Full wwPDB X-ray Structure Validation Report (i)

Sep 24, 2023 – 01:39 AM EDT

PDB ID	:	5SXA
Title	:	Crystal Structure of PI3Kalpha in complex with fragment 10
Authors	:	Gabelli, S.B.; Vogelstein, B.; Miller, M.S.; Amzel, L.M.
Deposited on	:	2016-08-09
Resolution	:	3.35 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 (i) 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 (i)) were used in the production of this report:

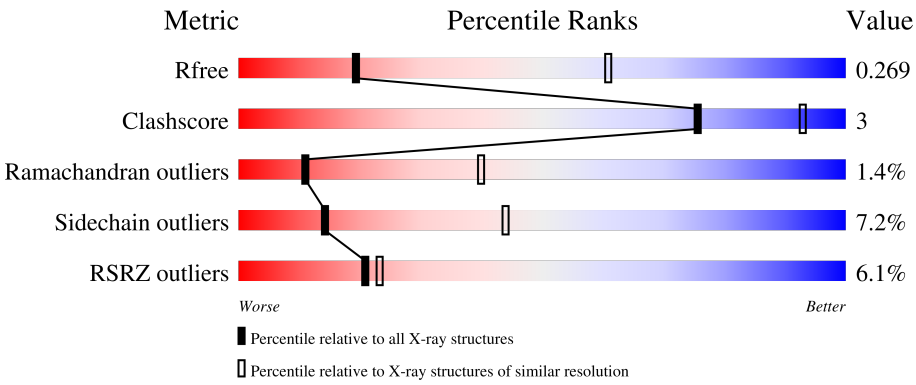
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.35 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	1558 (3.42-3.30)
Clashscore	141614	1627 (3.42-3.30)
Ramachandran outliers	138981	1599 (3.42-3.30)
Sidechain outliers	138945	1598 (3.42-3.30)
RSRZ outliers	127900	1507 (3.42-3.30)

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 <=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	3%, 82%, 12%, • •
2	B	279	16%, 65%, 13%, 22%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 10520 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	P	S			
1	A	1052	8629	5510	1474	1572	3	70	0	0	0

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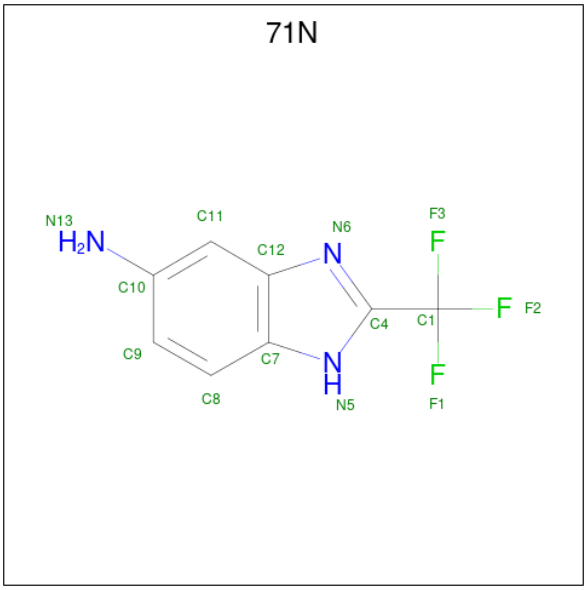
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P42336
A	0	SER	-	expression tag	UNP P42336

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	218	1877	1182	329	361	5	0	0	0

- Molecule 3 is 2-(trifluoromethyl)-1H-benzimidazol-5-amine (three-letter code: 71N) (formula: $C_8H_6F_3N_3$).



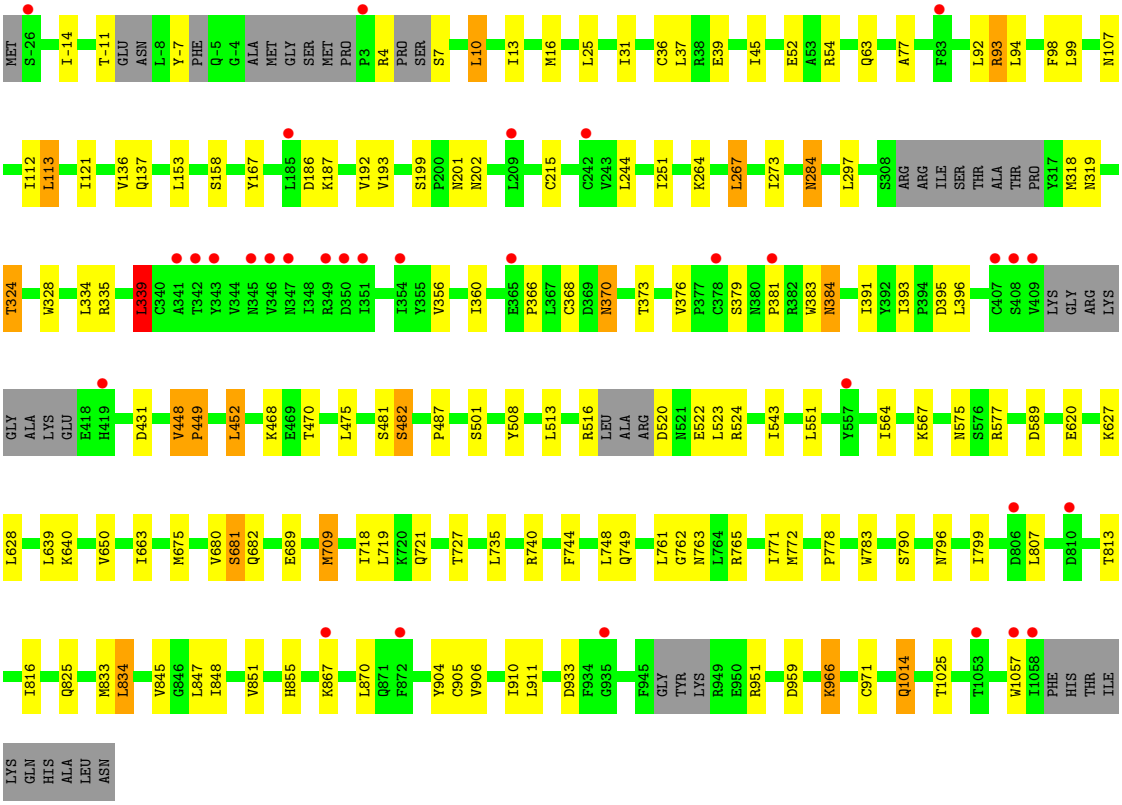
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	F	N		
3	A	1	14	8	3	3	0	0

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.

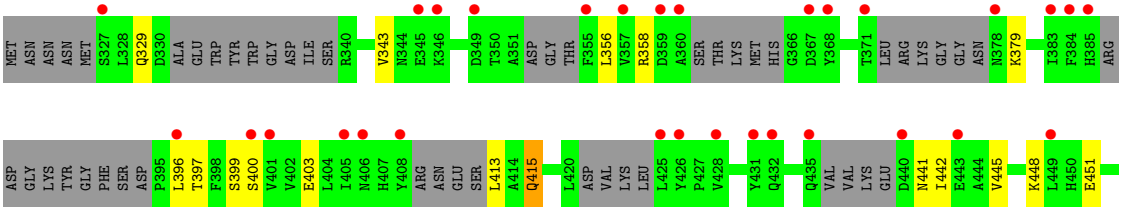
• Molecule 1: Phosphatidylinositol 4,5-bisphosphate 3-kinase catalytic subunit alpha isoform

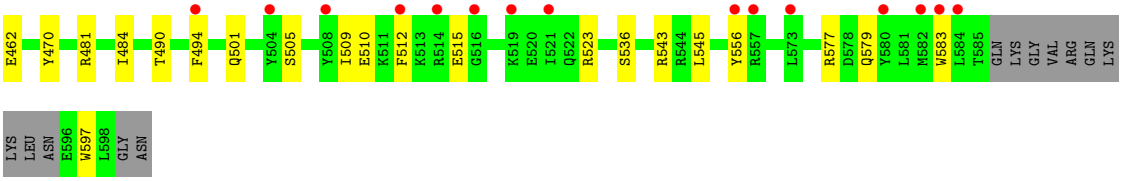
Chain A:



• Molecule 2: Phosphatidylinositol 3-kinase regulatory subunit alpha

Chain B:




E462
Y470
R481
I484
T490
F494
Q501
Y504
S505
Y508
I509
E510
K511
F512
K513
R514
E515
G516
K519
E520
I521
Q522
R523
S536
R543
R544
L545
Y556
R557
L573
R577
D578
Q579
Y580
L581
M582
W583
L584
T585
GLN
LYS
GLY
VAL
ARG
GLN
LYS
LYS
LEU
ASN
E596
W597
L598
GLY
ASN

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	115.22Å 117.37Å 152.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	92.91 – 3.35 49.45 – 3.34	Depositor EDS
% Data completeness (in resolution range)	99.9 (92.91-3.35) 99.9 (49.45-3.34)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.34 (at 3.33Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.222 , 0.275 0.220 , 0.269	Depositor DCC
R_{free} test set	1558 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å^2)	106.0	Xtriage
Anisotropy	0.379	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 87.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.013 for k,h,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	10520	wwPDB-VP
Average B, all atoms (Å^2)	136.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: 71N, 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	0.45	0/8794	0.68	2/11877 (0.0%)
2	B	0.46	0/1900	0.65	0/2533
All	All	0.45	0/10694	0.68	2/14410 (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	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	93	ARG	NE-CZ-NH1	8.45	124.53	120.30
1	A	339	LEU	CA-CB-CG	5.68	128.36	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	113	LEU	Peptide
1	A	448	VAL	Peptide

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	8629	0	8581	55	0
2	B	1877	0	1855	11	0
3	A	14	0	0	0	0
All	All	10520	0	10436	64	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 3.

All (64) 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:816:ILE:HG21	1:A:911:LEU:HD21	1.74	0.69
1:A:765:ARG:HD3	1:A:796:ASN:HD21	1.62	0.64
2:B:343:VAL:HG13	2:B:356:LEU:HD11	1.81	0.63
1:A:121:ILE:HD11	1:A:689:GLU:HA	1.82	0.62
1:A:910:ILE:O	1:A:1025:THR:HG21	2.07	0.54
1:A:10:LEU:HD13	1:A:16:MET:CE	2.37	0.54
1:A:639:LEU:HD22	1:A:650:VAL:HG22	1.90	0.53
1:A:543:ILE:HD11	1:A:567:LYS:HD3	1.91	0.53
1:A:334:LEU:HA	1:A:393:ILE:HD11	1.92	0.52
1:A:639:LEU:HD22	1:A:650:VAL:CG2	2.40	0.51
2:B:343:VAL:HG21	2:B:358:ARG:CD	2.41	0.51
1:A:370:ASN:N	1:A:370:ASN:HD22	2.08	0.51
1:A:376:VAL:HB	1:A:381:PRO:HA	1.94	0.50
1:A:799:ILE:CD1	1:A:847:LEU:HD22	2.41	0.49
1:A:136:VAL:HG21	1:A:682:GLN:HE22	1.77	0.49
1:A:628:LEU:HD23	1:A:663:ILE:HD13	1.94	0.49
1:A:1014:GLN:HE21	1:A:1014:GLN:HA	1.78	0.49
1:A:735:LEU:HD22	1:A:771:ILE:HG13	1.94	0.48
1:A:834:LEU:HD21	1:A:851:VAL:HG11	1.93	0.48
1:A:709:MET:CE	1:A:847:LEU:HD21	2.43	0.48
1:A:772:MET:HB2	1:A:778:PRO:HG2	1.96	0.48
1:A:284:ASN:HD22	1:A:284:ASN:N	2.12	0.47
2:B:501:GLN:O	2:B:505:SER:N	2.39	0.46
1:A:718:ILE:O	1:A:721:GLN:O	2.34	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:-11:THR:HB	1:A:-7:TYR:HB2	1.98	0.46
1:A:368:CYS:O	1:A:370:ASN:ND2	2.49	0.46
1:A:121:ILE:HD11	1:A:689:GLU:CA	2.45	0.45
1:A:719:LEU:HD13	1:A:771:ILE:HD13	1.97	0.45
1:A:376:VAL:HG12	1:A:379:SER:O	2.17	0.45
2:B:399:SER:HB2	2:B:403:GLU:HB3	1.97	0.45
1:A:675:MET:HE1	1:A:681:SER:O	2.17	0.45
1:A:833:MET:HE1	1:A:904:TYR:HA	1.99	0.45
1:A:324:THR:CG2	1:A:482:SER:HB3	2.47	0.45
1:A:749:GLN:O	1:A:762:GLY:O	2.35	0.45
1:A:356:VAL:HG23	1:A:383:TRP:CH2	2.52	0.45
1:A:25:LEU:HD21	2:B:494:PHE:CE1	2.53	0.44
1:A:448:VAL:N	1:A:449:PRO:HA	2.32	0.44
1:A:807:LEU:HD11	1:A:848:ILE:HD11	1.99	0.44
2:B:445:VAL:HG11	2:B:583:TRP:CE3	2.53	0.44
1:A:192:VAL:HG12	1:A:193:VAL:N	2.33	0.44
1:A:391:ILE:HD13	1:A:396:LEU:HD23	2.00	0.43
1:A:324:THR:HG21	1:A:482:SER:HB3	2.01	0.43
1:A:761:LEU:HD22	1:A:783:TRP:CE3	2.54	0.43
2:B:343:VAL:HG21	2:B:358:ARG:HD3	2.00	0.43
1:A:167:TYR:CE2	1:A:297:LEU:HD21	2.54	0.43
1:A:640:LYS:HE2	1:A:680:VAL:HG11	2.01	0.42
1:A:551:LEU:HD21	1:A:564:ILE:HD11	2.02	0.42
1:A:395:ASP:HA	1:A:577:ARG:HB2	2.00	0.42
1:A:749:GLN:HG2	1:A:763:ASN:HA	2.01	0.42
1:A:328:TRP:CE2	1:A:487:PRO:HG3	2.55	0.41
2:B:470:TYR:CE1	2:B:556:TYR:CE1	3.08	0.41
1:A:10:LEU:HD13	1:A:16:MET:HE2	2.02	0.41
1:A:383:TRP:O	1:A:384:ASN:ND2	2.52	0.41
1:A:136:VAL:HG21	1:A:682:GLN:NE2	2.34	0.41
1:A:98:PHE:CE2	2:B:490:THR:HG23	2.56	0.41
1:A:10:LEU:HB2	1:A:13:ILE:HD11	2.02	0.41
1:A:339:LEU:N	1:A:339:LEU:HD13	2.36	0.41
1:A:360:ILE:O	1:A:366:PRO:HD2	2.21	0.41
1:A:92:LEU:HB2	1:A:94:LEU:HG	2.03	0.40
1:A:267:LEU:HD22	1:A:273:ILE:HG13	2.02	0.40
1:A:906:VAL:HG12	1:A:910:ILE:HD12	2.03	0.40
2:B:413:LEU:HA	2:B:415:GLN:HE21	1.85	0.40
2:B:484:ILE:HG13	2:B:545:LEU:HD23	2.04	0.40
1:A:744:PHE:HD2	1:A:748:LEU:HD12	1.86	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	1032/1096 (94%)	923 (89%)	96 (9%)	13 (1%)	12 42
2	B	198/279 (71%)	177 (89%)	17 (9%)	4 (2%)	7 33
All	All	1230/1375 (90%)	1100 (89%)	113 (9%)	17 (1%)	11 40

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	202	ASN
1	A	452	LEU
2	B	400	SER
1	A	77	ALA
1	A	186	ASP
1	A	681	SER
1	A	264	LYS
1	A	522	GLU
1	A	966	LYS
2	B	442	ILE
2	B	515	GLU
1	A	201	ASN
1	A	481	SER
1	A	933	ASP
1	A	324	THR
2	B	512	PHE
1	A	449	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	961/996 (96%)	895 (93%)	66 (7%)	15	46
2	B	207/259 (80%)	189 (91%)	18 (9%)	10	35
All	All	1168/1255 (93%)	1084 (93%)	84 (7%)	14	44

All (84) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-14	ILE
1	A	4	ARG
1	A	10	LEU
1	A	31	ILE
1	A	36	CYS
1	A	37	LEU
1	A	39	GLU
1	A	45	ILE
1	A	52	GLU
1	A	54	ARG
1	A	63	GLN
1	A	93	ARG
1	A	99	LEU
1	A	107	ASN
1	A	112	ILE
1	A	113	LEU
1	A	137	GLN
1	A	153	LEU
1	A	187	LYS
1	A	199	SER
1	A	215	CYS
1	A	244	LEU
1	A	251	ILE
1	A	267	LEU
1	A	284	ASN
1	A	318	MET
1	A	319	ASN
1	A	335	ARG
1	A	339	LEU
1	A	370	ASN
1	A	373	THR
1	A	384	ASN
1	A	431	ASP
1	A	452	LEU

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Mol	Chain	Res	Type
1	A	468	LYS
1	A	470	THR
1	A	475	LEU
1	A	482	SER
1	A	501	SER
1	A	508	TYR
1	A	513	LEU
1	A	516	ARG
1	A	520	ASP
1	A	523	LEU
1	A	524	ARG
1	A	575	ASN
1	A	589	ASP
1	A	620	GLU
1	A	627	LYS
1	A	709	MET
1	A	727	THR
1	A	740	ARG
1	A	813	THR
1	A	825	GLN
1	A	834	LEU
1	A	845	VAL
1	A	855	HIS
1	A	867	LYS
1	A	870	LEU
1	A	905	CYS
1	A	951	ARG
1	A	959	ASP
1	A	966	LYS
1	A	971	CYS
1	A	1014	GLN
1	A	1057	TRP
2	B	329	GLN
2	B	379	LYS
2	B	396	LEU
2	B	397	THR
2	B	415	GLN
2	B	441	ASN
2	B	448	LYS
2	B	451	GLU
2	B	462	GLU
2	B	481	ARG

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Mol	Chain	Res	Type
2	B	509	ILE
2	B	510	GLU
2	B	523	ARG
2	B	536	SER
2	B	543	ARG
2	B	577	ARG
2	B	579	GLN
2	B	597	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	247	GLN
1	A	319	ASN
1	A	345	ASN
1	A	370	ASN
1	A	444	ASN
1	A	575	ASN
1	A	682	GLN
1	A	714	ASN
1	A	749	GLN
1	A	782	ASN
1	A	785	ASN
1	A	796	ASN
1	A	825	GLN
1	A	855	HIS
1	A	994	HIS
1	A	1014	GLN
1	A	1042	GLN
1	A	1047	HIS
2	B	415	GLN
2	B	457	GLN
2	B	564	ASN
2	B	579	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

3 non-standard protein/DNA/RNA residues 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	$\#\|Z\| > 2$
1	SEP	A	7	1	8,9,10	0.74	0	8,12,14	1.39	2 (25%)
1	SEP	A	790	1	8,9,10	0.66	0	8,12,14	2.12	2 (25%)
1	SEP	A	158	1	8,9,10	0.61	0	8,12,14	1.90	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	7	1	-	5/5/8/10	-
1	SEP	A	790	1	-	4/5/8/10	-
1	SEP	A	158	1	-	3/5/8/10	-

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	790	SEP	OG-CB-CA	5.17	113.18	108.14
1	A	158	SEP	OG-CB-CA	4.58	112.60	108.14
1	A	7	SEP	P-OG-CB	2.21	124.38	118.30
1	A	7	SEP	OG-CB-CA	2.13	110.22	108.14
1	A	790	SEP	P-OG-CB	2.06	123.97	118.30

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	7	SEP	N-CA-CB-OG

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Mol	Chain	Res	Type	Atoms
1	A	7	SEP	CB-OG-P-O2P
1	A	7	SEP	CB-OG-P-O3P
1	A	790	SEP	N-CA-CB-OG
1	A	790	SEP	CB-OG-P-O1P
1	A	790	SEP	CB-OG-P-O2P
1	A	790	SEP	CB-OG-P-O3P
1	A	7	SEP	CB-OG-P-O1P
1	A	158	SEP	CB-OG-P-O1P
1	A	158	SEP	N-CA-CB-OG
1	A	158	SEP	CB-OG-P-O2P
1	A	7	SEP	CA-CB-OG-P

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

1 ligand 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	$\#\|Z\| > 2$
3	71N	A	1101	-	13,15,15	1.16	1 (7%)	14,23,23	0.79	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	71N	A	1101	-	-	0/6/6/6	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1101	71N	C10-N13	3.04	1.48	1.38

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 (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	1049/1096 (95%)	0.22	32 (3%) 49 52	66, 117, 189, 248	0
2	B	218/279 (78%)	0.90	45 (20%) 1 1	110, 192, 249, 273	0
All	All	1267/1375 (92%)	0.34	77 (6%) 21 23	66, 125, 217, 273	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	385	HIS	5.7
1	A	347	ASN	5.5
1	A	346	VAL	5.3
1	A	-26	SER	5.1
2	B	432	GLN	4.7
2	B	514	ARG	4.3
2	B	368	TYR	3.8
1	A	354	ILE	3.8
2	B	383	ILE	3.7
2	B	405	ILE	3.6
1	A	378	CYS	3.6
2	B	435	GLN	3.6
2	B	349	ASP	3.5
1	A	381	PRO	3.5
2	B	367	ASP	3.4
2	B	583	TRP	3.4
2	B	371	THR	3.4
1	A	408	SER	3.4
2	B	360	ALA	3.4
1	A	345	ASN	3.3
2	B	378	ASN	3.3
1	A	3	PRO	3.3
2	B	521	ILE	3.3
1	A	872	PHE	3.3

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Mol	Chain	Res	Type	RSRZ
2	B	516	GLY	3.2
1	A	1053	THR	3.2
1	A	419	HIS	3.2
2	B	327	SER	3.2
2	B	428	VAL	3.1
1	A	350	ASP	3.1
2	B	425	LEU	3.1
2	B	440	ASP	3.0
2	B	512	PHE	3.0
2	B	384	PHE	3.0
2	B	426	TYR	3.0
1	A	409	VAL	2.9
2	B	519	LYS	2.9
2	B	408	TYR	2.9
2	B	508	TYR	2.9
1	A	935	GLY	2.9
1	A	1058	ILE	2.9
2	B	443	GLU	2.8
2	B	584	LEU	2.8
1	A	351	ILE	2.8
2	B	449	LEU	2.8
1	A	407	CYS	2.6
2	B	580	TYR	2.6
2	B	494	PHE	2.6
2	B	504	TYR	2.6
1	A	83	PHE	2.5
2	B	357	VAL	2.5
1	A	365	GLU	2.5
2	B	359	ASP	2.4
2	B	400	SER	2.4
2	B	573	LEU	2.4
2	B	345	GLU	2.4
2	B	406	ASN	2.3
2	B	401	VAL	2.3
1	A	342	THR	2.3
2	B	557	ARG	2.2
1	A	343	TYR	2.2
1	A	557	TYR	2.2
2	B	431	TYR	2.2
2	B	346	LYS	2.2
1	A	867	LYS	2.2
1	A	810	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	185	LEU	2.2
1	A	1057	TRP	2.1
1	A	209	LEU	2.1
1	A	349	ARG	2.1
2	B	396	LEU	2.1
2	B	556	TYR	2.1
2	B	582	MET	2.1
1	A	341	ALA	2.0
1	A	806	ASP	2.0
2	B	355	PHE	2.0
1	A	242	CYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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
1	SEP	A	7	10/11	0.63	0.33	154,178,227,235	0
1	SEP	A	158	10/11	0.92	0.15	133,165,229,235	0
1	SEP	A	790	10/11	0.93	0.16	110,126,204,217	0

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(Å^2)	Q<0.9
3	71N	A	1101	14/14	0.65	0.23	202,208,217,221	0

6.5 Other polymers (i)

There are no such residues in this entry.