



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 06:11 AM UTC

PDB ID : 7PG5 / pdb_00007pg5
Title : Crystal Structure of PI3Kalpha
Authors : Gong, G.; Pinotsis, N.; Williams, R.L.; Vanhaesebroeck, B.
Deposited on : 2021-08-13
Resolution : 2.20 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

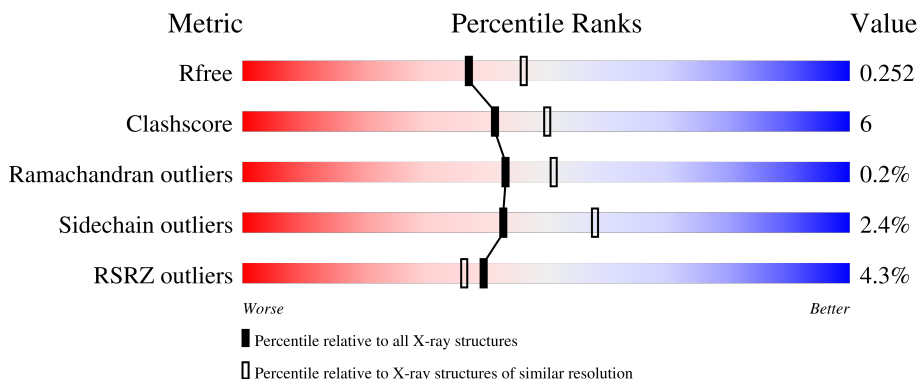
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	1068	
2	B	287	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 10787 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 Phosphatidylinositol 4,5-bisphosphate 3-kinase catalytic subunit alpha isoform.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	1028	8411	5377	1447	1520	67	0	4	0

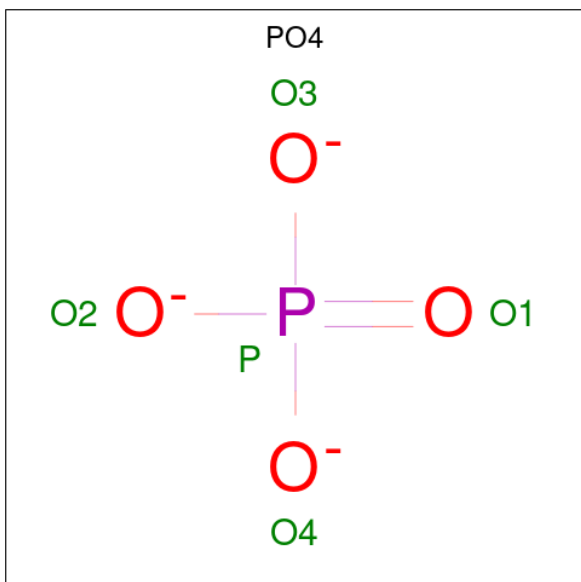
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	232	LYS	MET	engineered mutation	UNP P42336
A	233	LYS	LEU	engineered mutation	UNP P42336

- Molecule 2 is a protein called Phosphatidylinositol 3-kinase regulatory subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	266	2208	1381	392	430	5	0	0	0

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O P 5 4 1	0	0
3	A	1	Total O P 5 4 1	0	0

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

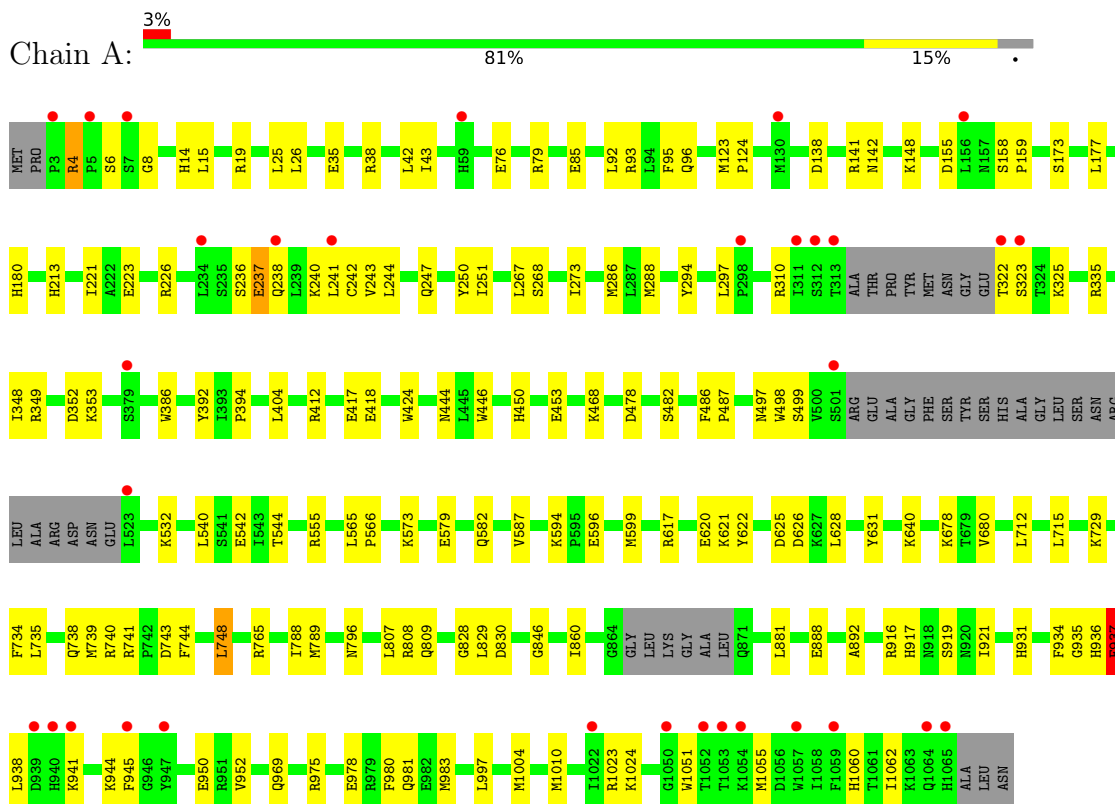
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	143	Total O 143 143	0	0
5	B	9	Total O 9 9	0	0

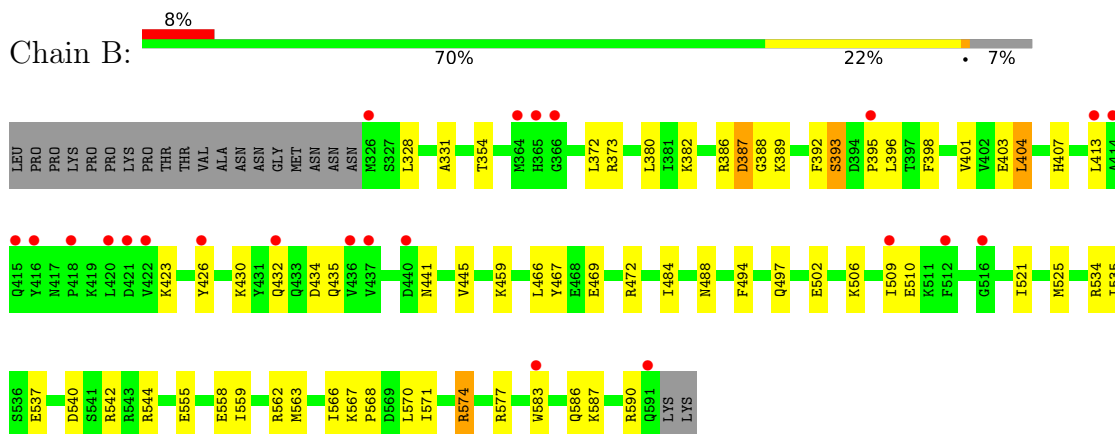
3 Residue-property plots [i](#)

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: Phosphatidylinositol 4,5-bisphosphate 3-kinase catalytic subunit alpha isoform



- Molecule 2: Phosphatidylinositol 3-kinase regulatory subunit alpha



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	105.43Å 105.24Å 136.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	57.17 – 2.20 57.17 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (57.17-2.20) 99.9 (57.17-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 2.20Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.196 , 0.246 0.208 , 0.252	Depositor DCC
R_{free} test set	3986 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å ²)	51.2	Xtrriage
Anisotropy	0.117	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 37.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.014 for k,h,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10787	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, GOL

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	0.56	1/8601 (0.0%)	0.82	9/11627 (0.1%)
2	B	0.50	0/2245	0.80	0/3016
All	All	0.55	1/10846 (0.0%)	0.81	9/14643 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	498	TRP	CA-C	-6.03	1.45	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	499	SER	N-CA-C	-6.42	105.76	113.97
1	A	952	VAL	CA-C-N	5.45	125.10	119.05
1	A	952	VAL	C-N-CA	5.45	125.10	119.05
1	A	937	PHE	N-CA-C	5.30	115.13	108.45
1	A	498	TRP	N-CA-C	-5.28	105.50	112.94
1	A	4	ARG	CA-C-N	-5.27	114.49	119.76
1	A	4	ARG	C-N-CA	-5.27	114.49	119.76
1	A	26	LEU	CA-C-N	5.11	125.18	119.87
1	A	26	LEU	C-N-CA	5.11	125.18	119.87

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	8411	0	8388	97	0
2	B	2208	0	2138	46	0
3	A	10	0	0	1	0
4	A	6	0	8	0	0
5	A	143	0	0	9	0
5	B	9	0	0	1	0
All	All	10787	0	10534	135	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 6.

All (135) 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:625:ASP:OD1	5:A:1201:HOH:O	1.82	0.98
1:A:888:GLU:O	5:A:1202:HOH:O	1.93	0.85
1:A:729:LYS:O	5:A:1203:HOH:O	1.98	0.80
2:B:558:GLU:OE1	5:B:601:HOH:O	1.99	0.78
1:A:453:GLU:OE2	2:B:574:ARG:NH2	2.19	0.75
1:A:223:GLU:OE1	5:A:1204:HOH:O	2.05	0.74
1:A:975:ARG:NH1	5:A:1207:HOH:O	2.19	0.73
2:B:534:ARG:NH1	2:B:537:GLU:OE2	2.21	0.73
1:A:322:THR:HG23	1:A:323:SER:H	1.53	0.71
1:A:1055:MET:HB3	1:A:1060:HIS:HA	1.73	0.70
1:A:741[A]:ARG:NH2	1:A:743:ASP:OD2	2.26	0.69
1:A:286:MET:HE2	1:A:288:MET:HE2	1.75	0.69
1:A:444:ASN:ND2	1:A:468:LYS:HA	2.06	0.69
1:A:236:SER:OG	1:A:240:LYS:NZ	2.28	0.67
1:A:892:ALA:O	5:A:1205:HOH:O	2.11	0.66
1:A:497:ASN:OD1	1:A:497:ASN:N	2.28	0.65
2:B:583:TRP:CZ3	2:B:587:LYS:HG3	2.33	0.64
1:A:808:ARG:HB3	1:A:1010:MET:HE2	1.80	0.64
1:A:142:ASN:ND2	5:A:1211:HOH:O	2.31	0.63
1:A:941:LYS:HD2	1:A:945:PHE:HE2	1.63	0.63
2:B:328:LEU:HD13	2:B:401:VAL:HB	1.80	0.63
1:A:15:LEU:HD11	1:A:738:GLN:NE2	2.14	0.62
2:B:540:ASP:OD2	2:B:544:ARG:NH1	2.33	0.62
1:A:807:LEU:HD12	1:A:846:GLY:HA3	1.82	0.61
1:A:916:ARG:HB3	1:A:921:ILE:HD11	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:741[A]:ARG:HH21	1:A:743:ASP:CG	2.10	0.59
1:A:599:MET:HE2	1:A:1004:MET:HE1	1.85	0.59
2:B:494:PHE:HB3	2:B:535:ILE:HG12	1.85	0.58
1:A:180:HIS:HD1	1:A:828:GLY:HA2	1.69	0.58
1:A:917:HIS:CE1	1:A:919:SER:HB2	2.40	0.57
2:B:413:LEU:HB2	2:B:423:LYS:H	1.70	0.57
1:A:712:LEU:HD21	1:A:748:LEU:HD21	1.86	0.56
2:B:331:ALA:O	2:B:430:LYS:NZ	2.38	0.56
1:A:1062:ILE:H	1:A:1062:ILE:HD12	1.71	0.56
1:A:138:ASP:OD1	1:A:141:ARG:NH2	2.36	0.56
1:A:236:SER:O	1:A:240:LYS:HG2	2.07	0.55
1:A:715:LEU:HD21	1:A:735:LEU:HD12	1.89	0.55
1:A:155:ASP:OD1	1:A:158:SER:OG	2.21	0.55
1:A:944:LYS:O	2:B:577:ARG:NH1	2.39	0.55
1:A:76:GLU:OE2	1:A:93:ARG:NH1	2.40	0.54
1:A:596:GLU:HG2	1:A:997:LEU:HD13	1.90	0.54
1:A:180:HIS:ND1	1:A:828:GLY:HA2	2.23	0.54
1:A:15:LEU:HD11	1:A:738:GLN:HE22	1.73	0.54
2:B:467:TYR:HD1	2:B:563:MET:HE1	1.72	0.54
1:A:325:LYS:HE2	1:A:482:SER:HB3	1.91	0.53
1:A:444:ASN:HD22	1:A:468:LYS:HA	1.72	0.53
2:B:434:ASP:OD1	2:B:583:TRP:NE1	2.40	0.53
1:A:941:LYS:HD2	1:A:945:PHE:CE2	2.42	0.53
1:A:444:ASN:ND2	1:A:468:LYS:HD3	2.25	0.52
1:A:599:MET:HE1	1:A:631:TYR:HB3	1.91	0.52
1:A:734:PHE:O	1:A:738:GLN:HG2	2.10	0.52
1:A:348:ILE:HD12	1:A:348:ILE:H	1.74	0.52
1:A:620:GLU:OE2	5:A:1206:HOH:O	2.19	0.52
2:B:393:SER:HB3	2:B:395:PRO:HD2	1.91	0.52
2:B:392:PHE:CE1	2:B:404:LEU:HD21	2.45	0.52
1:A:323:SER:O	1:A:482:SER:OG	2.26	0.52
1:A:788:ILE:HG23	1:A:789:MET:HG2	1.90	0.52
2:B:398:PHE:HZ	2:B:407:HIS:CD2	2.28	0.52
1:A:267:LEU:HG	1:A:273:ILE:HG13	1.92	0.51
1:A:38:ARG:HH12	1:A:741[A]:ARG:HE	1.59	0.51
2:B:509:ILE:HG13	2:B:510:GLU:N	2.26	0.50
2:B:445:VAL:HG21	2:B:583:TRP:CZ3	2.46	0.50
1:A:19:ARG:NH1	1:A:35:GLU:OE2	2.44	0.50
1:A:544:THR:HG21	2:B:382:LYS:HB2	1.93	0.50
1:A:213:HIS:HB2	1:A:268:SER:HB3	1.93	0.49
1:A:335:ARG:HG2	1:A:386:TRP:CE3	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:413:LEU:HB2	2:B:423:LYS:HA	1.95	0.49
2:B:445:VAL:HG21	2:B:583:TRP:CE3	2.48	0.49
1:A:555:ARG:NH2	1:A:582:GLN:OE1	2.46	0.48
1:A:937:PHE:CD1	1:A:938:LEU:HG	2.48	0.48
2:B:567:LYS:O	2:B:571:ILE:HG12	2.14	0.48
1:A:542:GLU:HB3	2:B:380:LEU:HD13	1.96	0.48
2:B:562:ARG:O	2:B:566:ILE:HG23	2.13	0.48
2:B:469:GLU:OE1	2:B:472:ARG:NH2	2.46	0.48
1:A:43:ILE:HG13	1:A:85:GLU:O	2.14	0.47
1:A:934:PHE:O	1:A:936:HIS:N	2.47	0.47
1:A:349:ARG:O	1:A:412:ARG:NH2	2.37	0.47
1:A:352:ASP:O	1:A:353:LYS:HG2	2.14	0.47
1:A:450:HIS:HB2	2:B:467:TYR:CZ	2.50	0.47
1:A:621:LYS:HB3	1:A:622:TYR:CD1	2.50	0.47
1:A:640:LYS:HG2	1:A:680:VAL:HG11	1.98	0.46
1:A:180:HIS:CD2	1:A:830:ASP:HB2	2.50	0.46
2:B:459:LYS:HG2	2:B:570:LEU:HD13	1.97	0.46
1:A:6:SER:HB3	1:A:14:HIS:CE1	2.51	0.46
1:A:916:ARG:HD2	1:A:931:HIS:ND1	2.31	0.46
2:B:559:ILE:O	2:B:563:MET:HG3	2.17	0.45
1:A:739:MET:HA	1:A:744:PHE:CD1	2.52	0.45
1:A:765:ARG:NH2	1:A:796:ASN:OD1	2.49	0.45
2:B:354:THR:HA	2:B:426:TYR:O	2.17	0.45
1:A:95:PHE:CD2	1:A:96:GLN:HG2	2.52	0.44
2:B:387:ASP:OD1	2:B:388:GLY:N	2.47	0.44
1:A:565:LEU:HB3	1:A:566:PRO:HD3	2.00	0.44
1:A:453:GLU:CD	2:B:574:ARG:HH22	2.23	0.44
1:A:540:LEU:HB2	1:A:1023:ARG:HD2	1.98	0.44
1:A:221:ILE:HD13	1:A:250:TYR:HB2	2.00	0.44
2:B:413:LEU:HB2	2:B:423:LYS:N	2.33	0.44
1:A:238:GLN:O	1:A:242:CYS:N	2.47	0.44
1:A:678:LYS:HA	1:A:678:LYS:HD2	1.85	0.43
1:A:981:GLN:OE1	1:A:1051:TRP:HA	2.18	0.43
1:A:740:ARG:HE	1:A:740:ARG:HB2	1.58	0.43
1:A:42:LEU:HD21	1:A:92:LEU:HD11	2.01	0.43
1:A:860:ILE:HA	3:A:1101:PO4:O2	2.19	0.43
1:A:226:ARG:NE	5:A:1204:HOH:O	2.18	0.43
1:A:617[B]:ARG:HH11	1:A:617[B]:ARG:HG2	1.84	0.43
2:B:441:ASN:OD1	2:B:441:ASN:N	2.46	0.42
1:A:294:TYR:HA	1:A:297:LEU:HD12	2.00	0.42
2:B:534:ARG:HA	2:B:534:ARG:HD2	1.78	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:389:LYS:HA	2:B:398:PHE:O	2.19	0.42
1:A:123:MET:HA	1:A:124:PRO:HD3	1.97	0.42
1:A:418:GLU:HB3	2:B:568:PRO:HB2	2.01	0.42
2:B:555:GLU:O	2:B:559:ILE:HG13	2.19	0.42
1:A:573:LYS:HD3	1:A:579:GLU:OE1	2.20	0.42
1:A:25:LEU:HB3	2:B:497:GLN:CD	2.45	0.41
1:A:392:TYR:CG	1:A:394:PRO:HD2	2.55	0.41
1:A:444:ASN:HD21	1:A:468:LYS:HD3	1.85	0.41
2:B:488:ASN:OD1	2:B:542:ARG:NH1	2.50	0.41
2:B:386:ARG:HB2	2:B:396:LEU:CD1	2.51	0.41
1:A:322:THR:HG23	1:A:323:SER:N	2.30	0.41
1:A:594:LYS:HA	1:A:594:LYS:HD3	1.98	0.41
2:B:413:LEU:CB	2:B:423:LYS:H	2.34	0.41
2:B:502:GLU:O	2:B:506:LYS:HG3	2.21	0.41
1:A:8:GLY:HA3	1:A:14:HIS:HA	2.02	0.41
1:A:937:PHE:CD1	1:A:937:PHE:C	2.99	0.41
1:A:980:PHE:O	1:A:983:MET:HB2	2.20	0.41
1:A:159:PRO:HG2	1:A:294:TYR:CD2	2.56	0.41
1:A:424:TRP:CE2	1:A:446:TRP:HB2	2.55	0.41
1:A:486:PHE:CG	1:A:487:PRO:HD2	2.55	0.41
1:A:950:GLU:OE2	1:A:1024:LYS:HE3	2.20	0.41
2:B:432:GLN:O	2:B:435:GLN:HG2	2.20	0.41
2:B:398:PHE:HD2	2:B:403:GLU:HB3	1.86	0.41
2:B:372:LEU:HD12	2:B:373:ARG:N	2.37	0.40
2:B:521:ILE:O	2:B:525:MET:HG3	2.21	0.40
2:B:586:GLN:HG3	2:B:590:ARG:HH21	1.85	0.40
1:A:237:GLU:O	1:A:241:LEU:HB2	2.21	0.40
1:A:404:LEU:C	1:A:404:LEU:HD12	2.46	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	1024/1068 (96%)	993 (97%)	29 (3%)	2 (0%)	43	51
2	B	264/287 (92%)	254 (96%)	9 (3%)	1 (0%)	30	34
All	All	1288/1355 (95%)	1247 (97%)	38 (3%)	3 (0%)	43	51

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	935	GLY
2	B	387	ASP
1	A	4	ARG

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	936/974 (96%)	913 (98%)	23 (2%)	42	56
2	B	235/266 (88%)	230 (98%)	5 (2%)	47	63
All	All	1171/1240 (94%)	1143 (98%)	28 (2%)	43	58

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	ARG
1	A	148	LYS
1	A	173	SER
1	A	177	LEU
1	A	237	GLU
1	A	243	VAL
1	A	244	LEU
1	A	247	GLN
1	A	251	ILE
1	A	310	ARG
1	A	417	GLU
1	A	478	ASP
1	A	532	LYS

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Mol	Chain	Res	Type
1	A	587	VAL
1	A	626	ASP
1	A	628	LEU
1	A	748	LEU
1	A	809	GLN
1	A	829	LEU
1	A	881	LEU
1	A	937	PHE
1	A	969	GLN
1	A	978	GLU
2	B	393	SER
2	B	404	LEU
2	B	466	LEU
2	B	484	ILE
2	B	574	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	145	ASN
1	A	180	HIS
1	A	189	GLN
1	A	247	GLN
1	A	384	ASN
1	A	738	GLN
1	A	825	GLN
1	A	855	HIS
1	A	861	GLN
1	A	1033	GLN
1	A	1042	GLN
2	B	385	HIS
2	B	406	ASN
2	B	407	HIS
2	B	432	GLN
2	B	501	GLN

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](#)

3 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
3	PO4	A	1102	-	4,4,4	0.64	0	6,6,6	0.44	0
4	GOL	A	1103	-	5,5,5	0.72	0	5,5,5	2.42	3 (60%)
3	PO4	A	1101	-	4,4,4	1.07	0	6,6,6	0.61	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	GOL	A	1103	-	-	0/4/4/4	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1103	GOL	O2-C2-C3	-3.45	94.91	109.18
4	A	1103	GOL	O3-C3-C2	-2.59	98.72	110.38
4	A	1103	GOL	O1-C1-C2	-2.39	99.62	110.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1101	PO4	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	1028/1068 (96%)	0.23	32 (3%) 51 48	28, 60, 91, 124	4 (0%)
2	B	266/287 (92%)	0.67	23 (8%) 16 14	48, 77, 120, 130	0
All	All	1294/1355 (95%)	0.32	55 (4%) 40 36	28, 64, 104, 130	4 (0%)

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	311	ILE	4.6
1	A	313	THR	4.3
1	A	1065	HIS	4.2
2	B	418	PRO	3.8
1	A	501	SER	3.7
1	A	234	LEU	3.6
1	A	947	TYR	3.6
1	A	1053	THR	3.4
2	B	422	VAL	3.4
1	A	1057	TRP	3.3
1	A	940[A]	HIS	3.3
1	A	941	LYS	3.2
2	B	591	GLN	3.2
2	B	436	VAL	3.2
1	A	523	LEU	3.2
2	B	420	LEU	3.0
1	A	322	THR	3.0
1	A	312	SER	3.0
1	A	1054	LYS	2.9
1	A	241	LEU	2.9
1	A	323	SER	2.9
1	A	238	GLN	2.8
2	B	414	ALA	2.8
2	B	437	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
2	B	326	MET	2.7
2	B	365	HIS	2.7
2	B	364	MET	2.6
1	A	7[A]	SER	2.6
2	B	395	PRO	2.6
2	B	416	TYR	2.6
2	B	583	TRP	2.5
2	B	426	TYR	2.5
1	A	939	ASP	2.4
1	A	1059	PHE	2.4
1	A	1064	GLN	2.4
1	A	130	MET	2.3
2	B	415	GLN	2.3
2	B	509	ILE	2.3
1	A	5	PRO	2.3
1	A	379	SER	2.3
1	A	156	LEU	2.3
1	A	59	HIS	2.2
1	A	1052	THR	2.2
2	B	432	GLN	2.2
1	A	945	PHE	2.2
1	A	3	PRO	2.2
2	B	421	ASP	2.2
2	B	512	PHE	2.1
2	B	413	LEU	2.1
1	A	1050	GLY	2.1
2	B	440	ASP	2.1
2	B	516	GLY	2.1
1	A	298	PRO	2.0
2	B	366	GLY	2.0
1	A	1022	ILE	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(Å ²)	Q<0.9
3	PO4	A	1101	5/5	0.85	0.17	77,82,83,85	5
3	PO4	A	1102	5/5	0.88	0.18	72,73,83,90	5
4	GOL	A	1103	6/6	0.90	0.41	66,72,79,79	6

6.5 Other polymers [i](#)

There are no such residues in this entry.