



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:31 PM UTC

PDB ID : 6MGI / pdb_00006mgi
Title : Photosynthetic phosphoenolpyruvate carboxylase isoenzyme from maize complexed with the allosteric activator glucose-6-phosphate in its allosteric site
Authors : Gonzalez-Segura, L.; Munoz-Clares, R.A.
Deposited on : 2018-09-13
Resolution : 2.99 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

<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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

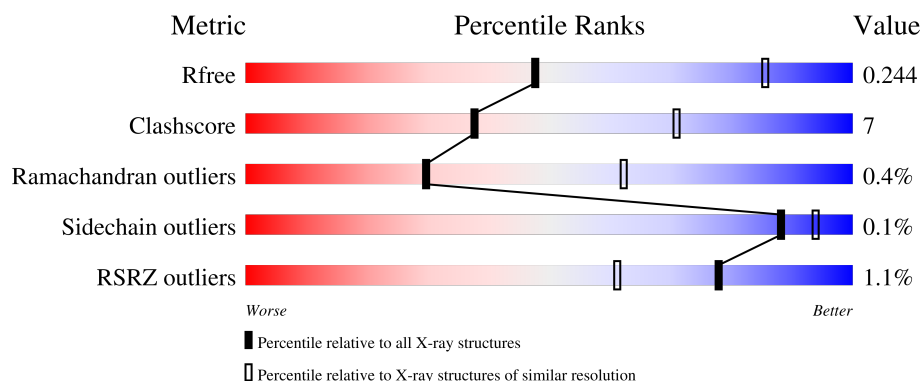
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.99 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}	180053	3580 (3.00-2.96)
Clashscore	190562	3904 (3.00-2.96)
Ramachandran outliers	187476	3761 (3.00-2.96)
Sidechain outliers	187428	3764 (3.00-2.96)
RSRZ outliers	180081	3579 (3.00-2.96)

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 ≥ 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	970	
1	B	970	

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
2	GLY	B	1001	-	X	-	-

2 Entry composition [i](#)

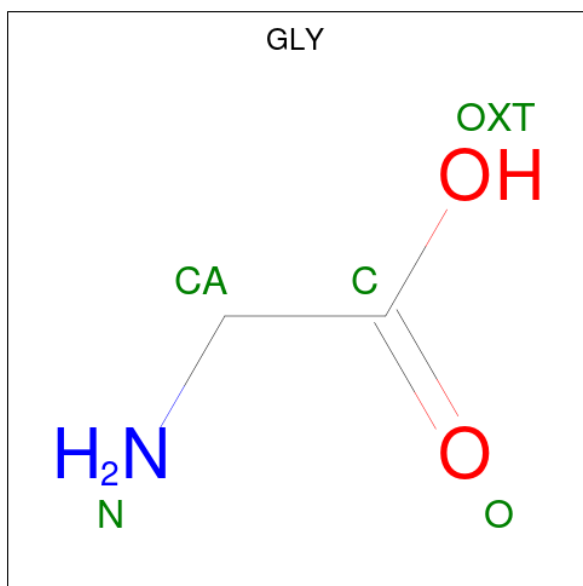
There are 7 unique types of molecules in this entry. The entry contains 14680 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 carboxylase.

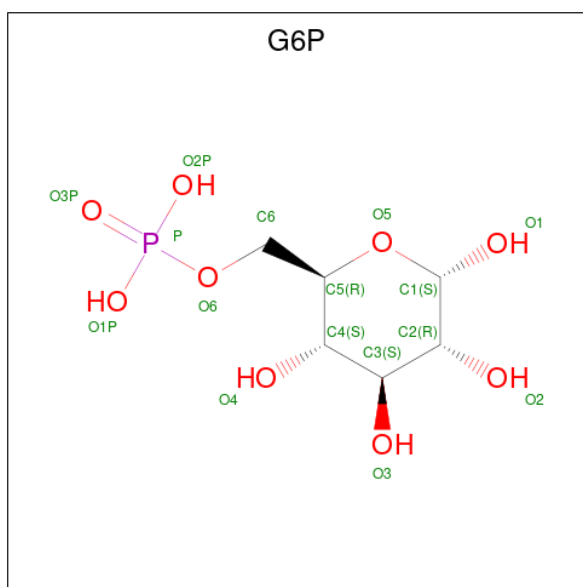
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	915	Total	C	N	O	S	0	0	0
			7298	4637	1267	1364	30			
1	B	911	Total	C	N	O	S	0	0	0
			7265	4617	1261	1357	30			

- Molecule 2 is GLYCINE (CCD ID: GLY) (formula: $C_2H_5NO_2$).



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

- Molecule 3 is 6-O-phosphono-alpha-D-glucopyranose (CCD ID: G6P) (formula: $C_6H_{13}O_9P$).



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

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



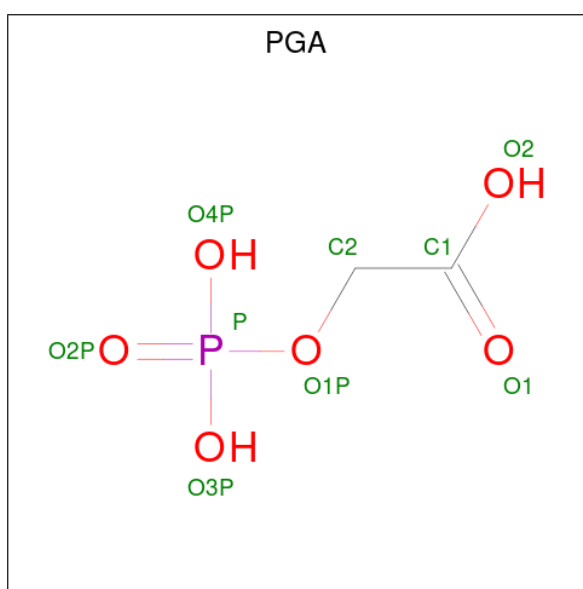
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		

Continued on next page...

Continued from previous page...

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

- Molecule 5 is 2-PHOSPHOGLYCOLIC ACID (CCD ID: PGA) (formula: $C_2H_5O_6P$).



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

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



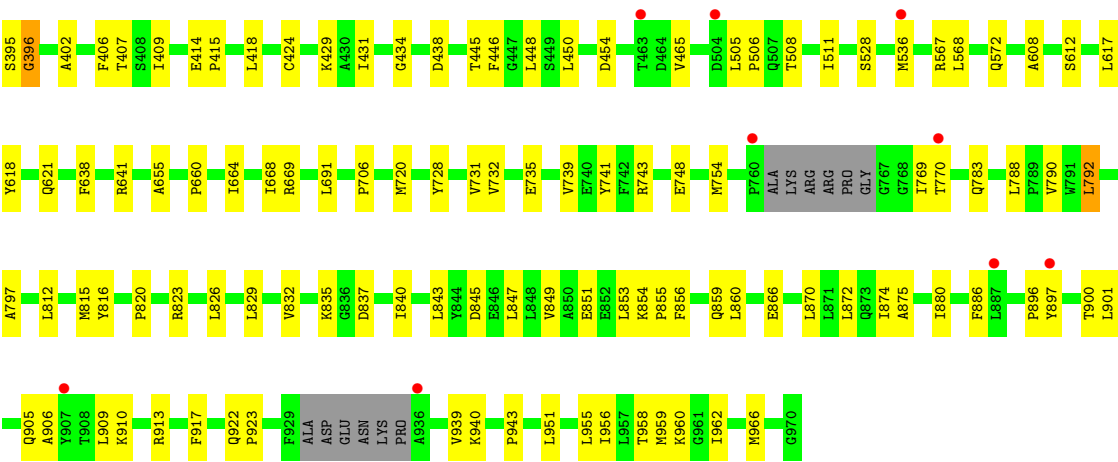
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			6	3	3		
6	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	13	Total	O	0	0
			13	13		
7	B	8	Total	O	0	0
			8	8		

- Molecule 1: Phosphoenolpyruvate carboxylase





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	152.09Å 170.68Å 244.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.62 – 2.99 29.62 – 2.99	Depositor EDS
% Data completeness (in resolution range)	97.1 (29.62-2.99) 97.0 (29.62-2.99)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 3.00Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.208 , 0.243 0.211 , 0.244	Depositor DCC
R_{free} test set	3168 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	61.8	Xtriage
Anisotropy	0.100	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 22.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	14680	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.31% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, GOL, PGA, G6P

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	0.15	0/7453	0.34	0/10096
1	B	0.13	0/7418	0.33	0/10046
All	All	0.14	0/14871	0.33	0/20142

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	7298	0	7294	87	0
1	B	7265	0	7261	124	0
2	A	5	0	2	0	0
2	B	5	0	2	0	0
3	A	16	0	11	0	0
3	B	16	0	11	1	0
4	A	16	0	24	0	0
4	B	8	0	12	0	0
5	A	9	0	2	0	0
5	B	9	0	2	0	0
6	B	12	0	16	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	13	0	0	0	0
7	B	8	0	0	0	0
All	All	14680	0	14637	206	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (206) 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:887:LEU:H	1:A:887:LEU:HD23	1.26	0.99
1:B:837:ASP:HB3	1:B:840:ILE:HD13	1.47	0.96
1:A:880:ILE:HG22	1:A:881:LEU:HG	1.57	0.87
1:A:887:LEU:HD23	1:A:887:LEU:N	1.93	0.84
1:B:359:ILE:HG22	1:B:360:GLU:HG2	1.67	0.77
1:A:403:GLU:OE1	1:A:403:GLU:N	2.19	0.75
1:B:434:GLY:HA3	6:B:1005:GOL:H31	1.71	0.71
1:B:728:TYR:HA	1:B:792:LEU:HD12	1.73	0.69
1:B:840:ILE:HD12	1:B:840:ILE:H	1.57	0.68
1:B:465:VAL:HG22	1:B:511:ILE:HG23	1.76	0.68
1:B:706:PRO:HB3	1:B:820:PRO:HG2	1.77	0.67
1:A:264:ASP:OD2	1:A:441:ARG:NH2	2.28	0.66
1:A:870:LEU:O	1:A:873:GLN:N	2.28	0.66
1:A:887:LEU:H	1:A:887:LEU:CD2	2.05	0.66
1:B:720:MET:HB3	1:B:797:ALA:HB1	1.78	0.65
1:A:264:ASP:OD1	1:A:276:LEU:N	2.22	0.64
1:B:905:GLN:HB2	1:B:958:THR:HG21	1.78	0.64
1:A:261:ARG:NH2	1:A:433:ASP:O	2.30	0.64
1:A:715:LYS:O	1:A:719:GLU:HG3	1.97	0.64
1:B:840:ILE:HD12	1:B:840:ILE:N	2.11	0.64
1:A:754:MET:HE2	1:A:756:ILE:HD11	1.79	0.63
1:A:878:LYS:HG2	1:A:882:GLU:CD	2.23	0.63
1:B:837:ASP:CB	1:B:840:ILE:HD13	2.26	0.63
1:A:184:ARG:HH22	1:B:360:GLU:CD	2.06	0.63
1:A:428:ASP:OD1	1:B:226:ARG:HD2	1.99	0.63
1:B:50:GLN:HG3	1:B:55:PRO:HA	1.79	0.62
1:A:395:SER:O	1:A:397:VAL:N	2.26	0.62
1:B:84:GLY:HA2	1:B:87:LEU:HD12	1.81	0.61
1:A:184:ARG:NH2	1:B:360:GLU:OE1	2.34	0.60
1:B:851:GLU:HA	1:B:854:LYS:HD2	1.84	0.60
1:B:741:TYR:HD1	1:B:849:VAL:HG21	1.66	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:754:MET:HE3	1:B:960:LYS:HG2	1.84	0.60
1:B:74:LYS:HG2	1:B:76:ASP:HB2	1.83	0.59
1:B:40:LEU:HD11	1:B:44:ARG:HH21	1.67	0.58
1:B:505:LEU:HD12	1:B:506:PRO:HD2	1.86	0.57
1:A:884:ASP:HB3	1:A:887:LEU:HD21	1.85	0.57
1:B:816:TYR:CE2	1:B:875:ALA:HA	2.39	0.57
1:B:896:PRO:O	1:B:900:THR:HG23	2.05	0.57
1:A:320:ILE:O	1:A:324:GLU:HG3	2.05	0.56
1:B:229:GLU:CD	1:B:229:GLU:H	2.14	0.56
1:B:872:LEU:HD21	1:B:880:ILE:HD13	1.87	0.56
1:B:568:LEU:HD22	1:B:608:ALA:HB2	1.88	0.56
1:B:409:ILE:HD13	1:B:448:LEU:HG	1.88	0.56
1:A:214:ASP:OD2	1:A:218:GLN:NE2	2.37	0.56
1:B:248:ILE:HA	1:B:252:VAL:HB	1.88	0.55
1:A:473:LEU:HB2	1:A:475:ILE:HD12	1.87	0.55
1:B:905:GLN:CB	1:B:958:THR:HG21	2.37	0.54
1:B:214:ASP:OD2	1:B:218:GLN:NE2	2.37	0.54
1:A:743:ARG:NH1	1:A:748:GLU:OE1	2.38	0.54
1:A:884:ASP:CB	1:A:887:LEU:HD21	2.38	0.54
1:A:256:VAL:HG12	1:A:691:LEU:HD11	1.89	0.54
1:B:956:ILE:HA	1:B:959:MET:HE3	1.90	0.54
1:B:335:CYS:HB2	1:B:339:LEU:HD23	1.90	0.54
1:B:951:LEU:HD12	1:B:951:LEU:H	1.73	0.53
1:A:110:ASN:O	1:A:114:GLU:HG3	2.08	0.53
1:B:184:ARG:HD2	1:B:243:TYR:HB2	1.89	0.53
1:A:502:PRO:HG2	1:A:505:LEU:HB2	1.91	0.52
1:B:446:PHE:HB3	1:B:450:LEU:HD23	1.92	0.52
1:A:284:ARG:HH12	1:A:452:LYS:HE3	1.74	0.52
1:A:716:LEU:HG	1:A:720:MET:HE2	1.91	0.52
1:A:41:LEU:HD12	1:A:112:ALA:HB2	1.91	0.52
1:B:256:VAL:HG11	1:B:445:THR:HG21	1.92	0.52
1:B:288:TRP:CD1	1:B:454:ASP:HB2	2.45	0.52
1:B:40:LEU:O	1:B:44:ARG:HD2	2.11	0.51
1:A:176:ALA:HB2	1:A:289:MET:HG2	1.91	0.51
1:A:337:ASP:O	1:A:341:VAL:HG12	2.10	0.51
1:B:845:ASP:O	1:B:849:VAL:HG12	2.10	0.51
1:B:395:SER:OG	1:B:396:GLY:N	2.35	0.51
1:A:324:GLU:OE1	1:A:383:TYR:OH	2.23	0.50
1:A:177:HIS:ND1	1:A:673:GLN:HG2	2.26	0.50
1:B:843:LEU:O	1:B:847:LEU:HD12	2.10	0.50
1:A:156:LYS:HE3	1:A:703:MET:HB3	1.92	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:THR:HG22	1:A:238:GLN:H	1.76	0.50
1:B:741:TYR:CD1	1:B:849:VAL:HG21	2.47	0.50
1:A:928:GLU:OE1	1:B:334:ARG:NH1	2.44	0.50
1:A:38:ASP:CG	1:A:893:LEU:HD21	2.36	0.50
1:A:465:VAL:HG22	1:A:511:ILE:HG23	1.94	0.49
1:A:942:ASN:OD1	1:A:945:SER:HB2	2.12	0.49
1:B:65:TYR:HD2	1:B:897:TYR:CE1	2.30	0.49
1:B:60:PHE:HE1	1:B:86:LYS:HG2	1.77	0.49
1:B:171:ASP:HB3	1:B:669:ARG:HG2	1.94	0.49
1:B:731:VAL:HG21	1:B:860:LEU:HD21	1.93	0.49
1:B:909:LEU:O	1:B:913:ARG:HG3	2.13	0.48
1:A:144:GLU:OE2	1:A:269:ASN:ND2	2.34	0.48
1:A:312:ARG:NH1	1:A:528:SER:HB3	2.28	0.48
1:B:840:ILE:H	1:B:840:ILE:CD1	2.24	0.48
1:B:568:LEU:O	1:B:572:GLN:HG3	2.12	0.48
1:A:746:THR:HB	1:A:780:SER:HB3	1.95	0.48
1:A:381:LYS:NZ	1:A:411:GLU:OE2	2.37	0.48
1:A:945:SER:OG	1:A:952:GLU:OE1	2.31	0.48
1:A:454:ASP:CG	1:A:531:PRO:HD2	2.39	0.48
1:B:293:ARG:NH2	1:B:304:THR:OG1	2.42	0.48
1:A:248:ILE:HA	1:A:252:VAL:HB	1.96	0.47
1:A:608:ALA:HA	1:A:725:THR:HG23	1.96	0.47
1:A:785:ARG:NH2	1:A:899:THR:OG1	2.46	0.47
1:B:44:ARG:HH12	1:B:214:ASP:CG	2.23	0.47
1:A:798:PHE:HZ	1:A:826:LEU:HD21	1.79	0.47
1:A:888:LYS:O	1:A:892:VAL:HG13	2.14	0.47
1:B:40:LEU:HD11	1:B:44:ARG:NH2	2.28	0.47
1:B:835:LYS:HB3	1:B:966:MET:HE1	1.96	0.47
1:A:162:PHE:CE2	1:A:166:LYS:HD2	2.49	0.47
1:A:409:ILE:HD12	1:A:409:ILE:H	1.80	0.47
1:B:402:ALA:HB1	1:B:407:THR:HG21	1.97	0.47
1:B:783:GLN:HE22	1:B:959:MET:HA	1.78	0.47
1:B:816:TYR:CZ	1:B:823:ARG:NH1	2.81	0.47
1:B:866:GLU:O	1:B:870:LEU:HD12	2.15	0.47
1:B:939:VAL:HG12	1:B:940:LYS:HD2	1.97	0.47
1:A:232:ARG:NH1	1:A:942:ASN:HB2	2.30	0.47
1:A:424:CYS:SG	1:A:429:LYS:HG2	2.55	0.47
1:B:379:ARG:HG2	1:B:383:TYR:CE1	2.50	0.46
1:B:812:LEU:HA	1:B:815:MET:HE2	1.96	0.46
1:A:111:LEU:O	1:A:115:VAL:HG23	2.15	0.46
1:A:142:ASP:OD1	1:A:143:ILE:N	2.48	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:940:LYS:HE3	1:B:940:LYS:HA	1.97	0.46
1:A:41:LEU:HD21	1:A:197:LEU:HD13	1.98	0.46
1:B:847:LEU:HD13	1:B:906:ALA:HB1	1.97	0.46
1:A:928:GLU:CD	1:B:334:ARG:HH12	2.24	0.46
1:B:283:ILE:HG21	1:B:691:LEU:HD22	1.98	0.46
1:A:173:VAL:HG22	1:A:286:SER:HB2	1.96	0.46
1:A:813:LYS:O	1:A:817:ASN:ND2	2.47	0.46
1:B:910:LYS:HG2	1:B:917:PHE:HD1	1.80	0.46
1:A:178:PRO:HB2	1:A:751:TYR:OH	2.15	0.46
1:B:57:LEU:O	1:B:61:VAL:HG23	2.15	0.46
1:A:709:PRO:HB3	1:A:819:TRP:CZ2	2.51	0.46
1:B:769:ILE:HG13	1:B:770:THR:N	2.31	0.45
1:B:910:LYS:HG2	1:B:917:PHE:CD1	2.52	0.45
1:B:739:VAL:HG22	1:B:769:ILE:HG21	1.99	0.45
1:A:598:MET:HA	1:A:638:PHE:HB3	1.98	0.45
1:B:332:MET:SD	1:B:332:MET:N	2.89	0.45
1:B:788:LEU:HD11	1:B:792:LEU:HD23	1.98	0.45
1:B:37:TYR:CE1	1:B:201:ASN:ND2	2.84	0.45
1:B:743:ARG:NH1	1:B:748:GLU:OE1	2.50	0.45
1:B:238:GLN:O	1:B:242:ARG:HG2	2.17	0.45
1:A:808:ASN:HA	1:A:811:VAL:HG12	1.99	0.45
1:B:342:ARG:HH22	1:B:414:GLU:CD	2.24	0.44
1:B:350:SER:O	1:B:364:GLN:NE2	2.50	0.44
1:B:956:ILE:O	1:B:960:LYS:HG3	2.18	0.44
1:B:837:ASP:CG	1:B:840:ILE:CD1	2.91	0.44
1:A:293:ARG:NH2	1:A:304:THR:OG1	2.48	0.44
1:B:843:LEU:HG	1:B:847:LEU:HD11	1.99	0.44
1:A:188:GLN:OE1	1:A:192:ARG:NE	2.51	0.44
1:A:501:LEU:HD12	1:A:502:PRO:HD2	2.00	0.43
1:B:200:LEU:HD11	1:B:210:LYS:HG2	2.00	0.43
1:A:171:ASP:HB3	1:A:669:ARG:HG2	2.00	0.43
1:B:42:VAL:O	1:B:46:LEU:HD23	2.17	0.43
1:B:664:ILE:HG23	1:B:668:ILE:HB	1.99	0.43
1:A:720:MET:HB3	1:A:797:ALA:HB1	2.01	0.43
1:B:200:LEU:HA	1:B:200:LEU:HD12	1.82	0.43
1:B:102:ILE:HG21	1:B:901:LEU:HG	2.00	0.43
1:B:147:LEU:HG	1:B:161:VAL:HG11	2.00	0.43
1:A:205:ILE:HD11	1:A:210:LYS:HG3	2.00	0.43
1:B:261:ARG:NH2	1:B:438:ASP:OD1	2.52	0.43
1:A:278:TYR:OH	1:A:417:GLU:OE1	2.30	0.43
1:A:568:LEU:HD22	1:A:608:ALA:HB2	2.01	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:874:ILE:H	1:A:874:ILE:HD12	1.84	0.43
1:B:339:LEU:HD13	1:B:418:LEU:HG	2.01	0.43
1:A:244:GLY:HA3	1:A:289:MET:SD	2.59	0.43
1:B:641:ARG:CZ	1:B:655:ALA:HB1	2.48	0.43
1:A:622:GLU:CD	1:A:714:ARG:HH22	2.27	0.43
1:B:318:LEU:HD12	1:B:450:LEU:HD11	2.01	0.43
1:B:332:MET:HE2	1:B:431:ILE:HD13	2.01	0.43
1:B:40:LEU:HD12	1:B:40:LEU:HA	1.69	0.42
1:B:508:THR:OG1	1:B:511:ILE:HD12	2.19	0.42
1:B:617:LEU:O	1:B:621:GLN:HG3	2.19	0.42
1:A:362:TRP:CD1	1:A:362:TRP:H	2.36	0.42
1:B:363:LYS:HE3	1:B:363:LYS:HB2	1.85	0.42
1:B:790:VAL:HG11	1:B:832:VAL:HG21	2.01	0.42
1:B:114:GLU:OE2	1:B:194:ARG:HD3	2.19	0.42
1:B:52:LEU:HD22	1:B:226:ARG:NH2	2.34	0.42
1:B:117:ILE:HD11	1:B:886:PHE:HE2	1.85	0.42
1:B:381:LYS:HB3	1:B:406:PHE:CE2	2.55	0.42
1:B:855:PRO:O	1:B:859:GLN:HG3	2.20	0.42
1:A:125:LYS:HB3	1:A:125:LYS:HE3	1.87	0.42
1:B:536:MET:HE1	1:B:567:ARG:HE	1.85	0.42
1:B:922:GLN:HB2	1:B:923:PRO:HD2	2.02	0.42
1:A:192:ARG:NH1	1:A:220:GLU:OE2	2.53	0.41
1:A:859:GLN:O	1:A:863:LYS:HG3	2.19	0.41
1:B:955:LEU:HG	1:B:959:MET:HE2	2.02	0.41
1:B:121:ARG:O	1:B:125:LYS:HG3	2.20	0.41
1:B:618:TYR:CG	1:B:660:PRO:HG3	2.54	0.41
1:B:917:PHE:C	1:B:917:PHE:CD2	2.98	0.41
1:A:276:LEU:HD12	1:A:277:PRO:HD2	2.01	0.41
1:A:954:THR:O	1:A:958:THR:HG23	2.21	0.41
1:B:173:VAL:HG21	1:B:638:PHE:CZ	2.55	0.41
1:B:424:CYS:SG	1:B:429:LYS:HD3	2.60	0.41
1:B:728:TYR:CE1	1:B:732:VAL:HG21	2.56	0.41
1:B:41:LEU:HD23	1:B:41:LEU:HA	1.94	0.41
1:B:188:GLN:HE21	1:B:188:GLN:HB3	1.71	0.41
1:B:254:LYS:HA	1:B:254:LYS:HD2	1.64	0.41
1:B:312:ARG:NH1	1:B:528:SER:HB3	2.36	0.41
1:B:870:LEU:O	1:B:874:ILE:HG12	2.21	0.41
1:A:788:LEU:HD11	1:A:792:LEU:HD22	2.03	0.41
1:B:231:ARG:NH1	3:B:1002:G6P:O2P	2.53	0.41
1:B:414:GLU:HB3	1:B:415:PRO:HD3	2.02	0.41
1:B:853:LEU:O	1:B:856:PHE:HB3	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:905:GLN:HB2	1:B:958:THR:CG2	2.50	0.41
1:A:170:VAL:HG11	1:A:694:PHE:HB3	2.03	0.41
1:B:951:LEU:HD12	1:B:951:LEU:N	2.36	0.41
1:B:958:THR:O	1:B:962:ILE:HD12	2.21	0.40
1:A:184:ARG:HD3	1:A:243:TYR:HD1	1.86	0.40
1:A:353:LYS:HA	1:A:353:LYS:HD2	1.84	0.40
1:B:79:LYS:HZ2	1:B:79:LYS:HG2	1.77	0.40
1:A:371:TYR:HA	1:A:374:ILE:HG22	2.03	0.40
1:B:826:LEU:HA	1:B:829:LEU:HD12	2.04	0.40
1:A:330:LEU:HD22	1:A:332:MET:HE1	2.04	0.40
1:A:878:LYS:HG3	1:A:879:ASP:N	2.37	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	909/970 (94%)	875 (96%)	32 (4%)	2 (0%)	43	73
1	B	903/970 (93%)	867 (96%)	31 (3%)	5 (1%)	21	53
All	All	1812/1940 (93%)	1742 (96%)	63 (4%)	7 (0%)	30	62

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	396	GLY
1	B	74	LYS
1	A	870	LEU
1	B	792	LEU
1	B	943	PRO
1	B	735	GLU
1	B	396	GLY

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	791/832 (95%)	791 (100%)	0	100	100
1	B	787/832 (95%)	785 (100%)	2 (0%)	86	91
All	All	1578/1664 (95%)	1576 (100%)	2 (0%)	88	94

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	206	THR
1	B	612	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	107	ASN
1	A	190	ASN
1	A	322	GLN
1	A	377	HIS
1	A	659	GLN
1	A	665	ASN
1	A	869	GLN
1	B	188	GLN
1	B	190	ASN
1	B	201	ASN
1	B	572	GLN
1	B	689	GLN
1	B	787	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

14 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	EDO	A	1004	-	3,3,3	0.38	0	2,2,2	0.39	0
4	EDO	A	1005	-	3,3,3	0.39	0	2,2,2	0.43	0
5	PGA	B	1007	-	8,8,8	1.05	0	10,11,11	0.99	0
3	G6P	A	1002	-	16,16,16	0.53	0	23,24,24	0.67	0
2	GLY	A	1001	-	4,4,4	1.11	1 (25%)	3,4,4	1.68	2 (66%)
4	EDO	B	1003	-	3,3,3	0.42	0	2,2,2	0.34	0
6	GOL	B	1005	-	5,5,5	0.96	0	5,5,5	1.00	0
5	PGA	A	1007	-	8,8,8	1.12	1 (12%)	10,11,11	0.98	0
4	EDO	B	1004	-	3,3,3	0.39	0	2,2,2	0.40	0
6	GOL	B	1006	-	5,5,5	0.93	0	5,5,5	1.07	0
2	GLY	B	1001	-	4,4,4	1.13	1 (25%)	3,4,4	1.69	2 (66%)
4	EDO	A	1006	-	3,3,3	0.33	0	2,2,2	0.58	0
3	G6P	B	1002	-	16,16,16	0.52	0	23,24,24	0.71	0
4	EDO	A	1003	-	3,3,3	0.43	0	2,2,2	0.34	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	EDO	A	1004	-	-	0/1/1/1	-
4	EDO	A	1005	-	-	0/1/1/1	-
5	PGA	B	1007	-	-	2/6/6/6	-
3	G6P	A	1002	-	-	0/6/26/26	0/1/1/1
2	GLY	A	1001	-	-	0/2/2/2	-
4	EDO	B	1003	-	-	0/1/1/1	-
6	GOL	B	1005	-	-	0/4/4/4	-
5	PGA	A	1007	-	-	5/6/6/6	-
4	EDO	B	1004	-	-	0/1/1/1	-
6	GOL	B	1006	-	-	0/4/4/4	-
2	GLY	B	1001	-	-	1/2/2/2	-
4	EDO	A	1006	-	-	1/1/1/1	-
3	G6P	B	1002	-	-	0/6/26/26	0/1/1/1
4	EDO	A	1003	-	-	0/1/1/1	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1001	GLY	OXT-C	-2.15	1.23	1.30
2	A	1001	GLY	OXT-C	-2.11	1.23	1.30
5	A	1007	PGA	O1P-C2	-2.09	1.41	1.43

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	GLY	OXT-C-O	-2.08	117.99	123.33
2	B	1001	GLY	OXT-C-CA	2.05	121.52	113.38
2	B	1001	GLY	OXT-C-O	-2.03	118.12	123.33
2	A	1001	GLY	OXT-C-CA	2.00	121.34	113.38

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1007	PGA	C2-O1P-P-O3P
5	B	1007	PGA	O1-C1-C2-O1P
5	B	1007	PGA	O2-C1-C2-O1P
5	A	1007	PGA	O1-C1-C2-O1P
5	A	1007	PGA	O2-C1-C2-O1P
4	A	1006	EDO	O1-C1-C2-O2
5	A	1007	PGA	C2-O1P-P-O2P
5	A	1007	PGA	C2-O1P-P-O4P

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
2	B	1001	GLY	OXT-C-CA-N

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	1005	GOL	1	0
3	B	1002	G6P	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	915/970 (94%)	-0.11	7 (0%) 82 67	34, 47, 73, 134	0
1	B	911/970 (93%)	0.05	13 (1%) 73 55	32, 55, 86, 109	0
All	All	1826/1940 (94%)	-0.03	20 (1%) 78 61	32, 50, 82, 134	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	35	ILE	3.8
1	A	35	ILE	3.4
1	B	887	LEU	3.3
1	B	504	ASP	3.3
1	A	685	HIS	3.1
1	B	936	ALA	2.8
1	B	204	ASP	2.8
1	B	463	THR	2.6
1	A	474	GLY	2.6
1	B	907	TYR	2.5
1	A	887	LEU	2.4
1	B	536	MET	2.3
1	B	760	PRO	2.3
1	A	126	LEU	2.3
1	B	897	TYR	2.3
1	A	934	LYS	2.2
1	B	40	LEU	2.2
1	A	886	PHE	2.1
1	B	770	THR	2.0
1	B	264	ASP	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 oligosaccharides 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, 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	B-factors($\text{\AA}^2$)	Q<0.9
4	EDO	A	1006	4/4	0.69	0.23	43,43,43,43	0
5	PGA	B	1007	9/9	0.75	0.22	41,62,78,98	9
5	PGA	A	1007	9/9	0.78	0.23	38,50,70,80	9
4	EDO	B	1004	4/4	0.83	0.29	44,47,52,54	0
4	EDO	A	1005	4/4	0.86	0.17	50,50,50,50	0
2	GLY	B	1001	5/5	0.88	0.22	60,65,79,105	0
4	EDO	A	1003	4/4	0.89	0.18	31,39,46,50	0
2	GLY	A	1001	5/5	0.89	0.15	36,47,51,52	0
6	GOL	B	1005	6/6	0.90	0.12	37,44,46,51	0
4	EDO	A	1004	4/4	0.91	0.18	40,40,41,44	0
6	GOL	B	1006	6/6	0.91	0.11	40,46,49,50	0
4	EDO	B	1003	4/4	0.92	0.18	41,43,44,47	0
3	G6P	B	1002	16/16	0.93	0.07	50,57,67,69	0
3	G6P	A	1002	16/16	0.94	0.07	37,47,53,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.