



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 9, 2017 – 10:47 AM EST

PDB ID : 5SWG
Title : Crystal Structure of PI3Kalpha in complex with fragments 5 and 21
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Deposited on : 2016-08-08
Resolution : 3.11 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20028442
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20028442

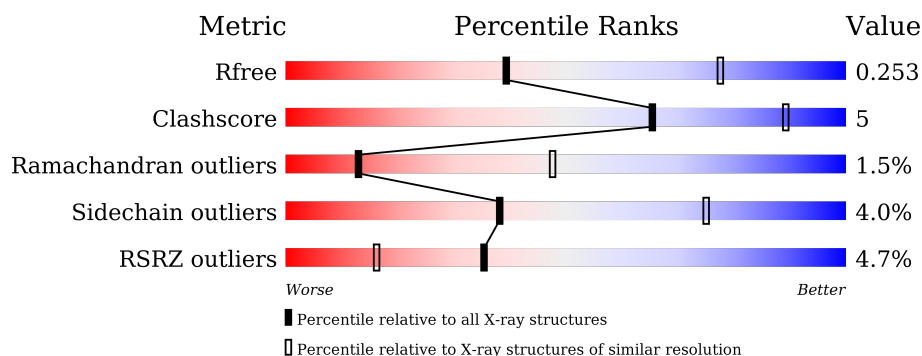
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.11 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1112 (3.16-3.08)
Clashscore	102246	1218 (3.16-3.08)
Ramachandran outliers	100387	1175 (3.16-3.08)
Sidechain outliers	100360	1175 (3.16-3.08)
RSRZ outliers	91569	1114 (3.16-3.08)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1096	<div> <div>3%</div> <div>82%</div> <div>11%</div> <div>5%</div> </div>
2	B	279	<div> <div>11%</div> <div>77%</div> <div>11%</div> <div>9%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	AX7	A	1101	-	-	-	X
4	CAQ	B	701	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10723 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 sub-unit alpha isoform.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	1042	Total	C	N	O	P	S	0	0	0
			8530	5453	1458	1549	1	69			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-27	MET	-	initiating methionine	UNP P42336
A	-26	SER	-	expression tag	UNP P42336
A	-25	TYR	-	expression tag	UNP P42336
A	-24	TYR	-	expression tag	UNP P42336
A	-23	HIS	-	expression tag	UNP P42336
A	-22	HIS	-	expression tag	UNP P42336
A	-21	HIS	-	expression tag	UNP P42336
A	-20	HIS	-	expression tag	UNP P42336
A	-19	HIS	-	expression tag	UNP P42336
A	-18	HIS	-	expression tag	UNP P42336
A	-17	ASP	-	expression tag	UNP P42336
A	-16	TYR	-	expression tag	UNP P42336
A	-15	ASP	-	expression tag	UNP P42336
A	-14	ILE	-	expression tag	UNP P42336
A	-13	PRO	-	expression tag	UNP P42336
A	-12	THR	-	expression tag	UNP P42336
A	-11	THR	-	expression tag	UNP P42336
A	-10	GLU	-	expression tag	UNP P42336
A	-9	ASN	-	expression tag	UNP P42336
A	-8	LEU	-	expression tag	UNP P42336
A	-7	TYR	-	expression tag	UNP P42336
A	-6	PHE	-	expression tag	UNP P42336
A	-5	GLN	-	expression tag	UNP P42336
A	-4	GLY	-	expression tag	UNP P42336
A	-3	ALA	-	expression tag	UNP P42336
A	-2	MET	-	expression tag	UNP P42336

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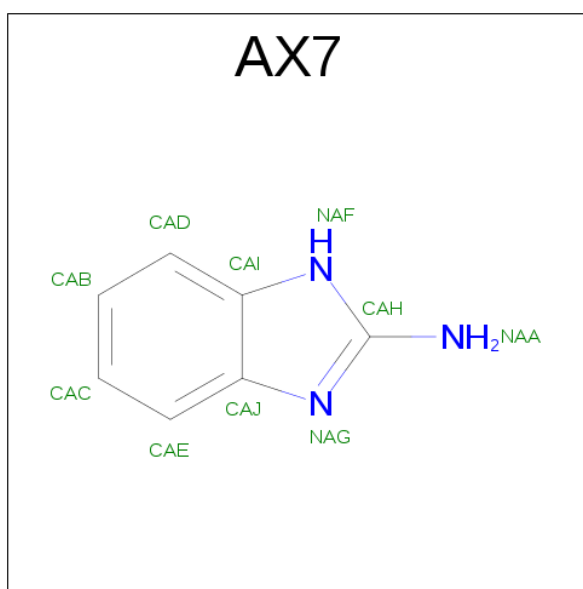
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Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P42336
A	0	SER	-	expression tag	UNP P42336

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

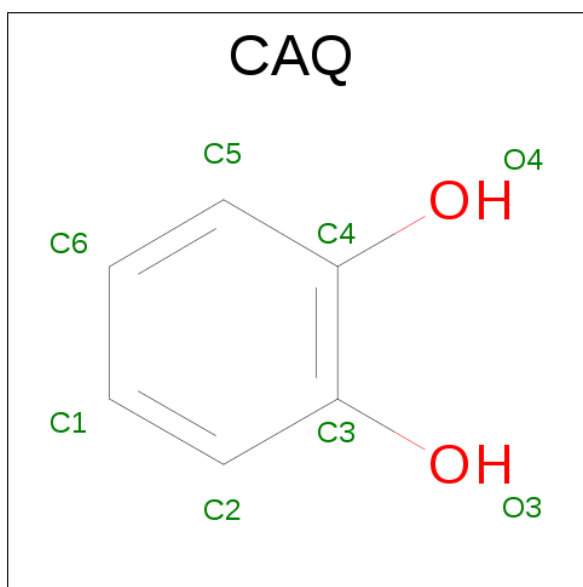
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	255	Total	C	N	O	S	0	0	0
			2175	1371	385	413	6			

- Molecule 3 is 1H-benzimidazol-2-amine (three-letter code: AX7) (formula: C₇H₇N₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			10	7	3		

- Molecule 4 is CATECHOL (three-letter code: CAQ) (formula: C₆H₆O₂).

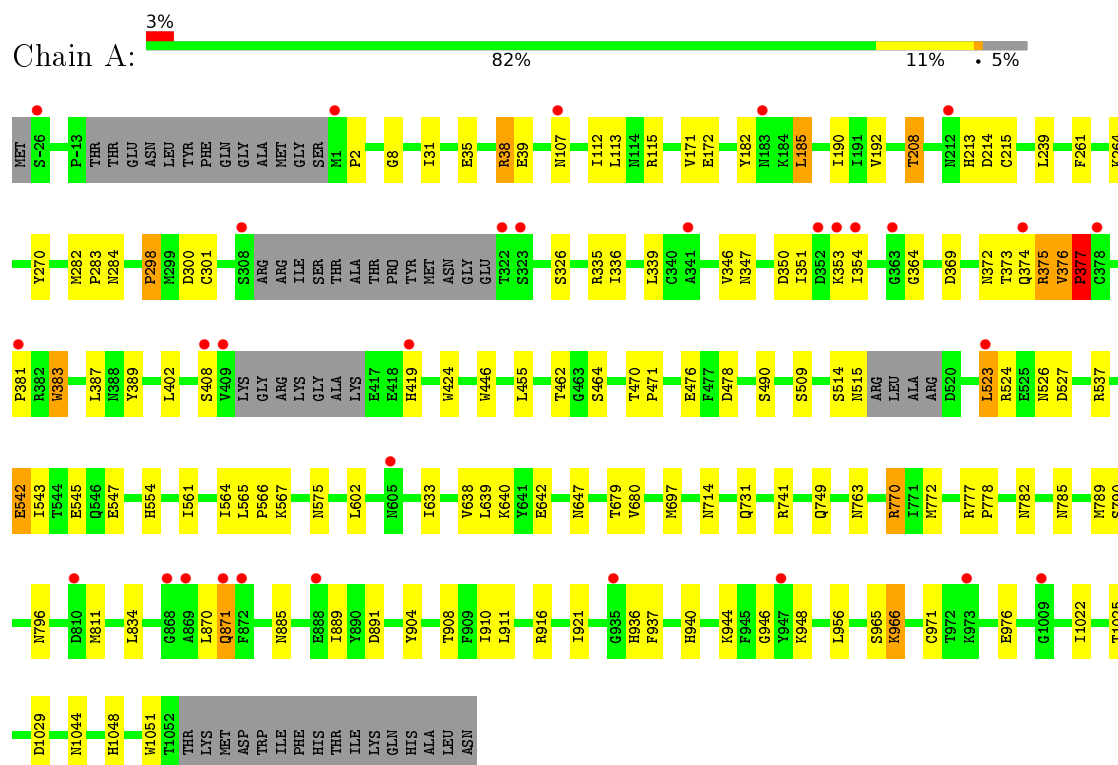


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			8	6	2		

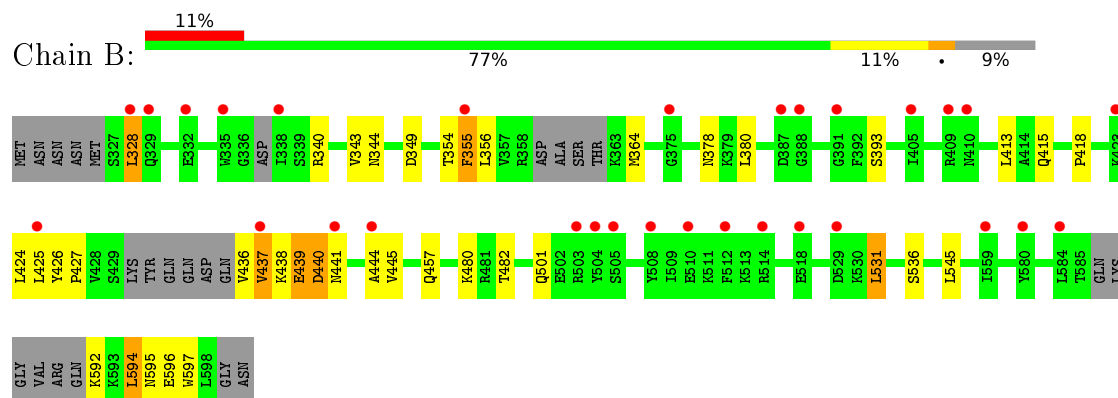
3 Residue-property plots

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: 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	114.51Å 116.68Å 149.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	91.97 – 3.11 49.10 – 3.11	Depositor EDS
% Data completeness (in resolution range)	99.8 (91.97-3.11) 99.8 (49.10-3.11)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 3.12Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.206 , 0.262 0.205 , 0.253	Depositor DCC
R_{free} test set	1866 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	89.8	Xtriage
Anisotropy	0.175	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 62.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.010 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10723	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.99% 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: CAQ, AX7, SEP

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	0.50	0/8719	0.75	1/11786 (0.0%)
2	B	0.46	0/2210	0.70	1/2955 (0.0%)
All	All	0.49	0/10929	0.74	2/14741 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	364	GLY	N-CA-C	-5.16	100.19	113.10
2	B	328	LEU	CA-CB-CG	5.08	126.98	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	113	LEU	Peptide

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	8530	0	8495	80	0
2	B	2175	0	2161	32	0
3	A	10	0	7	0	0
4	B	8	0	5	0	0
All	All	10723	0	10668	107	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 5.

All (107) 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:353:LYS:HB3	1:A:377:PRO:HG3	1.26	1.08
1:A:298:PRO:CG	1:A:697:MET:HG3	1.97	0.93
1:A:346:VAL:HG13	1:A:351:ILE:HG13	1.58	0.84
2:B:594:LEU:HD12	2:B:594:LEU:H	1.46	0.80
1:A:298:PRO:HG2	1:A:697:MET:HG3	1.63	0.78
1:A:353:LYS:CB	1:A:377:PRO:HG3	2.14	0.76
1:A:523:LEU:CD1	1:A:527:ASP:HB2	2.16	0.76
1:A:353:LYS:HB3	1:A:377:PRO:CG	2.13	0.75
1:A:523:LEU:HD12	1:A:524:ARG:N	2.04	0.72
2:B:438:LYS:O	2:B:440:ASP:N	2.24	0.70
2:B:439:GLU:C	2:B:441:ASN:H	2.01	0.64
1:A:353:LYS:HD3	1:A:374:GLN:HB3	1.80	0.64
1:A:298:PRO:CB	1:A:697:MET:HG3	2.29	0.63
1:A:870:LEU:O	1:A:871:GLN:HG3	1.98	0.62
1:A:911:LEU:O	1:A:937:PHE:O	2.18	0.62
2:B:439:GLU:O	2:B:441:ASN:N	2.33	0.62
1:A:372:ASN:HB3	1:A:387:LEU:HD21	1.83	0.61
1:A:300:ASP:OD1	1:A:301:CYS:N	2.33	0.61
1:A:213:HIS:CE1	1:A:214:ASP:HB3	2.36	0.61
1:A:298:PRO:HG2	1:A:697:MET:CG	2.29	0.60
1:A:376:VAL:CG2	1:A:381:PRO:HA	2.33	0.58
1:A:31:ILE:HD11	2:B:531:LEU:HD13	1.85	0.58
2:B:594:LEU:O	2:B:597:TRP:N	2.37	0.57
1:A:543:ILE:HD11	1:A:567:LYS:HD3	1.87	0.57
1:A:336:ILE:HD13	1:A:402:LEU:HD22	1.86	0.56
1:A:346:VAL:HG13	1:A:351:ILE:CG1	2.33	0.56
1:A:339:LEU:O	1:A:383:TRP:O	2.23	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:436:VAL:O	2:B:436:VAL:HG23	2.06	0.55
1:A:642:GLU:HG2	1:A:647:ASN:CG	2.26	0.55
2:B:437:VAL:HB	2:B:439:GLU:HG3	1.89	0.55
1:A:870:LEU:O	1:A:871:GLN:CB	2.55	0.54
1:A:523:LEU:C	1:A:523:LEU:HD12	2.27	0.54
2:B:355:PHE:CD2	2:B:355:PHE:C	2.82	0.53
1:A:353:LYS:HA	1:A:377:PRO:HB3	1.91	0.53
1:A:936:HIS:HB3	1:A:940:HIS:HB3	1.90	0.53
1:A:870:LEU:O	1:A:871:GLN:CG	2.57	0.52
2:B:355:PHE:C	2:B:355:PHE:HD2	2.13	0.52
2:B:594:LEU:O	2:B:596:GLU:N	2.42	0.52
1:A:523:LEU:HD12	1:A:524:ARG:O	2.10	0.52
1:A:537:ARG:NH1	1:A:547:GLU:OE1	2.42	0.51
1:A:772:MET:HB2	1:A:778:PRO:HG2	1.92	0.51
1:A:514:SER:O	1:A:515:ASN:ND2	2.44	0.50
1:A:523:LEU:HD22	1:A:554:HIS:CD2	2.46	0.50
2:B:355:PHE:CZ	2:B:427:PRO:HB3	2.47	0.50
2:B:340:ARG:O	2:B:344:ASN:ND2	2.45	0.49
1:A:354:ILE:HG22	1:A:408:SER:HA	1.93	0.49
1:A:602:LEU:HB3	1:A:638:VAL:HG11	1.94	0.49
1:A:373:THR:HG21	1:A:375:ARG:NH1	2.27	0.49
1:A:561:ILE:O	1:A:564:ILE:HG22	2.12	0.48
2:B:439:GLU:C	2:B:441:ASN:N	2.66	0.48
2:B:355:PHE:CD1	2:B:424:LEU:HB3	2.49	0.48
2:B:437:VAL:CG1	2:B:439:GLU:HG3	2.44	0.48
1:A:782:ASN:OD1	1:A:796:ASN:ND2	2.47	0.48
1:A:346:VAL:CG1	1:A:351:ILE:HG13	2.37	0.47
1:A:910:ILE:O	1:A:1025:THR:HG21	2.15	0.47
1:A:916:ARG:HB3	1:A:921:ILE:HD11	1.97	0.47
1:A:336:ILE:HD12	1:A:389:TYR:CE2	2.50	0.46
1:A:640:LYS:CE	1:A:680:VAL:HG11	2.46	0.46
1:A:731:GLN:OE1	1:A:777:ARG:NH2	2.47	0.46
1:A:171:VAL:HG12	1:A:172:GLU:O	2.15	0.46
1:A:885:ASN:HB3	1:A:889:ILE:HG22	1.98	0.46
2:B:439:GLU:HB2	2:B:445:VAL:CG2	2.45	0.46
1:A:353:LYS:HD3	1:A:374:GLN:CB	2.44	0.46
1:A:904:TYR:O	1:A:908:THR:HB	2.16	0.45
2:B:355:PHE:CE1	2:B:424:LEU:HB3	2.50	0.45
1:A:789:MET:O	1:A:790:SEP:C	2.64	0.45
1:A:298:PRO:HB2	1:A:697:MET:HG3	1.98	0.45
1:A:192:VAL:HG13	1:A:283:PRO:HB2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:261:PHE:HA	1:A:270:TYR:CE2	2.52	0.44
1:A:542:GLU:OE2	2:B:340:ARG:NE	2.50	0.44
2:B:594:LEU:CD1	2:B:594:LEU:H	2.21	0.44
2:B:592:LYS:N	2:B:594:LEU:HD13	2.33	0.44
1:A:602:LEU:O	1:A:638:VAL:HG13	2.17	0.44
1:A:8:GLY:HA3	1:A:714:ASN:HD21	1.83	0.44
2:B:436:VAL:C	2:B:437:VAL:HG23	2.38	0.44
1:A:642:GLU:HG2	1:A:647:ASN:ND2	2.33	0.44
1:A:870:LEU:O	1:A:871:GLN:HB2	2.18	0.44
1:A:542:GLU:HG2	2:B:380:LEU:HD22	2.00	0.43
1:A:1022:ILE:HA	1:A:1025:THR:HG22	2.01	0.43
2:B:439:GLU:HB2	2:B:445:VAL:HG23	1.99	0.43
1:A:419:HIS:HB3	1:A:455:LEU:HD21	2.01	0.43
1:A:1044:ASN:ND2	1:A:1051:TRP:O	2.52	0.43
2:B:354:THR:HA	2:B:426:TYR:HB2	2.01	0.43
1:A:944:LYS:HD2	2:B:457:GLN:HG2	2.01	0.43
1:A:347:ASN:HB2	1:A:350:ASP:HB2	2.00	0.43
1:A:376:VAL:N	1:A:377:PRO:CD	2.80	0.43
1:A:749:GLN:HG2	1:A:763:ASN:HA	2.01	0.43
1:A:373:THR:HG21	1:A:375:ARG:CZ	2.49	0.42
1:A:1029:ASP:OD1	2:B:340:ARG:NH2	2.51	0.42
1:A:777:ARG:N	1:A:778:PRO:CD	2.82	0.42
1:A:965:SER:HA	1:A:976:GLU:HG3	2.02	0.42
2:B:344:ASN:OD1	2:B:378:ASN:ND2	2.53	0.42
1:A:565:LEU:HB3	1:A:566:PRO:HD3	2.01	0.42
1:A:182:TYR:O	1:A:185:LEU:HD12	2.20	0.42
1:A:38:ARG:HD3	1:A:741:ARG:NH1	2.35	0.41
1:A:770:ARG:HG3	1:A:772:MET:CE	2.50	0.41
1:A:772:MET:CE	1:A:772:MET:HA	2.51	0.41
2:B:441:ASN:HB2	2:B:444:ALA:HB3	2.02	0.41
1:A:282:MET:O	1:A:284:ASN:ND2	2.54	0.41
1:A:213:HIS:ND1	1:A:214:ASP:N	2.68	0.41
2:B:480:LYS:HG2	2:B:545:LEU:HD11	2.04	0.40
1:A:523:LEU:HD13	1:A:527:ASP:HB2	1.98	0.40
2:B:343:VAL:HG13	2:B:356:LEU:HD11	2.02	0.40
1:A:523:LEU:CD1	1:A:524:ARG:O	2.69	0.40
1:A:424:TRP:CE2	1:A:446:TRP:HB2	2.56	0.40
1:A:633:ILE:HD11	1:A:811:MET:HE2	2.04	0.40
2:B:413:LEU:HD11	2:B:424:LEU:HD21	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	1031/1096 (94%)	941 (91%)	77 (8%)	13 (1%)	15	51
2	B	245/279 (88%)	226 (92%)	13 (5%)	6 (2%)	7	34
All	All	1276/1375 (93%)	1167 (92%)	90 (7%)	19 (2%)	13	47

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	208	THR
1	A	377	PRO
1	A	871	GLN
1	A	946	GLY
2	B	437	VAL
2	B	440	ASP
1	A	2	PRO
1	A	966	LYS
2	B	439	GLU
2	B	595	ASN
1	A	264	LYS
1	A	298	PRO
1	A	948	LYS
1	A	376	VAL
1	A	956	LEU
1	A	38	ARG
2	B	393	SER
1	A	471	PRO
2	B	418	PRO

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	954/998 (96%)	917 (96%)	37 (4%)	39	75
2	B	238/259 (92%)	227 (95%)	11 (5%)	33	70
All	All	1192/1257 (95%)	1144 (96%)	48 (4%)	38	75

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	GLU
1	A	39	GLU
1	A	107	ASN
1	A	112	ILE
1	A	115	ARG
1	A	185	LEU
1	A	190	ILE
1	A	208	THR
1	A	215	CYS
1	A	239	LEU
1	A	326	SER
1	A	335	ARG
1	A	369	ASP
1	A	375	ARG
1	A	377	PRO
1	A	383	TRP
1	A	462	THR
1	A	464	SER
1	A	470	THR
1	A	476	GLU
1	A	478	ASP
1	A	490	SER
1	A	509	SER
1	A	523	LEU
1	A	526	ASN
1	A	542	GLU
1	A	545	GLU
1	A	575	ASN
1	A	639	LEU
1	A	679	THR
1	A	770	ARG
1	A	785	ASN
1	A	834	LEU

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Mol	Chain	Res	Type
1	A	891	ASP
1	A	966	LYS
1	A	971	CYS
1	A	1048	HIS
2	B	328	LEU
2	B	349	ASP
2	B	355	PHE
2	B	364	MET
2	B	415	GLN
2	B	425	LEU
2	B	482	THR
2	B	501	GLN
2	B	531	LEU
2	B	536	SER
2	B	594	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	-21	HIS
1	A	137	GLN
1	A	515	ASN
1	A	554	HIS
1	A	575	ASN
1	A	605	ASN
1	A	714	ASN
1	A	749	GLN
1	A	782	ASN
1	A	785	ASN
1	A	795	GLN
1	A	796	ASN
1	A	878	HIS
1	A	885	ASN
1	A	996	ASN
1	A	1047	HIS
2	B	417	ASN
2	B	517	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
1	SEP	A	790	1	7,9,10	0.74	0	8,12,14	1.20	1 (12%)

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
1	SEP	A	790	1	-	0/5/8/10	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	790	SEP	O3P-P-O2P	2.22	115.61	107.44

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
1	A	790	SEP	1	0

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

2 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	AX7	A	1101	-	11,11,11	1.33	1 (9%)	10,15,15	1.13	0
4	CAQ	B	701	-	8,8,8	0.59	0	10,10,10	0.33	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	AX7	A	1101	-	-	0/0/0/0	0/2/2/2
4	CAQ	B	701	-	-	0/0/0/0	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1101	AX7	CAJ-CAI	2.48	1.50	1.42

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

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	1041/1096 (94%)	0.21	31 (2%) 54 30	48, 87, 145, 199	0
2	B	255/279 (91%)	0.67	30 (11%) 6 2	80, 138, 178, 202	0
All	All	1296/1375 (94%)	0.30	61 (4%) 35 16	48, 94, 162, 202	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	947	TYR	6.1
2	B	335	TRP	4.5
2	B	510	GLU	4.0
1	A	872	PHE	3.8
2	B	584	LEU	3.7
1	A	341	ALA	3.6
2	B	409	ARG	3.3
2	B	425	LEU	3.3
1	A	-26	SER	3.3
2	B	437	VAL	3.1
2	B	388	GLY	3.1
1	A	374	GLN	3.1
1	A	419	HIS	3.1
1	A	353	LYS	3.0
1	A	408	SER	3.0
2	B	423	LYS	2.8
2	B	410	ASN	2.8
1	A	409	VAL	2.7
1	A	973	LYS	2.7
2	B	441	ASN	2.7
2	B	405	ILE	2.6
1	A	308	SER	2.6
2	B	328	LEU	2.6
1	A	1	MET	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	559	ILE	2.5
1	A	523	LEU	2.5
2	B	508	TYR	2.5
1	A	322	THR	2.4
2	B	580	TYR	2.4
1	A	1009	GLY	2.4
1	A	107	ASN	2.4
2	B	518	GLU	2.3
1	A	378	CYS	2.3
1	A	381	PRO	2.3
1	A	871	GLN	2.3
1	A	810	ASP	2.3
1	A	869	ALA	2.2
1	A	212	ASN	2.2
2	B	387	ASP	2.2
2	B	444	ALA	2.2
1	A	363	GLY	2.2
2	B	514	ARG	2.2
2	B	391	GLY	2.2
1	A	354	ILE	2.2
2	B	332	GLU	2.2
2	B	355	PHE	2.1
2	B	503	ARG	2.1
2	B	529	ASP	2.1
2	B	329	GLN	2.1
1	A	868	GLY	2.1
2	B	504	TYR	2.1
2	B	512	PHE	2.1
1	A	935	GLY	2.1
1	A	605	ASN	2.1
1	A	183	ASN	2.0
2	B	338	ILE	2.0
2	B	375	GLY	2.0
1	A	323	SER	2.0
1	A	888	GLU	2.0
1	A	352	ASP	2.0
2	B	505	SER	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	SEP	A	790	10/11	0.90	0.16	-	69,76,138,138	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	CAQ	B	701	8/8	0.87	0.88	6.23	122,140,154,156	0
3	AX7	A	1101	10/10	0.84	0.71	4.83	96,106,125,126	0

6.5 Other polymers [i](#)

There are no such residues in this entry.