11	1.84	0.42
1:C:610:LEU:HA	1:C:613:VAL:HG13	2.02	0.42
1:D:660:ASP:O	6:D:5174:HOH:O	2.21	0.42
1:B:570:LEU:O	1:B:597:SER:HA	2.20	0.42
1:C:396:ARG:HD3	6:C:5029:HOH:O	2.18	0.42
1:A:396:ARG:HD3	6:C:5155:HOH:O	2.19	0.42
1:B:609:ILE:O	1:B:613:VAL:HG13	2.19	0.41
1:A:437:ARG:HH21	1:A:464:TYR:HE2	1.68	0.41
1:D:392:ILE:HG23	1:D:445:THR:HG21	2.03	0.41
1:B:383:GLY:O	1:B:386:LYS:HB2	2.20	0.41
5:D:5005:UD1:H5'2	5:D:5005:UD1:H8'2	2.01	0.41
1:D:671:THR:CG2	1:D:676:GLN:HA	2.51	0.41
1:C:450:SER:O	1:C:476:VAL:HA	2.21	0.41
1:C:569:SER:HA	1:C:596:HIS:O	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:569:SER:HA	1:B:596:HIS:O	2.21	0.41
1:B:595:MET:HE2	1:B:595:MET:HB2	1.90	0.41
1:D:448:PHE:CZ	1:D:462:LEU:HA	2.56	0.40
1:D:422:GLU:OE2	1:D:528:ARG:NH2	2.52	0.40
1:C:400:MET:CE	6:C:5190:HOH:O	2.69	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	337/367 (92%)	332 (98%)	5 (2%)	0	100	100
1	B	356/367 (97%)	349 (98%)	5 (1%)	2 (1%)	33	15
1	C	352/367 (96%)	349 (99%)	3 (1%)	0	100	100
1	D	341/367 (93%)	337 (99%)	3 (1%)	1 (0%)	50	31
All	All	1386/1468 (94%)	1367 (99%)	16 (1%)	3 (0%)	56	38

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	386	LYS
1	D	386	LYS
1	B	394	ARG

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	298/319 (93%)	297 (100%)	1 (0%)	96	94
1	B	312/319 (98%)	310 (99%)	2 (1%)	92	90
1	C	308/319 (97%)	304 (99%)	4 (1%)	80	71
1	D	301/319 (94%)	297 (99%)	4 (1%)	80	71
All	All	1219/1276 (96%)	1208 (99%)	11 (1%)	87	83

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	511	TYR
1	B	511	TYR
1	B	563	SER
1	C	437	ARG
1	C	511	TYR
1	C	532	HIS
1	C	660	ASP
1	D	397	ARG
1	D	511	TYR
1	D	532	HIS
1	D	566	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	561	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 14 ligands modelled in this entry, 4 are monoatomic - leaving 10 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	UD1	A	5002	3	41,41,41	0.90	3 (7%)	58,62,62	1.14	2 (3%)
2	BG6	A	713	-	16,16,16	1.09	2 (12%)	24,24,24	1.24	3 (12%)
5	UD1	B	5003	3	41,41,41	1.06	4 (9%)	58,62,62	1.31	6 (10%)
2	BG6	B	713	-	16,16,16	0.85	0	24,24,24	1.07	2 (8%)
4	ACT	B	714	-	1,3,3	1.55	0	0,3,3	0.00	-
5	UD1	C	5004	3	41,41,41	0.86	1 (2%)	58,62,62	1.20	3 (5%)
2	BG6	C	713	-	16,16,16	1.23	2 (12%)	24,24,24	0.66	0
4	ACT	C	714	-	1,3,3	1.80	0	0,3,3	0.00	-
5	UD1	D	5005	3	41,41,41	0.83	2 (4%)	58,62,62	1.09	3 (5%)
2	BG6	D	713	-	16,16,16	1.27	2 (12%)	24,24,24	0.78	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
5	UD1	A	5002	3	-	0/25/63/63	0/3/3/3
2	BG6	A	713	-	1/1/6/6	0/6/26/26	0/1/1/1
5	UD1	B	5003	3	-	0/25/63/63	0/3/3/3
2	BG6	B	713	-	1/1/6/6	0/6/26/26	0/1/1/1
4	ACT	B	714	-	-	0/0/0/0	0/0/0/0
5	UD1	C	5004	3	-	0/25/63/63	0/3/3/3
2	BG6	C	713	-	1/1/6/6	0/6/26/26	0/1/1/1
4	ACT	C	714	-	-	0/0/0/0	0/0/0/0
5	UD1	D	5005	3	-	0/25/63/63	0/3/3/3
2	BG6	D	713	-	1/1/6/6	0/6/26/26	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	713	BG6	O5-C1	3.48	1.50	1.43
2	D	713	BG6	P-O3P	3.24	1.62	1.51
5	A	5002	UD1	O4B-C1B	3.13	1.46	1.41
5	B	5003	UD1	C6-C5	3.07	1.41	1.36
5	B	5003	UD1	C2-N1	2.98	1.41	1.38
5	A	5002	UD1	C6-C5	2.95	1.40	1.36
2	D	713	BG6	O1-C1	2.72	1.49	1.39
2	C	713	BG6	P-O3P	2.65	1.60	1.51
5	D	5005	UD1	O4B-C1B	2.61	1.45	1.41
5	B	5003	UD1	C2B-C1B	-2.52	1.49	1.53
5	B	5003	UD1	O4B-C1B	2.52	1.45	1.41
2	A	713	BG6	O5-C1	2.23	1.47	1.43
5	A	5002	UD1	C2B-C1B	-2.21	1.50	1.53
2	A	713	BG6	P-O2P	2.17	1.62	1.54
5	D	5005	UD1	C2B-C1B	-2.13	1.50	1.53
5	C	5004	UD1	O4B-C1B	2.00	1.44	1.41

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	5002	UD1	N3-C2-N1	5.44	120.51	115.97
5	D	5005	UD1	N3-C2-N1	4.18	119.46	115.97
5	B	5003	UD1	N3-C2-N1	4.17	119.46	115.97
5	C	5004	UD1	N3-C2-N1	3.93	119.25	115.97
5	C	5004	UD1	O5'-C1'-O1'	-3.38	106.95	111.36
5	C	5004	UD1	C2-N1-C1B	3.31	120.28	118.21
5	B	5003	UD1	C3'-C4'-C5'	3.29	116.08	110.20
5	B	5003	UD1	O5'-C5'-C4'	3.28	115.83	109.76
5	B	5003	UD1	C4'-C3'-C2'	3.10	114.84	110.44
2	B	713	BG6	O5-C1-C2	-2.83	105.48	109.86
5	B	5003	UD1	C5-C4-N3	2.77	121.92	116.70
2	A	713	BG6	O1-C1-C2	-2.58	102.16	109.47
5	A	5002	UD1	O4B-C1B-N1	-2.47	102.86	108.06
2	A	713	BG6	O1P-P-O6	2.33	113.07	106.65
5	D	5005	UD1	O5'-C5'-C4'	2.30	114.02	109.76
2	A	713	BG6	O2P-P-O3P	-2.12	103.50	110.44
2	B	713	BG6	C3-C4-C5	-2.10	106.45	110.20
5	D	5005	UD1	C6'-C5'-C4'	-2.05	108.04	113.00
5	B	5003	UD1	O3'-C3'-C4'	-2.01	105.85	110.35

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	D	713	BG6	C1
2	C	713	BG6	C1
2	A	713	BG6	C1
2	B	713	BG6	C1

There are no torsion outliers.

There are no ring outliers.

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, 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	339/367 (92%)	-0.05	14 (4%) 35 28	14, 24, 41, 53	0
1	B	352/367 (95%)	-0.05	13 (3%) 39 32	15, 24, 47, 59	0
1	C	352/367 (95%)	0.06	21 (5%) 21 16	14, 25, 52, 65	0
1	D	340/367 (92%)	-0.08	13 (3%) 38 31	14, 24, 39, 49	0
All	All	1383/1468 (94%)	-0.03	61 (4%) 33 26	14, 24, 46, 65	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	659	ASN	5.7
1	B	563	SER	5.2
1	A	501	VAL	5.0
1	C	660	ASP	4.9
1	C	607	HIS	4.9
1	C	615	GLU	4.9
1	C	616	ASP	4.5
1	A	348	GLY	4.3
1	A	349	PRO	4.2
1	D	386	LYS	4.2
1	B	607	HIS	4.0
1	A	658	SER	3.8
1	C	655	ALA	3.8
1	C	613	VAL	3.7
1	A	499	ILE	3.7
1	B	566	ASP	3.7
1	D	349	PRO	3.6
1	D	659	ASN	3.5
1	C	611	ALA	3.4
1	A	642	ARG	3.4
1	C	606	LYS	3.3

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Mol	Chain	Res	Type	RSRZ
1	C	658	SER	3.3
1	C	662	VAL	3.2
1	B	386	LYS	3.2
1	B	594	TYR	3.1
1	C	594	TYR	3.0
1	D	376	GLU	3.0
1	C	568	LYS	3.0
1	A	662	VAL	3.0
1	C	642	ARG	2.9
1	A	605	LEU	2.9
1	C	566	ASP	2.9
1	C	349	PRO	2.8
1	A	353	PHE	2.8
1	B	606	LYS	2.7
1	B	387	SER	2.6
1	D	658	SER	2.6
1	B	499	ILE	2.6
1	A	563	SER	2.5
1	A	386	LYS	2.5
1	B	353	PHE	2.4
1	D	568	LYS	2.4
1	D	563	SER	2.4
1	C	612	LEU	2.3
1	D	387	SER	2.3
1	A	641	ALA	2.3
1	C	657	ILE	2.3
1	D	353	PHE	2.2
1	C	614	ASP	2.2
1	B	564	LEU	2.2
1	D	605	LEU	2.2
1	D	499	ILE	2.2
1	D	698	ASP	2.1
1	C	696	GLY	2.1
1	D	642	ARG	2.1
1	A	655	ALA	2.1
1	B	376	GLU	2.1
1	C	441	PHE	2.1
1	B	501	VAL	2.0
1	B	349	PRO	2.0
1	A	698	ASP	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	BG6	A	713	16/16	0.14	0.52	21,28,40,41	0
5	UD1	B	5003	39/39	0.12	0.23	18,26,58,59	0
5	UD1	D	5005	39/39	0.13	0.21	21,26,47,51	0
2	BG6	C	713	16/16	0.08	0.01	19,24,38,38	0
5	UD1	A	5002	39/39	0.12	-0.03	20,25,46,49	0
3	NA	A	5001	1/1	0.11	-0.08	26,26,26,26	0
2	BG6	D	713	16/16	0.09	-0.26	20,28,40,41	0
2	BG6	B	713	16/16	0.08	-0.33	17,22,35,38	0
4	ACT	C	714	4/4	0.13	-0.34	40,40,41,41	0
5	UD1	C	5004	39/39	0.07	-0.50	18,21,28,33	0
4	ACT	B	714	4/4	0.12	-0.56	34,35,35,35	0
3	NA	D	5004	1/1	0.07	-1.23	27,27,27,27	0
3	NA	C	5003	1/1	0.04	-2.80	21,21,21,21	0
3	NA	B	5002	1/1	0.04	-4.21	21,21,21,21	0

6.5 Other polymers ⓘ

There are no such residues in this entry.