



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 04:38 PM BST

PDB ID : 2CUN
Title : Crystal structure of Phosphoglycerate Kinase from *Pyrococcus horikoshii* OT3
Authors : Mizutani, H.; Kunishima, N.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2005-05-27
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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

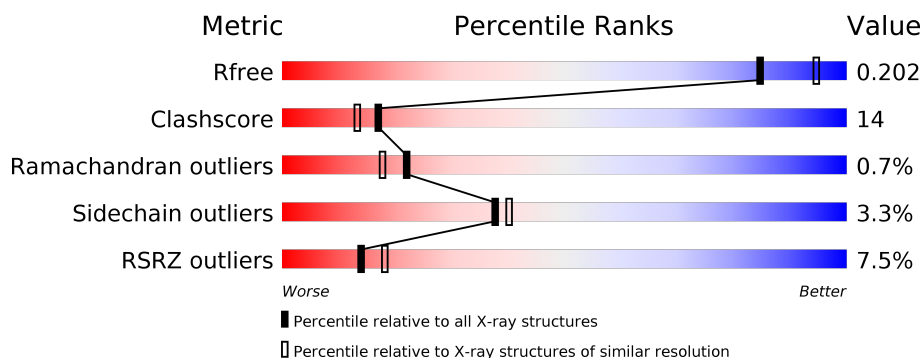
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (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 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	410	<div> <div>5%</div> <div> <div></div> <div>78%</div> <div>20%</div> <div>..</div> </div> </div>
1	B	410	<div> <div>10%</div> <div> <div></div> <div>64%</div> <div>33%</div> <div>..</div> </div> </div>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6890 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.

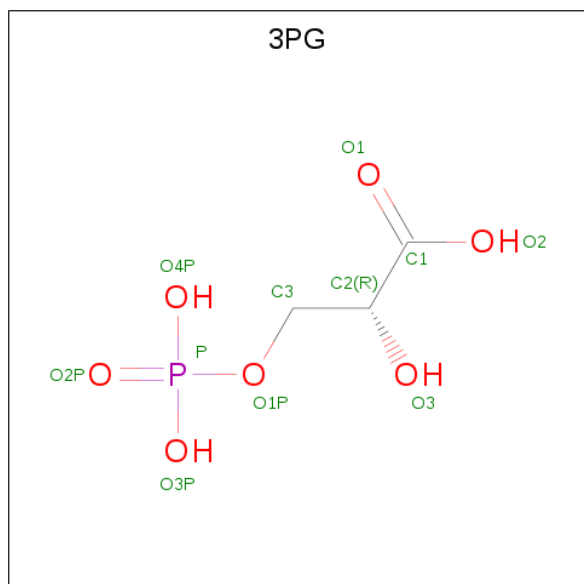
- Molecule 1 is a protein called Phosphoglycerate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	405	Total	C	N	O	S	0	0	0
			3229	2078	544	595	12			
1	B	405	Total	C	N	O	S	0	1	0
			3230	2078	544	596	12			

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 3 is 3-PHOSPHOGLYCERIC ACID (three-letter code: 3PG) (formula: C₃H₇O₇P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			11	3	7	1		
3	B	1	Total	C	O	P	0	0
			11	3	7	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 5 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			8	6	2		
5	A	1	Total	C	O	0	0
			8	6	2		
5	B	1	Total	C	O	0	0
			8	6	2		

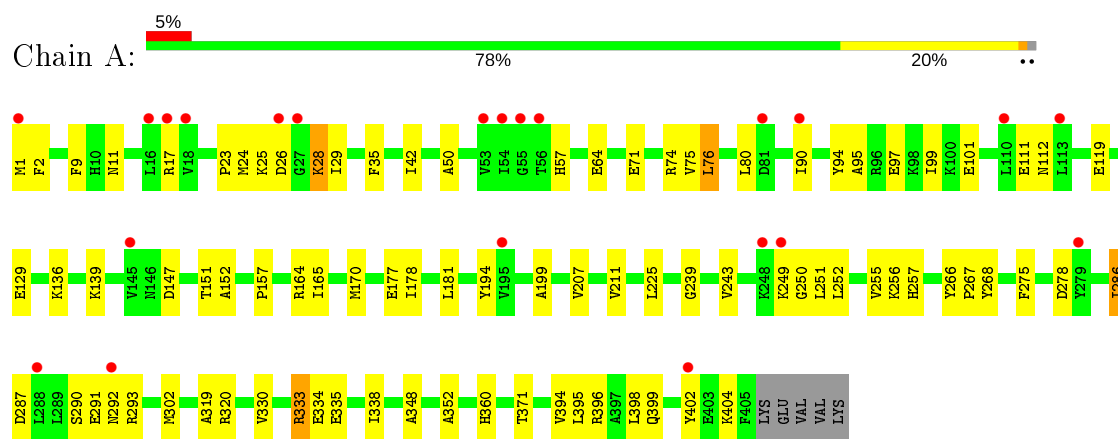
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	218	Total	O	0	0
			218	218		
6	B	153	Total	O	0	0
			153	153		

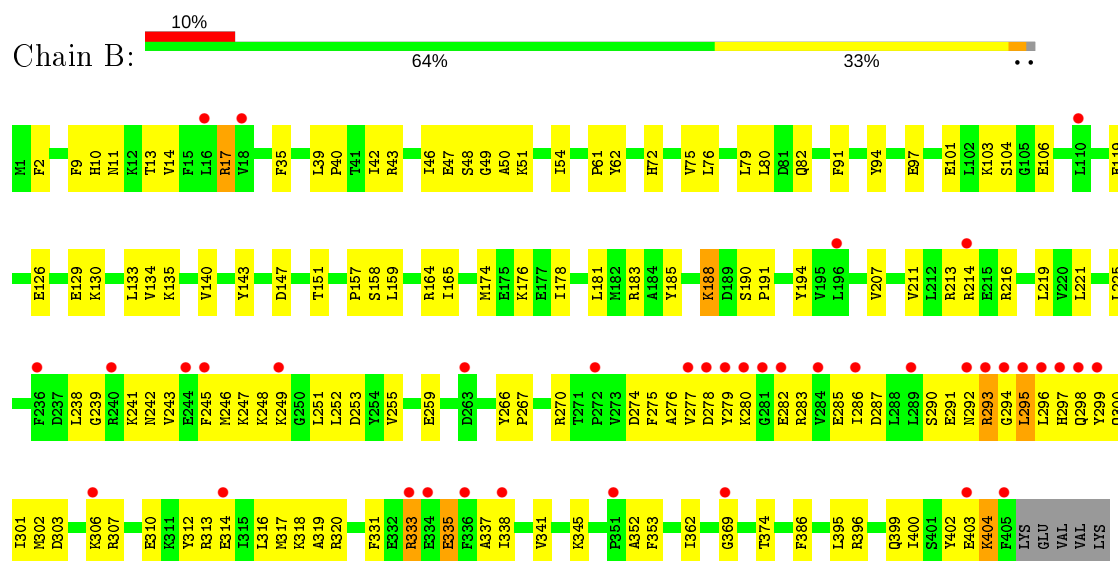
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Phosphoglycerate kinase



• Molecule 1: Phosphoglycerate kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.85Å 123.33Å 82.94Å 90.00° 95.42° 90.00°	Depositor
Resolution (Å)	30.00 – 2.10 29.48 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.8 (30.00-2.10) 99.9 (29.48-2.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.78 (at 2.10Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.210 , 0.244 0.207 , 0.202	Depositor DCC
R_{free} test set	2945 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	35.5	Xtriage
Anisotropy	0.493	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 53.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6890	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.97% 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: GOL, MPD, CL, 3PG

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	0.34	0/3289	0.60	0/4423
1	B	0.32	0/3295	0.57	0/4431
All	All	0.33	0/6584	0.58	0/8854

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	3229	0	3292	73	0
1	B	3230	0	3293	112	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	11	0	4	0	0
3	B	11	0	4	0	0
4	A	6	0	8	1	0
4	B	6	0	8	0	0
5	A	16	0	28	1	0
5	B	8	0	14	1	0
6	A	218	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	153	0	0	2	0
All	All	6890	0	6651	184	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 14.

All (184) 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:188:LYS:H	1:B:188:LYS:HD3	1.34	0.92
1:A:338:ILE:HD12	1:A:338:ILE:H	1.33	0.92
1:B:333:ARG:HH11	1:B:333:ARG:HB3	1.42	0.85
1:A:35:PHE:HD1	1:A:76:LEU:HD13	1.43	0.83
1:A:178:ILE:HD12	1:A:395:LEU:HD21	1.61	0.82
1:B:119:GLU:HA	1:B:157:PRO:HG2	1.64	0.79
1:A:136:LYS:HA	1:A:139:LYS:HE3	1.65	0.77
1:B:247:LYS:HG2	1:B:252:LEU:HD12	1.65	0.77
1:B:318:LYS:HA	1:B:318:LYS:HE2	1.67	0.77
1:B:188:LYS:HD3	1:B:188:LYS:N	2.00	0.76
1:A:396:ARG:HA	1:A:399:GLN:HE21	1.52	0.74
1:B:296:LEU:HD12	1:B:296:LEU:H	1.50	0.74
1:B:319:ALA:O	1:B:352:ALA:HB2	1.88	0.73
1:A:249:LYS:HB3	1:A:251:LEU:HD13	1.71	0.73
1:B:17:ARG:HD2	1:B:158:SER:OG	1.91	0.70
1:B:245:PHE:O	1:B:248:LYS:HB2	1.91	0.70
1:B:287:ASP:OD2	1:B:307:ARG:HD3	1.90	0.70
1:A:333:ARG:HH11	1:A:333:ARG:HB3	1.56	0.70
1:A:24:MET:HE3	1:A:29:ILE:HG13	1.73	0.69
1:B:225:LEU:HD13	1:B:251:LEU:HD21	1.73	0.69
1:A:338:ILE:H	1:A:338:ILE:CD1	2.05	0.69
1:B:306:LYS:O	1:B:310:GLU:HG3	1.93	0.69
1:B:313:ARG:O	1:B:317:MET:HG2	1.93	0.68
1:B:188:LYS:CD	1:B:188:LYS:H	2.05	0.68
1:A:97:GLU:O	1:A:101:GLU:HG3	1.94	0.67
1:A:129:GLU:OE1	1:A:164:ARG:HD2	1.93	0.67
1:A:338:ILE:HD12	1:A:338:ILE:N	2.10	0.66
1:A:35:PHE:CD1	1:A:76:LEU:HD13	2.29	0.66
1:B:255:VAL:O	1:B:259:GLU:HG3	1.96	0.66
1:A:250:GLY:HA2	6:A:666:HOH:O	1.97	0.65
1:B:14:VAL:HG22	1:B:143:TYR:HB2	1.78	0.64
1:B:306:LYS:HE3	1:B:310:GLU:OE1	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:293:ARG:HB2	1:B:293:ARG:HH11	1.63	0.63
1:B:287:ASP:HB3	1:B:290:SER:HB2	1.80	0.63
1:A:119:GLU:HA	1:A:157:PRO:HG2	1.82	0.61
1:B:241:LYS:HB3	1:B:299:TYR:O	2.00	0.61
1:B:54:ILE:HD13	1:B:76:LEU:HD21	1.81	0.61
1:A:256:LYS:HG3	1:A:257:HIS:H	1.65	0.60
1:B:43:ARG:HH11	1:B:43:ARG:HG3	1.65	0.60
1:A:178:ILE:CD1	1:A:395:LEU:HD21	2.30	0.59
1:B:97:GLU:O	1:B:101:GLU:HG3	2.01	0.59
1:B:296:LEU:CD1	1:B:296:LEU:H	2.15	0.59
1:B:337:ALA:O	1:B:341:VAL:HG23	2.03	0.59
1:A:29:ILE:HG13	1:A:71:GLU:OE1	2.03	0.59
1:A:23:PRO:HG3	1:A:360:HIS:CG	2.38	0.59
1:B:320:ARG:HG3	1:B:320:ARG:HH11	1.68	0.59
1:A:9:PHE:HB3	1:A:50:ALA:HB2	1.84	0.58
1:B:296:LEU:HD12	1:B:296:LEU:N	2.17	0.58
1:B:280:LYS:O	1:B:280:LYS:HG2	2.03	0.58
1:B:252:LEU:HA	1:B:255:VAL:HG23	1.86	0.58
1:A:267:PRO:HG2	1:A:268:TYR:CD1	2.39	0.57
1:A:286:ILE:HD13	1:A:286:ILE:H	1.69	0.57
1:B:239:GLY:O	1:B:243:VAL:HG23	2.05	0.57
1:A:275:PHE:HB2	1:A:286:ILE:HD11	1.86	0.57
1:B:310:GLU:O	1:B:314:GLU:HG3	2.05	0.57
1:B:295:LEU:HD22	1:B:299:TYR:CE2	2.40	0.56
1:B:42:ILE:O	1:B:46:ILE:HG13	2.05	0.56
1:B:130:LYS:HA	1:B:135:LYS:HE3	1.86	0.56
1:B:75:VAL:O	1:B:79:LEU:HD13	2.06	0.56
1:B:129:GLU:OE1	1:B:164:ARG:HD2	2.06	0.56
1:B:35:PHE:HD1	1:B:76:LEU:HD13	1.71	0.56
1:B:239:GLY:HA3	1:B:296:LEU:HB3	1.87	0.56
1:B:2:PHE:CE1	1:B:402:TYR:HA	2.41	0.55
1:A:252:LEU:HA	1:A:255:VAL:HG23	1.88	0.55
1:A:74:ARG:HG3	1:A:75:VAL:N	2.22	0.55
1:A:335:GLU:CD	1:A:335:GLU:H	2.10	0.55
1:A:287:ASP:HB3	1:A:290:SER:CB	2.37	0.54
1:A:239:GLY:O	1:A:243:VAL:HG23	2.08	0.54
1:B:341:VAL:HG12	1:B:345:LYS:HE3	1.89	0.54
1:A:333:ARG:HB3	1:A:333:ARG:NH1	2.23	0.54
1:B:277:VAL:HG22	1:B:278:ASP:N	2.23	0.54
1:B:362:ILE:HG13	5:B:507:MPD:HM1	1.90	0.54
1:A:287:ASP:HB3	1:A:290:SER:HB2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:242:ASN:O	1:B:246:MET:HG2	2.08	0.53
1:B:43:ARG:O	1:B:47:GLU:HG2	2.09	0.53
1:B:17:ARG:HD3	1:B:159:LEU:HD23	1.90	0.53
1:A:24:MET:HE3	1:A:29:ILE:CG1	2.37	0.53
1:B:213:ARG:HG3	6:B:635:HOH:O	2.09	0.53
1:B:395:LEU:O	1:B:399:GLN:HG3	2.09	0.53
1:B:293:ARG:CB	1:B:293:ARG:HH11	2.21	0.52
1:A:28:LYS:HB3	1:A:28:LYS:NZ	2.24	0.52
1:B:396:ARG:O	1:B:400:ILE:HG13	2.09	0.52
1:B:292:ASN:O	1:B:293:ARG:HG3	2.10	0.52
1:B:10:HIS:HA	1:B:48:SER:O	2.09	0.52
1:A:1:MET:HE1	1:A:170:MET:SD	2.50	0.52
1:A:256:LYS:HG3	1:A:257:HIS:N	2.25	0.52
1:A:178:ILE:HD11	1:A:395:LEU:HD11	1.92	0.51
1:B:335:GLU:H	1:B:335:GLU:CD	2.12	0.51
1:B:9:PHE:HB3	1:B:50:ALA:HB2	1.93	0.51
1:B:249:LYS:HB2	1:B:251:LEU:HD13	1.91	0.51
1:B:295:LEU:HB3	1:B:299:TYR:HD2	1.76	0.51
1:A:293:ARG:HG3	1:A:293:ARG:HH11	1.76	0.51
1:B:207:VAL:O	1:B:211:VAL:HG22	2.11	0.51
1:B:276:ALA:HA	1:B:285:GLU:HA	1.92	0.51
1:B:103:LYS:HG2	1:B:106:GLU:OE1	2.11	0.51
1:A:94:TYR:HB2	1:B:94:TYR:HB2	1.93	0.51
1:A:278:ASP:HB2	1:A:302:MET:HG3	1.92	0.50
1:B:238:LEU:O	1:B:242:ASN:HB2	2.11	0.50
1:B:183:ARG:HD3	1:B:353:PHE:CE2	2.47	0.50
1:A:348:ALA:HB1	1:A:371:THR:O	2.12	0.50
1:A:225:LEU:HD13	1:A:251:LEU:HD21	1.93	0.50
1:A:319:ALA:O	1:A:352:ALA:HB2	2.12	0.50
1:A:64:GLU:HG2	6:A:707:HOH:O	2.11	0.50
1:A:119:GLU:HA	1:A:157:PRO:CG	2.43	0.49
1:B:275:PHE:HD2	1:B:301:ILE:HG21	1.77	0.49
1:A:2:PHE:CE1	1:A:402:TYR:HA	2.48	0.49
1:B:221:LEU:HD23	1:B:270:ARG:HB2	1.95	0.48
1:B:333:ARG:HH11	1:B:333:ARG:CB	2.19	0.48
1:B:126:GLU:CD	1:B:126:GLU:H	2.16	0.48
1:B:295:LEU:HB3	1:B:299:TYR:CD2	2.48	0.48
1:A:177:GLU:O	1:A:181:LEU:HD13	2.13	0.48
1:A:207:VAL:O	1:A:211:VAL:HG22	2.14	0.48
1:A:194:TYR:CE2	1:A:211:VAL:HG11	2.48	0.48
1:B:190:SER:HB2	1:B:191:PRO:HA	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:225:LEU:HD13	1:B:251:LEU:CD2	2.42	0.48
1:A:266:TYR:N	1:A:267:PRO:HD2	2.29	0.47
1:B:294:GLY:O	1:B:296:LEU:N	2.47	0.47
1:B:247:LYS:CG	1:B:252:LEU:HD12	2.40	0.47
1:B:292:ASN:C	1:B:293:ARG:HG3	2.34	0.47
1:A:396:ARG:HA	1:A:399:GLN:NE2	2.24	0.47
1:A:249:LYS:CB	1:A:251:LEU:HD13	2.43	0.47
1:B:183:ARG:HD3	1:B:353:PHE:CZ	2.49	0.47
1:B:190:SER:O	1:B:320:ARG:HG3	2.15	0.47
1:B:13:THR:HG21	1:B:140:VAL:O	2.14	0.47
1:B:279:TYR:O	1:B:280:LYS:HB3	2.15	0.46
1:A:199:ALA:CB	1:A:225:LEU:HD12	2.45	0.46
1:A:330:VAL:CG1	1:A:333:ARG:HG3	2.45	0.46
1:B:266:TYR:N	1:B:267:PRO:HD2	2.30	0.46
1:B:13:THR:HA	1:B:51:LYS:HB2	1.97	0.46
1:B:103:LYS:HB2	1:B:106:GLU:HG3	1.96	0.46
1:B:181:LEU:HB3	1:B:386:PHE:CE2	2.51	0.46
1:A:275:PHE:O	1:A:286:ILE:HD13	2.16	0.46
1:B:312:TYR:O	1:B:316:LEU:HG	2.16	0.46
1:A:152:ALA:HA	1:A:394:VAL:CG2	2.46	0.46
1:B:277:VAL:HG23	1:B:300:GLN:O	2.15	0.45
1:B:49:GLY:O	1:B:104:SER:HB2	2.15	0.45
1:B:134:VAL:CG1	1:B:165:ILE:HD12	2.47	0.45
1:A:95:ALA:O	1:A:99:ILE:HG13	2.17	0.45
1:B:402:TYR:C	1:B:404:LYS:H	2.20	0.45
1:B:276:ALA:HB2	1:B:285:GLU:HG2	1.99	0.44
1:B:61:PRO:O	1:B:62:TYR:HB2	2.16	0.44
1:B:185:TYR:O	1:B:216:ARG:HD2	2.18	0.44
1:B:353:PHE:CD1	1:B:374:THR:HG21	2.52	0.44
1:B:287:ASP:HB3	1:B:290:SER:CB	2.46	0.44
1:A:57:HIS:HE2	4:A:501:GOL:H11	1.83	0.44
1:A:291:GLU:C	1:A:293:ARG:H	2.20	0.44
1:B:126:GLU:O	1:B:130:LYS:HD3	2.17	0.43
1:A:25:LYS:HD2	1:A:26:ASP:OD2	2.18	0.43
1:A:333:ARG:NH1	1:A:335:GLU:OE2	2.51	0.43
1:A:330:VAL:HG11	1:A:333:ARG:HG3	2.01	0.43
1:B:341:VAL:CG1	1:B:345:LYS:HE3	2.47	0.43
1:B:302:MET:O	1:B:303:ASP:HB2	2.19	0.43
1:A:293:ARG:NE	6:A:570:HOH:O	2.52	0.43
1:B:277:VAL:CG2	1:B:278:ASP:N	2.82	0.43
1:A:1:MET:CE	1:A:398:LEU:HD13	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:ILE:HG21	1:A:80:LEU:HD11	2.01	0.42
1:B:341:VAL:O	1:B:345:LYS:HG3	2.18	0.42
1:B:119:GLU:HA	1:B:157:PRO:CG	2.42	0.42
1:A:139:LYS:HE2	6:A:563:HOH:O	2.19	0.42
1:B:293:ARG:CB	1:B:293:ARG:NH1	2.83	0.42
1:B:286:ILE:HG21	1:B:293:ARG:HG2	2.02	0.42
1:B:353:PHE:HD1	1:B:374:THR:HG21	1.84	0.42
1:B:43:ARG:HG3	1:B:43:ARG:NH1	2.33	0.42
1:A:320:ARG:HA	1:A:320:ARG:HD3	1.86	0.42
1:A:136:LYS:HA	1:A:139:LYS:CE	2.44	0.42
1:A:24:MET:HE2	1:A:28:LYS:C	2.39	0.42
1:A:28:LYS:HB3	1:A:28:LYS:HZ3	1.85	0.41
1:B:283:ARG:NE	1:B:338:ILE:HD12	2.35	0.41
1:B:320:ARG:HG3	1:B:320:ARG:NH1	2.34	0.41
1:A:1:MET:HE3	1:A:398:LEU:HD13	2.02	0.41
1:B:274:ASP:HB2	1:B:286:ILE:O	2.21	0.41
1:A:90:ILE:HD12	5:A:509:MPD:CM	2.51	0.41
1:B:91:PHE:CD2	1:B:133:LEU:HA	2.56	0.41
1:B:219:LEU:HD11	1:B:270:ARG:NH1	2.36	0.41
1:B:72:HIS:O	1:B:76:LEU:HB2	2.21	0.41
1:B:194:TYR:CE2	1:B:211:VAL:HG11	2.55	0.41
1:B:39:LEU:HB2	1:B:40:PRO:HD3	2.03	0.41
1:B:176:LYS:HG3	6:B:589:HOH:O	2.20	0.41
1:A:111:GLU:HG3	1:A:112:ASN:N	2.36	0.40
1:B:174:MET:O	1:B:178:ILE:HG13	2.21	0.40
1:B:80:LEU:C	1:B:82:GLN:H	2.24	0.40
1:A:129:GLU:HB2	1:A:165:ILE:HD11	2.04	0.40
1:A:334:GLU:HG2	6:A:681:HOH:O	2.21	0.40
1:B:133:LEU:HD23	1:B:133:LEU:C	2.42	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	403/410 (98%)	392 (97%)	10 (2%)	1 (0%)	47	49
1	B	404/410 (98%)	375 (93%)	24 (6%)	5 (1%)	13	8
All	All	807/820 (98%)	767 (95%)	34 (4%)	6 (1%)	22	18

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	295	LEU
1	B	298	GLN
1	B	369	GLY
1	A	292	ASN
1	B	291	GLU
1	B	403	GLU

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	345/350 (99%)	336 (97%)	9 (3%)	46	50
1	B	346/350 (99%)	332 (96%)	14 (4%)	31	32
All	All	691/700 (99%)	668 (97%)	23 (3%)	38	40

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	17	ARG
1	A	28	LYS
1	A	76	LEU
1	A	147	ASP
1	A	151	THR
1	A	286	ILE
1	A	333	ARG
1	A	404	LYS
1	B	11	ASN

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Mol	Chain	Res	Type
1	B	17	ARG
1	B	147	ASP
1	B	151	THR
1	B	188	LYS
1	B	214	ARG
1	B	253	ASP
1	B	282	GLU
1	B	293	ARG
1	B	297	HIS
1	B	331	PHE
1	B	333	ARG
1	B	335	GLU
1	B	404	LYS

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	11	ASN
1	A	399	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 9 ligands modelled in this entry, 2 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
3	3PG	B	504	-	7,10,10	2.98	3 (42%)	10,14,14	1.06	1 (10%)
5	MPD	B	507	-	7,7,7	0.60	0	9,10,10	0.31	0
5	MPD	A	508	-	7,7,7	0.58	0	9,10,10	0.34	0
3	3PG	A	503	-	7,10,10	3.09	4 (57%)	10,14,14	1.14	1 (10%)
5	MPD	A	509	-	7,7,7	0.76	0	9,10,10	0.38	0
4	GOL	A	501	-	5,5,5	0.27	0	5,5,5	0.38	0
4	GOL	B	502	-	5,5,5	0.25	0	5,5,5	0.32	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
3	3PG	B	504	-	-	2/6/10/10	-
5	MPD	B	507	-	-	0/5/5/5	-
5	MPD	A	508	-	-	0/5/5/5	-
3	3PG	A	503	-	-	2/6/10/10	-
5	MPD	A	509	-	-	0/5/5/5	-
4	GOL	A	501	-	-	0/4/4/4	-
4	GOL	B	502	-	-	0/4/4/4	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	503	3PG	P-O2P	5.68	1.68	1.50
3	B	504	3PG	P-O2P	5.63	1.68	1.50
3	A	503	3PG	P-O4P	3.81	1.69	1.54
3	A	503	3PG	P-O3P	3.57	1.68	1.54
3	B	504	3PG	P-O4P	3.51	1.68	1.54
3	B	504	3PG	P-O3P	3.34	1.67	1.54
3	A	503	3PG	P-O1P	2.01	1.66	1.60

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	B	504	3PG	O1P-C3-C2	2.05	113.85	107.94
3	A	503	3PG	O1P-C3-C2	2.03	113.79	107.94

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	504	3PG	C1-C2-C3-O1P
3	A	503	3PG	C1-C2-C3-O1P
3	B	504	3PG	O3-C2-C3-O1P
3	A	503	3PG	O3-C2-C3-O1P

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	507	MPD	1	0
5	A	509	MPD	1	0
4	A	501	GOL	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	405/410 (98%)	0.19	22 (5%) 25 31	22, 36, 60, 78	0
1	B	405/410 (98%)	0.60	39 (9%) 8 10	22, 46, 81, 93	0
All	All	810/820 (98%)	0.40	61 (7%) 14 18	22, 40, 72, 93	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	279	TYR	6.9
1	B	297	HIS	6.7
1	B	299	TYR	6.4
1	B	284	VAL	5.3
1	B	280	LYS	4.3
1	B	245	PHE	4.2
1	A	54	ILE	3.4
1	B	296	LEU	3.4
1	B	281	GLY	3.4
1	A	27	GLY	3.3
1	B	405	PHE	3.3
1	B	282	GLU	3.2
1	B	214	ARG	3.1
1	B	286	ILE	3.1
1	B	292	ASN	3.1
1	B	336	PHE	3.0
1	A	16	LEU	3.0
1	A	113	LEU	2.9
1	A	288	LEU	2.9
1	B	244	GLU	2.8
1	B	351	PRO	2.8
1	A	26	ASP	2.7
1	A	402	TYR	2.6
1	B	298	GLN	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	272	PRO	2.6
1	B	240	ARG	2.6
1	B	263	ASP	2.6
1	A	18	VAL	2.6
1	A	279	TYR	2.5
1	B	403	GLU	2.5
1	A	55	GLY	2.5
1	A	248	LYS	2.4
1	B	110	LEU	2.4
1	A	90	ILE	2.4
1	B	294	GLY	2.4
1	A	292	ASN	2.3
1	A	195	VAL	2.3
1	A	53	VAL	2.3
1	B	249	LYS	2.3
1	A	145	VAL	2.3
1	B	369	GLY	2.3
1	B	295	LEU	2.2
1	A	110	LEU	2.2
1	B	333	ARG	2.2
1	B	306	LYS	2.2
1	B	278	ASP	2.2
1	B	196	LEU	2.2
1	B	334	GLU	2.2
1	B	338	ILE	2.2
1	B	289	LEU	2.2
1	B	236	PHE	2.2
1	B	293	ARG	2.2
1	B	16	LEU	2.1
1	A	249	LYS	2.1
1	B	18	VAL	2.1
1	B	277	VAL	2.1
1	A	56	THR	2.1
1	A	17	ARG	2.1
1	A	81	ASP	2.0
1	A	1	MET	2.0
1	B	314	GLU	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 [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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(\AA^2)	Q<0.9
5	MPD	A	509	8/8	0.82	0.29	31,42,46,49	0
4	GOL	B	502	6/6	0.85	0.17	33,40,43,46	0
2	CL	B	506	1/1	0.87	0.09	68,68,68,68	0
5	MPD	B	507	8/8	0.88	0.17	42,44,45,45	0
4	GOL	A	501	6/6	0.88	0.17	39,47,51,54	0
5	MPD	A	508	8/8	0.88	0.17	41,45,46,46	0
2	CL	A	505	1/1	0.90	0.13	64,64,64,64	0
3	3PG	A	503	11/11	0.93	0.15	58,63,69,70	0
3	3PG	B	504	11/11	0.94	0.17	37,47,59,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.