



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:26 PM UTC

PDB ID : 3ZVV / pdb_00003zvv
Title : Fragment Bound to PI3Kinase gamma
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Deposited on : 2011-07-27
Resolution : 2.50 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

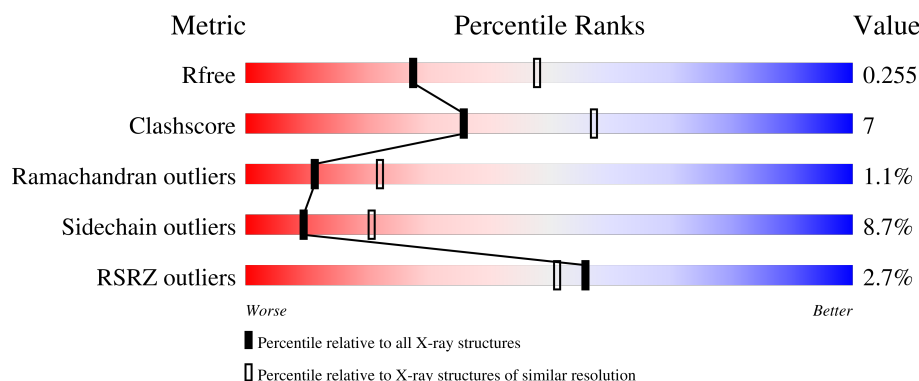
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

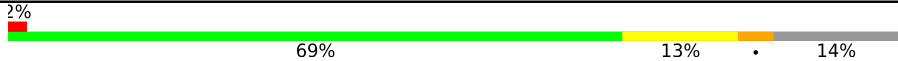
The reported resolution of this entry is 2.50 Å.

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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	966	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	XAZ	A	1500	-	X	-	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7023 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.

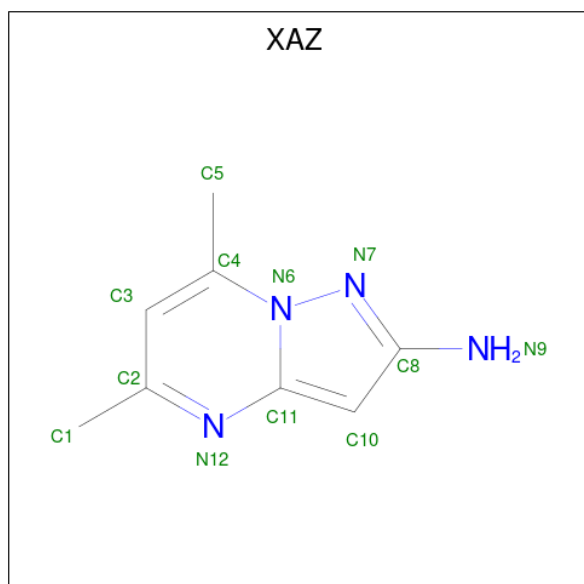
- Molecule 1 is a protein called PHOSPHATIDYLINOSITOL-4,5-BISPHOSPHATE 3-KINASE CATALYTIC SUBUNIT GAMMA ISOFORM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	829	Total	C	N	O	S	0	0	0
			6724	4319	1150	1220	35			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	143	MET	-	expression tag	UNP P48736
A	1103	HIS	-	expression tag	UNP P48736
A	1104	HIS	-	expression tag	UNP P48736
A	1105	HIS	-	expression tag	UNP P48736
A	1106	HIS	-	expression tag	UNP P48736
A	1107	HIS	-	expression tag	UNP P48736
A	1108	HIS	-	expression tag	UNP P48736

- Molecule 2 is 5,7-dimethylpyrazolo[1,5-a]pyrimidin-2-amine (CCD ID: XAZ) (formula: C₈H₁₀N₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			12	8	4		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	287	Total	O	0	0
			287	287		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	143.49Å 68.03Å 106.25Å 90.00° 95.16° 90.00°	Depositor
Resolution (Å)	16.62 – 2.50 16.62 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.1 (16.62-2.50) 98.8 (16.62-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 2.51Å)	Xtriage
Refinement program	BUSTER 2.9.6	Depositor
R, R_{free}	0.192 , 0.242 0.203 , 0.255	Depositor DCC
R_{free} test set	1755 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	53.4	Xtriage
Anisotropy	0.369	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 74.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7023	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.53% 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: XAZ

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	0.79	1/6868 (0.0%)	1.17	24/9288 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	804	MET	SD-CE	-7.21	1.61	1.79

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	377	THR	CA-C-N	7.08	131.24	121.33
1	A	377	THR	C-N-CA	7.08	131.24	121.33
1	A	966	GLY	CA-C-N	6.44	133.30	121.70
1	A	966	GLY	C-N-CA	6.44	133.30	121.70
1	A	964	ASP	CA-CB-CG	6.39	118.99	112.60
1	A	898	ASN	CA-C-N	6.27	133.52	121.54
1	A	898	ASN	C-N-CA	6.27	133.52	121.54
1	A	1041	GLN	N-CA-C	6.04	118.30	108.63
1	A	906	VAL	N-CA-C	5.85	117.28	110.62
1	A	948	HIS	CA-C-N	5.85	130.04	120.63
1	A	948	HIS	C-N-CA	5.85	130.04	120.63
1	A	1040	PRO	N-CA-C	5.80	124.42	112.47
1	A	895	THR	CA-C-N	5.70	132.23	121.97
1	A	895	THR	C-N-CA	5.70	132.23	121.97
1	A	373	LEU	N-CA-C	5.62	122.23	109.81
1	A	1040	PRO	CB-CA-C	5.52	120.67	111.56
1	A	1039	MET	O-C-N	5.51	126.37	121.36
1	A	191	ARG	N-CA-C	5.48	118.85	110.20
1	A	521	ASP	CA-CB-CG	5.46	118.06	112.60
1	A	202	VAL	N-CA-CB	-5.46	103.26	112.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	957	THR	CB-CA-C	5.45	119.81	110.01
1	A	904	ASP	N-CA-C	5.36	117.20	111.36
1	A	412	VAL	N-CA-C	5.04	115.84	108.53
1	A	498	ASN	N-CA-C	5.01	118.18	110.42

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	6724	0	6768	98	0
2	A	12	0	10	2	0
3	A	287	0	0	3	0
All	All	7023	0	6778	98	0

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

All (98) 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:804:MET:HE1	1:A:831:ILE:HG12	1.33	1.07
1:A:939:THR:HG23	1:A:945:GLY:HA2	1.49	0.94
1:A:225:HIS:HE1	1:A:304:HIS:HD2	1.14	0.93
1:A:689:LYS:HG2	1:A:728:MET:HE2	1.52	0.92
1:A:689:LYS:HG2	1:A:728:MET:CE	2.00	0.91
1:A:839:ARG:HA	1:A:842:MET:HE2	1.53	0.90
1:A:235:VAL:HG21	1:A:244:ILE:HG12	1.56	0.87
1:A:935:TYR:O	1:A:939:THR:HB	1.77	0.84
1:A:939:THR:HG23	1:A:945:GLY:CA	2.14	0.77
1:A:804:MET:CE	1:A:831:ILE:HG12	2.14	0.75
1:A:804:MET:HE1	1:A:831:ILE:CG1	2.16	0.74
1:A:949:ASN:H	1:A:1083:GLN:HE22	1.34	0.74
1:A:887:THR:HG22	1:A:890:LYS:H	1.53	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:HIS:HD2	3:A:2028:HOH:O	1.72	0.72
1:A:225:HIS:HE1	1:A:304:HIS:CD2	2.04	0.72
1:A:196:TYR:OH	1:A:728:MET:HE3	1.90	0.71
1:A:689:LYS:CG	1:A:728:MET:HE2	2.21	0.70
1:A:558:ILE:O	1:A:561:THR:HG22	1.92	0.69
1:A:271:VAL:CG2	1:A:282:VAL:HG11	2.22	0.68
1:A:214:LYS:HD3	1:A:297:LEU:O	1.97	0.65
1:A:391:GLN:NE2	1:A:633:CYS:HB2	2.12	0.65
1:A:1043:THR:C	1:A:1045:LYS:H	2.05	0.64
1:A:689:LYS:HG2	1:A:728:MET:HE1	1.78	0.63
1:A:614:ARG:HG2	1:A:617:TRP:HB3	1.82	0.62
1:A:1040:PRO:HB3	3:A:2112:HOH:O	1.98	0.61
1:A:804:MET:HE2	1:A:812:TRP:HB2	1.83	0.60
1:A:939:THR:CG2	1:A:945:GLY:HA2	2.25	0.60
1:A:561:THR:HG21	1:A:565:ASN:HD22	1.67	0.60
1:A:271:VAL:CG2	1:A:282:VAL:CG1	2.81	0.59
1:A:840:GLN:HG2	1:A:1039:MET:HE2	1.85	0.59
1:A:240:THR:HG23	1:A:243:ALA:H	1.67	0.59
1:A:981:GLU:HA	1:A:982:ARG:CZ	2.36	0.56
1:A:689:LYS:CD	1:A:728:MET:HE2	2.37	0.55
1:A:381:VAL:CG2	1:A:433:ILE:HG23	2.38	0.54
1:A:882:VAL:HG23	2:A:1500:XAZ:C5	2.37	0.54
1:A:768:LYS:O	1:A:772:GLU:HG2	2.07	0.54
1:A:460:LEU:HD21	1:A:501:LYS:HZ3	1.73	0.54
1:A:593:PHE:CZ	1:A:611:LEU:HD21	2.42	0.53
1:A:882:VAL:HG23	2:A:1500:XAZ:H53C	1.90	0.53
1:A:246:GLN:HE21	1:A:246:GLN:HA	1.74	0.53
1:A:235:VAL:CG2	1:A:244:ILE:HG12	2.35	0.52
1:A:905:GLU:HG3	1:A:993:PHE:CE2	2.45	0.52
1:A:903:LYS:HB3	1:A:906:VAL:HG23	1.91	0.52
1:A:271:VAL:HG22	1:A:282:VAL:CG1	2.38	0.52
1:A:1044:SER:HA	1:A:1048:ILE:HD12	1.93	0.51
1:A:651:LEU:HD22	1:A:655:ASP:HB3	1.92	0.51
1:A:168:VAL:HG13	1:A:170:ASP:O	2.10	0.51
1:A:225:HIS:CE1	1:A:304:HIS:HD2	2.07	0.51
1:A:181:VAL:HG12	1:A:185:MET:HE2	1.91	0.51
1:A:949:ASN:H	1:A:1083:GLN:NE2	2.07	0.50
1:A:549:ASN:O	1:A:553:LYS:HG3	2.11	0.49
1:A:196:TYR:OH	1:A:728:MET:CE	2.61	0.48
1:A:1042:LEU:CD1	1:A:1042:LEU:H	2.26	0.48
1:A:317:GLU:O	1:A:726:THR:CG2	2.62	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:748:ASP:HB3	1:A:770:LYS:NZ	2.29	0.48
1:A:749:ILE:CD1	1:A:770:LYS:HD2	2.44	0.47
1:A:1002:THR:HG23	1:A:1006:PHE:HD2	1.79	0.47
1:A:432:GLN:OE1	1:A:501:LYS:NZ	2.39	0.47
1:A:583:LEU:HD22	1:A:610:LEU:HD22	1.96	0.47
1:A:476:ARG:HB3	1:A:520:LEU:HD23	1.96	0.47
1:A:239:ASP:O	1:A:287:ILE:HG13	2.13	0.47
1:A:184:ARG:HD3	1:A:719:ALA:O	2.14	0.47
1:A:547:MET:HG2	1:A:578:PHE:CD1	2.50	0.47
1:A:1042:LEU:H	1:A:1042:LEU:HD13	1.80	0.46
1:A:317:GLU:O	1:A:726:THR:HG23	2.16	0.46
1:A:735:GLN:O	1:A:739:ILE:HG23	2.16	0.46
1:A:1036:MET:HA	1:A:1042:LEU:HD11	1.98	0.45
1:A:410:TRP:HB3	1:A:412:VAL:HG22	1.98	0.45
1:A:366:ARG:NH1	1:A:479:GLU:OE1	2.49	0.45
1:A:391:GLN:HE22	1:A:502:LEU:HD11	1.82	0.45
1:A:887:THR:HG21	1:A:950:ASP:OD1	2.18	0.44
1:A:379:LEU:HD23	1:A:380:THR:HG22	1.99	0.44
1:A:406:GLU:CD	1:A:406:GLU:H	2.26	0.44
1:A:373:LEU:O	1:A:374:PRO:C	2.61	0.44
1:A:749:ILE:HD12	1:A:770:LYS:HD2	2.00	0.44
1:A:1040:PRO:O	1:A:1042:LEU:HD12	2.18	0.43
1:A:233:ILE:HD12	1:A:248:PHE:HD1	1.83	0.43
1:A:363:VAL:HG23	1:A:520:LEU:CD1	2.48	0.43
1:A:476:ARG:HG3	1:A:480:TYR:OH	2.19	0.43
1:A:1043:THR:C	1:A:1045:LYS:N	2.73	0.43
1:A:593:PHE:HZ	1:A:611:LEU:HD21	1.83	0.43
1:A:739:ILE:HG13	1:A:740:GLU:N	2.33	0.43
1:A:236:SER:O	1:A:287:ILE:HD11	2.19	0.42
1:A:898:ASN:HB2	1:A:901:ALA:HB3	2.01	0.42
1:A:271:VAL:HG21	1:A:282:VAL:HG11	2.00	0.42
1:A:366:ARG:HH12	1:A:479:GLU:CD	2.27	0.42
1:A:391:GLN:HE22	1:A:633:CYS:HB2	1.84	0.42
1:A:578:PHE:HA	3:A:2135:HOH:O	2.20	0.42
1:A:363:VAL:HG12	1:A:416:PHE:HE1	1.85	0.41
1:A:903:LYS:NZ	1:A:905:GLU:HB3	2.36	0.41
1:A:1043:THR:O	1:A:1045:LYS:N	2.54	0.41
1:A:983:VAL:HG13	1:A:985:PHE:O	2.20	0.41
1:A:198:MET:SD	1:A:271:VAL:HG21	2.61	0.40
1:A:853:SER:O	1:A:857:THR:HG23	2.21	0.40
1:A:1086:TRP:CD1	1:A:1090:LEU:HG	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1042:LEU:CD1	1:A:1042:LEU:N	2.85	0.40
1:A:1041:GLN:HE21	1:A:1041:GLN:HA	1.87	0.40
1:A:547:MET:HE3	1:A:578:PHE:CD2	2.57	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	811/966 (84%)	777 (96%)	25 (3%)	9 (1%)	11	22

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	216	ALA
1	A	374	PRO
1	A	899	THR
1	A	1040	PRO
1	A	896	VAL
1	A	1044	SER
1	A	999	GLY
1	A	227	SER
1	A	966	GLY

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	744/864 (86%)	679 (91%)	65 (9%)	9 21

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	152	ARG
1	A	165	VAL
1	A	168	VAL
1	A	202	VAL
1	A	207	LEU
1	A	214	LYS
1	A	235	VAL
1	A	254	LYS
1	A	271	VAL
1	A	281	LEU
1	A	287	ILE
1	A	320	LYS
1	A	321	GLU
1	A	357	CYS
1	A	366	ARG
1	A	369	ASP
1	A	370	ILE
1	A	373	LEU
1	A	377	THR
1	A	379	LEU
1	A	391	GLN
1	A	392	GLN
1	A	406	GLU
1	A	421	LYS
1	A	435	CYS
1	A	461	LEU
1	A	476	ARG
1	A	498	ASN
1	A	520	LEU
1	A	527	ILE
1	A	574	LEU
1	A	601	GLN
1	A	610	LEU
1	A	613	ARG
1	A	615	GLU
1	A	682	LEU
1	A	707	ARG
1	A	726	THR

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Mol	Chain	Res	Type
1	A	728	MET
1	A	739	ILE
1	A	749	ILE
1	A	766	GLN
1	A	767	LEU
1	A	776	ASN
1	A	799	GLU
1	A	807	LYS
1	A	823	LEU
1	A	865	LEU
1	A	887	THR
1	A	893	GLN
1	A	895	THR
1	A	903	LYS
1	A	907	LEU
1	A	939	THR
1	A	957	THR
1	A	982	ARG
1	A	983	VAL
1	A	1001	LYS
1	A	1011	ASP
1	A	1039	MET
1	A	1040	PRO
1	A	1041	GLN
1	A	1042	LEU
1	A	1052	ARG
1	A	1090	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	225	HIS
1	A	246	GLN
1	A	291	GLN
1	A	304	HIS
1	A	386	ASN
1	A	389	HIS
1	A	391	GLN
1	A	392	GLN
1	A	396	GLN
1	A	430	ASN
1	A	565	ASN

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Mol	Chain	Res	Type
1	A	601	GLN
1	A	629	GLN
1	A	634	ASN
1	A	639	ASN
1	A	658	HIS
1	A	743	GLN
1	A	834	HIS
1	A	1041	GLN
1	A	1071	GLN
1	A	1083	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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
2	XAZ	A	1500	-	12,13,13	3.26	8 (66%)	15,19,19	5.71	10 (66%)

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
2	XAZ	A	1500	-	-	-	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1500	XAZ	C2-N12	5.55	1.41	1.33
2	A	1500	XAZ	C5-C4	-4.89	1.41	1.49
2	A	1500	XAZ	C11-N12	4.16	1.44	1.36
2	A	1500	XAZ	C1-C2	4.02	1.56	1.50
2	A	1500	XAZ	C11-N6	3.46	1.44	1.39
2	A	1500	XAZ	C8-N9	3.41	1.42	1.33
2	A	1500	XAZ	N6-N7	2.56	1.42	1.38
2	A	1500	XAZ	C10-C8	2.27	1.44	1.41

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1500	XAZ	C11-N12-C2	11.60	124.07	117.35
2	A	1500	XAZ	C1-C2-N12	8.56	126.39	116.84
2	A	1500	XAZ	C3-C4-N6	8.10	125.05	115.03
2	A	1500	XAZ	C3-C2-N12	-7.49	115.34	122.71
2	A	1500	XAZ	N6-C11-N12	-7.06	116.38	121.49
2	A	1500	XAZ	C4-N6-N7	6.40	135.65	125.74
2	A	1500	XAZ	C8-N7-N6	5.33	111.06	104.12
2	A	1500	XAZ	C11-N6-N7	-4.26	105.89	111.96
2	A	1500	XAZ	C10-C11-N12	3.05	137.43	132.64
2	A	1500	XAZ	C10-C8-N9	2.94	130.82	127.11

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1500	XAZ	2	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	829/966 (85%)	0.02	22 (2%)	56 51	34, 76, 138, 196	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1044	SER	7.5
1	A	1092	LEU	4.0
1	A	528	ALA	3.8
1	A	374	PRO	3.8
1	A	1040	PRO	3.2
1	A	357	CYS	3.1
1	A	1086	TRP	3.1
1	A	1091	VAL	3.0
1	A	896	VAL	2.8
1	A	216	ALA	2.8
1	A	1087	PHE	2.8
1	A	379	LEU	2.7
1	A	1090	LEU	2.5
1	A	435	CYS	2.4
1	A	144	SER	2.3
1	A	525	HIS	2.2
1	A	1089	HIS	2.2
1	A	308	ASP	2.2
1	A	1002	THR	2.1
1	A	395	CYS	2.1
1	A	1066	LYS	2.0
1	A	997	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	XAZ	A	1500	12/12	0.96	0.06	51,59,60,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.