



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:19 AM UTC

PDB ID : 3IBE / pdb_00003ibe
Title : Crystal Structure of a Pyrazolopyrimidine Inhibitor Bound to PI3 Kinase Gamma
Authors : Bard, J.; Svenson, K.
Deposited on : 2009-07-15
Resolution : 2.80 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

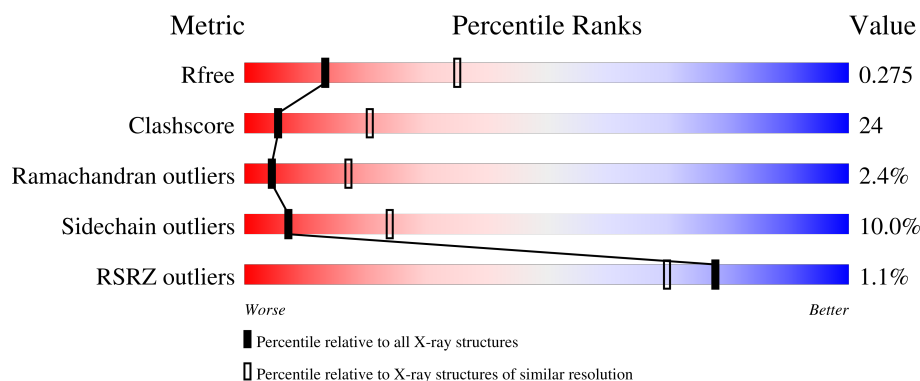
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6622 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.

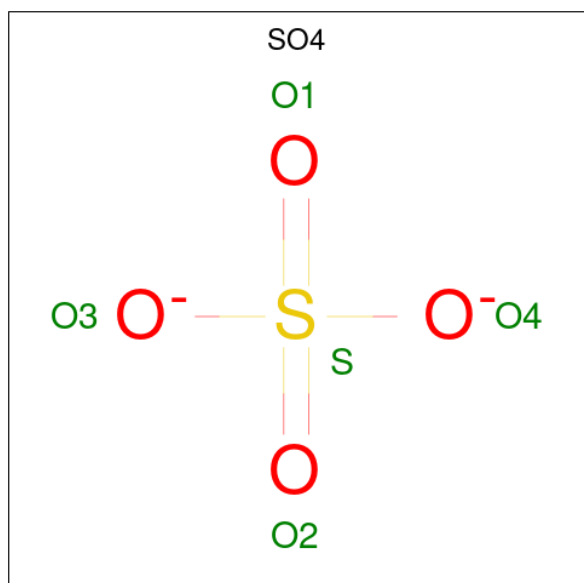
- Molecule 1 is a protein called Phosphatidylinositol-4,5-bisphosphate 3-kinase catalytic subunit gamma isoform.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	850	Total	C	N	O	S	0	0	0
			6530	4195	1102	1199	34			

There are 8 discrepancies between the modelled and reference sequences:

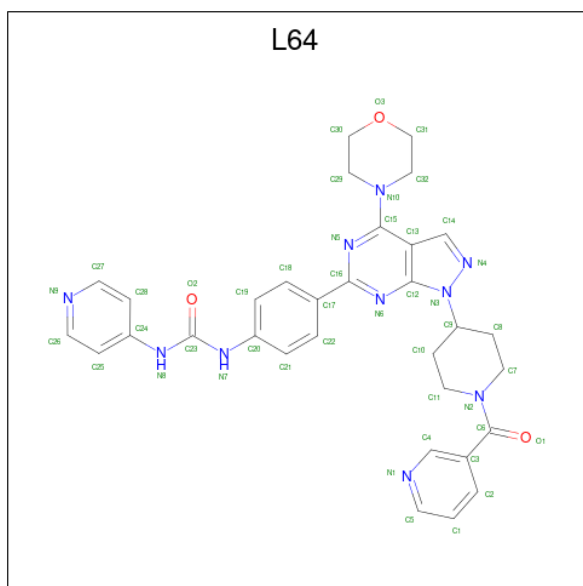
Chain	Residue	Modelled	Actual	Comment	Reference
A	143	MET	-	expression tag	UNP P48736
A	459	ARG	GLN	variant	UNP P48736
A	1103	HIS	-	expression tag	UNP P48736
A	1104	HIS	-	expression tag	UNP P48736
A	1105	HIS	-	expression tag	UNP P48736
A	1106	HIS	-	expression tag	UNP P48736
A	1107	HIS	-	expression tag	UNP P48736
A	1108	HIS	-	expression tag	UNP P48736

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 1-(4-{4-morpholin-4-yl-1-[1-(pyridin-3-ylcarbonyl)piperidin-4-yl]-1H-pyrazolo[3,4-d]pyrimidin-6-yl}phenyl)-3-pyridin-4-ylurea (CCD ID: L64) (formula: C₃₂H₃₂N₁₀O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			45	32	10	3		

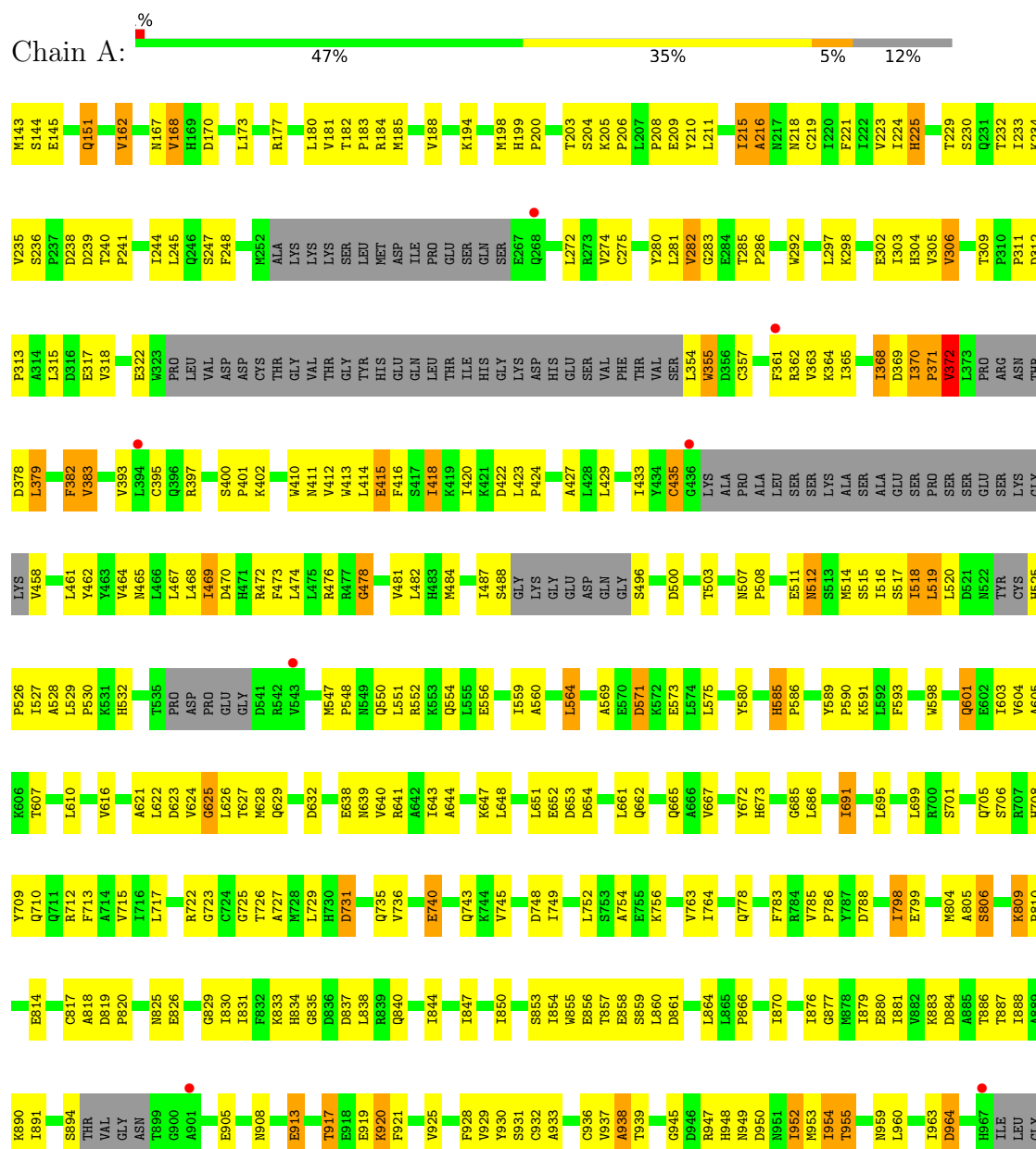
- Molecule 4 is water.

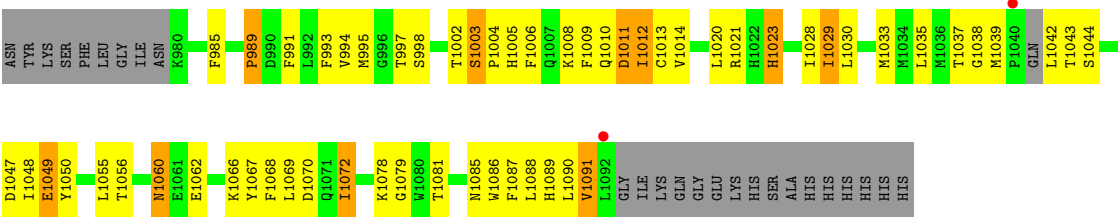
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	42	Total	O	0	0
			42	42		

3 Residue-property plots

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 gamma isoform





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	145.16Å 68.46Å 106.90Å 90.00° 94.80° 90.00°	Depositor
Resolution (Å)	44.71 – 2.80 44.71 – 2.80	Depositor EDS
% Data completeness (in resolution range)	83.1 (44.71-2.80) 83.1 (44.71-2.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 2.81Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.210 , 0.277 0.204 , 0.275	Depositor DCC
R_{free} test set	1117 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å ²)	69.4	Xtriage
Anisotropy	0.181	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 58.6	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	6622	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.94% 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: L64, SO4

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	0.73	0/6672	1.02	25/9081 (0.3%)

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	478	GLY	N-CA-C	12.41	124.10	111.56
1	A	809	LYS	CA-C-N	7.96	129.34	120.66
1	A	809	LYS	C-N-CA	7.96	129.34	120.66
1	A	603	ILE	CB-CA-C	-6.87	103.23	111.81
1	A	418	ILE	N-CA-C	6.75	117.36	107.37
1	A	778	GLN	N-CA-C	6.67	118.21	111.07
1	A	788	ASP	CA-C-N	6.65	126.78	119.87
1	A	788	ASP	C-N-CA	6.65	126.78	119.87
1	A	695	LEU	N-CA-C	-6.64	103.97	111.14
1	A	1003	SER	CA-C-N	5.89	125.57	119.56
1	A	1003	SER	C-N-CA	5.89	125.57	119.56
1	A	840	GLN	N-CA-C	-5.89	106.22	113.41
1	A	1012	ILE	CB-CA-C	-5.88	104.44	111.97
1	A	856	GLU	N-CA-C	-5.63	106.55	113.41
1	A	616	VAL	CB-CA-C	-5.48	104.89	112.24
1	A	312	ASP	CA-C-N	5.47	125.14	119.56
1	A	312	ASP	C-N-CA	5.47	125.14	119.56
1	A	644	ALA	N-CA-C	-5.42	105.45	111.36
1	A	571	ASP	N-CA-C	-5.39	105.41	111.28
1	A	844	ILE	CB-CA-C	-5.38	105.08	111.97
1	A	151	GLN	N-CA-C	-5.34	106.03	112.54
1	A	297	LEU	N-CA-C	5.27	116.71	110.97
1	A	743	GLN	N-CA-C	-5.21	105.29	110.97
1	A	625	GLY	N-CA-C	-5.10	106.60	112.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	890	LYS	N-CA-C	-5.04	106.40	112.54

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	6530	0	6206	301	0
2	A	5	0	0	0	0
3	A	45	0	32	9	0
4	A	42	0	0	1	0
All	All	6622	0	6238	303	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 24.

All (303) 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:180:LEU:O	1:A:183:PRO:HD2	1.83	0.79
1:A:953:MET:O	1:A:960:LEU:HD12	1.85	0.76
1:A:429:LEU:HB2	1:A:468:LEU:HD21	1.70	0.72
1:A:274:VAL:HG21	1:A:292:TRP:CD1	2.26	0.70
1:A:181:VAL:O	1:A:185:MET:HG3	1.94	0.68
1:A:627:THR:HG21	1:A:648:LEU:HD21	1.76	0.68
1:A:685:GLY:HA2	1:A:691:ILE:HG22	1.77	0.67
1:A:219:CYS:SG	1:A:234:LYS:HG2	2.35	0.66
1:A:397:ARG:HD2	1:A:415:GLU:O	1.95	0.66
1:A:731:ASP:O	1:A:735:GLN:HG3	1.95	0.66
1:A:936:CYS:SG	1:A:985:PHE:CD1	2.89	0.65
1:A:925:VAL:O	1:A:929:VAL:HG23	1.96	0.65
1:A:1013:CYS:HB3	1:A:1068:PHE:CE2	2.31	0.65
1:A:564:LEU:HD21	1:A:1048:ILE:HG22	1.77	0.65
1:A:514:MET:HG3	1:A:515:SER:N	2.11	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:833:LYS:O	1:A:876:ILE:HG13	1.96	0.65
1:A:370:ILE:HG12	1:A:371:PRO:O	1.97	0.65
1:A:378:ASP:O	1:A:379:LEU:HG	1.97	0.64
1:A:571:ASP:O	1:A:575:LEU:HD13	1.97	0.64
1:A:461:LEU:C	1:A:462:TYR:CD2	2.75	0.64
1:A:355:TRP:CE3	1:A:530:PRO:HG2	2.33	0.64
1:A:467:LEU:HD13	1:A:672:TYR:CE2	2.33	0.63
1:A:627:THR:HG21	1:A:648:LEU:CD2	2.29	0.63
1:A:435:CYS:O	1:A:458:VAL:HA	1.98	0.63
1:A:933:ALA:O	1:A:937:VAL:HG23	1.98	0.62
1:A:143:MET:C	1:A:145:GLU:H	2.07	0.62
1:A:354:LEU:O	1:A:527:ILE:HD11	1.99	0.62
1:A:507:ASN:OD1	1:A:508:PRO:CD	2.48	0.62
1:A:224:ILE:O	1:A:230:SER:HA	1.99	0.62
1:A:888:ILE:O	1:A:891:ILE:HB	2.00	0.61
1:A:709:TYR:HB3	1:A:713:PHE:CE1	2.35	0.61
1:A:357:CYS:HB2	1:A:527:ILE:CD1	2.30	0.61
1:A:413:TRP:CH2	1:A:519:LEU:HD23	2.35	0.61
1:A:640:VAL:O	1:A:643:ILE:HG12	2.00	0.61
1:A:1048:ILE:C	1:A:1050:TYR:H	2.07	0.61
1:A:921:PHE:O	1:A:925:VAL:HG23	2.01	0.61
1:A:1048:ILE:C	1:A:1050:TYR:N	2.55	0.61
1:A:1081:THR:O	1:A:1085:ASN:OD1	2.18	0.61
1:A:629:GLN:HG2	1:A:1029:ILE:HG13	1.83	0.60
1:A:173:LEU:HD23	1:A:673:HIS:CD2	2.37	0.60
1:A:464:VAL:HG22	1:A:465:ASN:H	1.65	0.60
1:A:1003:SER:HB2	1:A:1004:PRO:CD	2.31	0.60
1:A:955:THR:OG1	1:A:959:ASN:HB3	2.02	0.60
1:A:469:ILE:N	1:A:469:ILE:HD12	2.17	0.59
1:A:930:TYR:CE2	1:A:1012:ILE:HD13	2.37	0.59
1:A:500:ASP:HB3	1:A:708:HIS:CD2	2.36	0.59
1:A:726:THR:HA	1:A:729:LEU:HD12	1.84	0.59
1:A:952:ILE:HG22	1:A:953:MET:N	2.17	0.59
1:A:198:MET:HE2	1:A:311:PRO:HD2	1.85	0.59
1:A:662:GLN:HG2	1:A:1030:LEU:CD2	2.32	0.58
1:A:548:PRO:HG2	1:A:551:LEU:HD12	1.84	0.58
1:A:953:MET:SD	1:A:963:ILE:HD13	2.43	0.58
1:A:302:GLU:O	1:A:304:HIS:CD2	2.55	0.58
1:A:936:CYS:HB3	1:A:985:PHE:CE1	2.39	0.58
1:A:952:ILE:CG2	1:A:960:LEU:HD11	2.33	0.58
1:A:424:PRO:HD2	1:A:427:ALA:HB2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1069:LEU:HA	1:A:1072:ILE:HG13	1.85	0.58
1:A:272:LEU:HB3	1:A:305:VAL:HG11	1.85	0.57
1:A:1006:PHE:CE2	1:A:1010:GLN:CD	2.82	0.57
1:A:424:PRO:HG3	1:A:598:TRP:O	2.04	0.57
1:A:180:LEU:C	1:A:183:PRO:HD2	2.30	0.57
1:A:804:MET:HE2	3:A:1109:L64:H8	1.85	0.57
1:A:368:ILE:HD13	1:A:514:MET:HE1	1.87	0.57
1:A:1044:SER:HB3	1:A:1047:ASP:OD2	2.04	0.57
1:A:215:ILE:O	1:A:216:ALA:HB3	2.05	0.56
1:A:585:HIS:ND1	1:A:585:HIS:N	2.53	0.56
1:A:282:VAL:CG2	1:A:283:GLY:N	2.69	0.56
1:A:928:PHE:HZ	1:A:991:PHE:CD1	2.22	0.56
1:A:275:CYS:SG	1:A:306:VAL:HG12	2.46	0.56
1:A:928:PHE:O	1:A:932:CYS:HB3	2.06	0.56
1:A:357:CYS:HB2	1:A:527:ILE:HD12	1.87	0.56
1:A:624:VAL:O	1:A:628:MET:HG2	2.05	0.55
1:A:745:VAL:O	1:A:749:ILE:HG12	2.05	0.55
1:A:1035:LEU:O	1:A:1042:LEU:CB	2.53	0.55
1:A:225:HIS:N	1:A:225:HIS:CD2	2.74	0.55
1:A:236:SER:HB3	1:A:239:ASP:OD1	2.07	0.55
1:A:625:GLY:O	1:A:629:GLN:HG3	2.06	0.55
1:A:422:ASP:OD1	1:A:601:GLN:HB3	2.07	0.55
1:A:1089:HIS:C	1:A:1091:VAL:H	2.15	0.55
1:A:1003:SER:HB2	1:A:1004:PRO:HD2	1.88	0.55
1:A:413:TRP:CZ2	1:A:519:LEU:HD23	2.41	0.55
1:A:831:ILE:HG22	1:A:831:ILE:O	2.05	0.54
1:A:1044:SER:HB3	1:A:1047:ASP:CG	2.32	0.54
1:A:952:ILE:HG21	1:A:960:LEU:HD11	1.90	0.54
1:A:685:GLY:HA2	1:A:691:ILE:CG2	2.38	0.54
1:A:831:ILE:HG13	1:A:881:ILE:HG13	1.90	0.54
1:A:853:SER:O	1:A:857:THR:HG23	2.07	0.54
1:A:804:MET:HE2	3:A:1109:L64:C8	2.37	0.54
1:A:913:GLU:O	1:A:913:GLU:HG2	2.08	0.54
1:A:1006:PHE:C	1:A:1006:PHE:CD2	2.85	0.54
1:A:280:TYR:HB3	1:A:282:VAL:HG12	1.90	0.54
1:A:371:PRO:O	1:A:372:VAL:C	2.50	0.53
1:A:1048:ILE:O	1:A:1050:TYR:N	2.40	0.53
1:A:547:MET:HE2	1:A:552:ARG:N	2.22	0.53
1:A:887:THR:HA	1:A:953:MET:HG2	1.89	0.53
1:A:317:GLU:HG2	1:A:318:VAL:N	2.24	0.53
1:A:754:ALA:C	1:A:756:LYS:N	2.67	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:837:ASP:O	3:A:1109:L64:C28	2.57	0.53
1:A:478:GLY:H	1:A:520:LEU:HB2	1.75	0.52
1:A:482:LEU:HB2	1:A:516:ILE:HG22	1.92	0.52
1:A:837:ASP:O	3:A:1109:L64:C27	2.57	0.52
1:A:838:LEU:HD12	1:A:877:GLY:CA	2.39	0.52
1:A:224:ILE:C	1:A:225:HIS:CD2	2.88	0.52
1:A:208:PRO:HD2	1:A:211:LEU:HB2	1.90	0.52
1:A:1078:LYS:O	1:A:1079:GLY:C	2.53	0.52
1:A:948:HIS:CD2	1:A:1086:TRP:HB2	2.45	0.52
1:A:1035:LEU:HB3	1:A:1042:LEU:HA	1.90	0.52
1:A:507:ASN:OD1	1:A:508:PRO:HD2	2.10	0.52
1:A:151:GLN:OE1	1:A:722:ARG:NH2	2.43	0.51
1:A:547:MET:HE2	1:A:552:ARG:CA	2.41	0.51
1:A:274:VAL:HG21	1:A:292:TRP:NE1	2.26	0.51
1:A:496:SER:HB3	1:A:1043:THR:HB	1.93	0.51
1:A:879:ILE:HG22	1:A:880:GLU:N	2.25	0.51
1:A:361:PHE:HD2	1:A:416:PHE:HD1	1.58	0.51
1:A:686:LEU:CD2	1:A:723:GLY:C	2.83	0.51
1:A:876:ILE:O	1:A:876:ILE:HG23	2.10	0.51
1:A:365:ILE:HG22	1:A:365:ILE:O	2.09	0.50
1:A:368:ILE:HD12	1:A:368:ILE:C	2.36	0.50
1:A:1009:PHE:C	1:A:1009:PHE:CD2	2.89	0.50
1:A:274:VAL:CG2	1:A:292:TRP:CD1	2.94	0.50
1:A:838:LEU:HD12	1:A:877:GLY:HA3	1.93	0.50
1:A:559:ILE:O	1:A:559:ILE:HG22	2.12	0.50
1:A:580:TYR:CE2	1:A:610:LEU:CD1	2.94	0.50
1:A:315:LEU:O	1:A:727:ALA:HB2	2.10	0.50
1:A:205:LYS:HB3	1:A:206:PRO:HD2	1.93	0.50
1:A:825:ASN:O	1:A:826:GLU:C	2.53	0.50
1:A:182:THR:HB	1:A:183:PRO:HD3	1.93	0.49
1:A:469:ILE:N	1:A:469:ILE:CD1	2.74	0.49
1:A:464:VAL:HG12	1:A:484:MET:HE3	1.94	0.49
1:A:302:GLU:O	1:A:304:HIS:N	2.46	0.49
1:A:238:ASP:HA	1:A:286:PRO:HB3	1.93	0.49
1:A:954:ILE:HA	1:A:959:ASN:O	2.12	0.49
1:A:667:VAL:HG12	1:A:667:VAL:O	2.11	0.49
1:A:809:LYS:N	1:A:810:PRO:HD3	2.28	0.49
1:A:736:VAL:O	1:A:740:GLU:HB2	2.13	0.49
1:A:167:ASN:C	1:A:168:VAL:CG1	2.85	0.49
1:A:936:CYS:SG	1:A:985:PHE:CE1	3.05	0.49
1:A:952:ILE:CD1	1:A:952:ILE:N	2.75	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1056:THR:HG23	1:A:1056:THR:O	2.11	0.49
1:A:857:THR:C	1:A:858:GLU:HG2	2.38	0.48
1:A:1013:CYS:HB3	1:A:1068:PHE:HE2	1.76	0.48
1:A:1035:LEU:CD1	1:A:1048:ILE:HG12	2.42	0.48
1:A:854:ILE:O	1:A:854:ILE:HG22	2.12	0.48
1:A:1005:HIS:O	1:A:1008:LYS:HB3	2.12	0.48
1:A:368:ILE:HD13	1:A:514:MET:SD	2.53	0.48
1:A:948:HIS:ND1	1:A:950:ASP:HB2	2.28	0.48
1:A:363:VAL:O	1:A:363:VAL:HG13	2.14	0.48
1:A:465:ASN:HB2	1:A:503:THR:O	2.13	0.48
1:A:550:GLN:O	1:A:554:GLN:HG3	2.14	0.48
1:A:837:ASP:O	3:A:1109:L64:H27	2.13	0.48
1:A:673:HIS:HA	1:A:712:ARG:HE	1.78	0.48
1:A:799:GLU:CD	1:A:799:GLU:H	2.22	0.48
1:A:233:ILE:HD11	1:A:248:PHE:HD1	1.78	0.48
1:A:170:ASP:OD1	1:A:170:ASP:C	2.57	0.47
1:A:1023:HIS:ND1	1:A:1023:HIS:N	2.62	0.47
1:A:464:VAL:HG22	1:A:465:ASN:N	2.28	0.47
1:A:1067:TYR:O	1:A:1070:ASP:HB2	2.14	0.47
1:A:274:VAL:CG1	1:A:275:CYS:N	2.76	0.47
1:A:420:ILE:CG2	1:A:420:ILE:O	2.62	0.47
1:A:699:LEU:HD13	1:A:717:LEU:HB2	1.97	0.47
1:A:143:MET:C	1:A:145:GLU:N	2.72	0.47
1:A:184:ARG:HH11	1:A:184:ARG:HB2	1.79	0.47
1:A:361:PHE:HB2	1:A:420:ILE:HG13	1.97	0.47
1:A:995:MET:HE1	1:A:1009:PHE:CD1	2.50	0.47
1:A:1078:LYS:O	1:A:1081:THR:HG23	2.13	0.47
1:A:725:GLY:O	1:A:729:LEU:HG	2.15	0.47
1:A:917:THR:OG1	1:A:919:GLU:HB3	2.15	0.47
1:A:379:LEU:HD13	1:A:435:CYS:SG	2.55	0.46
1:A:383:VAL:HG22	1:A:433:ILE:HG12	1.98	0.46
1:A:952:ILE:CG2	1:A:953:MET:N	2.76	0.46
1:A:355:TRP:HE3	1:A:530:PRO:HG2	1.78	0.46
1:A:601:GLN:CD	1:A:601:GLN:C	2.83	0.46
1:A:204:SER:HB2	1:A:652:GLU:OE2	2.15	0.46
1:A:1010:GLN:O	1:A:1014:VAL:HG23	2.15	0.46
1:A:298:LYS:HE3	1:A:861:ASP:OD2	2.16	0.46
1:A:604:VAL:O	1:A:607:THR:HB	2.15	0.46
1:A:382:PHE:HD1	1:A:382:PHE:H	1.64	0.46
1:A:414:LEU:HD23	1:A:414:LEU:HA	1.80	0.46
1:A:472:ARG:O	1:A:473:PHE:HB2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:368:ILE:HD13	1:A:514:MET:CE	2.46	0.46
1:A:748:ASP:O	1:A:752:LEU:HD13	2.16	0.46
1:A:274:VAL:CG2	1:A:292:TRP:NE1	2.79	0.46
1:A:240:THR:O	1:A:244:ILE:HG13	2.15	0.45
1:A:783:PHE:CD2	1:A:783:PHE:C	2.93	0.45
1:A:199:HIS:O	1:A:200:PRO:C	2.58	0.45
1:A:834:HIS:ND1	1:A:835:GLY:N	2.64	0.45
1:A:837:ASP:O	3:A:1109:L64:H28	2.17	0.45
1:A:1042:LEU:O	1:A:1043:THR:C	2.59	0.45
1:A:223:VAL:HG12	1:A:225:HIS:NE2	2.31	0.45
1:A:623:ASP:OD1	1:A:623:ASP:C	2.59	0.45
1:A:1008:LYS:O	1:A:1011:ASP:HB2	2.17	0.45
1:A:569:ALA:O	1:A:573:GLU:HG3	2.16	0.45
1:A:1089:HIS:C	1:A:1091:VAL:N	2.74	0.45
1:A:997:THR:HG22	1:A:998:SER:N	2.31	0.45
1:A:198:MET:O	1:A:199:HIS:C	2.60	0.45
1:A:552:ARG:O	1:A:556:GLU:HG3	2.16	0.45
1:A:917:THR:HG23	1:A:920:LYS:CD	2.47	0.45
1:A:306:VAL:HG22	1:A:306:VAL:O	2.17	0.45
1:A:547:MET:HE2	1:A:552:ARG:HA	1.98	0.45
1:A:834:HIS:HB2	1:A:876:ILE:HD12	1.99	0.45
1:A:936:CYS:SG	1:A:985:PHE:HD1	2.36	0.45
1:A:184:ARG:O	1:A:188:VAL:HG23	2.17	0.45
1:A:511:GLU:C	1:A:512:ASN:OD1	2.60	0.45
1:A:947:ARG:HH21	1:A:947:ARG:HG3	1.82	0.45
3:A:1109:L64:O2	3:A:1109:L64:H19	2.16	0.45
1:A:838:LEU:HD12	1:A:877:GLY:N	2.32	0.44
1:A:947:ARG:NH1	1:A:964:ASP:O	2.49	0.44
1:A:224:ILE:C	1:A:225:HIS:HD2	2.25	0.44
1:A:393:VAL:HG23	1:A:393:VAL:O	2.16	0.44
1:A:551:LEU:HA	1:A:551:LEU:HD23	1.75	0.44
1:A:686:LEU:HD22	1:A:723:GLY:O	2.17	0.44
1:A:209:GLU:HB2	1:A:859:SER:HB3	2.00	0.44
1:A:210:TYR:HE2	1:A:859:SER:HA	1.82	0.44
1:A:525:HIS:N	1:A:526:PRO:CD	2.81	0.44
1:A:372:VAL:O	1:A:372:VAL:HG12	2.17	0.44
1:A:496:SER:HA	1:A:1043:THR:O	2.17	0.44
1:A:955:THR:OG1	1:A:959:ASN:CB	2.65	0.44
1:A:210:TYR:HE2	1:A:860:LEU:N	2.16	0.44
1:A:661:LEU:O	1:A:665:GLN:HG2	2.17	0.44
1:A:917:THR:HG1	1:A:919:GLU:HB3	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:831:ILE:HG13	1:A:881:ILE:CG1	2.48	0.44
1:A:931:SER:OG	1:A:932:CYS:N	2.50	0.44
1:A:550:GLN:O	1:A:550:GLN:HG3	2.17	0.44
1:A:514:MET:CG	1:A:515:SER:N	2.80	0.43
1:A:354:LEU:O	1:A:527:ILE:CG1	2.67	0.43
1:A:855:TRP:HZ3	1:A:864:LEU:HD11	1.82	0.43
1:A:210:TYR:CE2	1:A:860:LEU:N	2.86	0.43
1:A:378:ASP:O	1:A:379:LEU:CG	2.65	0.43
1:A:481:VAL:HA	1:A:516:ILE:O	2.18	0.43
1:A:830:ILE:HG22	1:A:831:ILE:N	2.33	0.43
1:A:931:SER:OG	1:A:960:LEU:HB3	2.19	0.43
1:A:939:THR:HB	1:A:945:GLY:HA2	2.00	0.43
1:A:1088:LEU:O	1:A:1091:VAL:HB	2.19	0.43
1:A:215:ILE:O	1:A:216:ALA:CB	2.66	0.43
1:A:221:PHE:HD2	1:A:232:THR:CG2	2.32	0.43
1:A:589:TYR:O	1:A:593:PHE:CD1	2.72	0.43
1:A:855:TRP:O	1:A:860:LEU:O	2.36	0.43
1:A:908:ASN:OD1	1:A:994:VAL:HG22	2.17	0.43
1:A:401:PRO:O	1:A:402:LYS:HG3	2.19	0.43
1:A:925:VAL:O	1:A:925:VAL:HG12	2.17	0.43
1:A:517:SER:C	1:A:518:ILE:HG23	2.44	0.42
1:A:1033:MET:O	1:A:1037:THR:HG23	2.19	0.42
1:A:423:LEU:HA	1:A:424:PRO:HD3	1.82	0.42
1:A:653:ASP:O	1:A:654:ASP:C	2.62	0.42
1:A:805:ALA:O	1:A:806:SER:C	2.62	0.42
1:A:304:HIS:C	1:A:305:VAL:HG23	2.44	0.42
1:A:362:ARG:HG3	1:A:413:TRP:CZ3	2.54	0.42
1:A:605:ALA:C	1:A:607:THR:N	2.75	0.42
1:A:829:GLY:O	1:A:881:ILE:HB	2.19	0.42
1:A:412:VAL:O	1:A:412:VAL:HG23	2.20	0.42
1:A:474:LEU:HD23	1:A:528:ALA:HB2	2.02	0.42
1:A:512:ASN:OD1	1:A:512:ASN:N	2.53	0.42
1:A:210:TYR:CE2	1:A:859:SER:HA	2.54	0.42
1:A:218:ASN:O	1:A:218:ASN:ND2	2.53	0.42
1:A:667:VAL:O	1:A:667:VAL:CG1	2.67	0.42
1:A:1066:LYS:O	1:A:1067:TYR:C	2.63	0.42
1:A:361:PHE:HD2	1:A:416:PHE:CD1	2.37	0.42
1:A:368:ILE:HG23	1:A:410:TRP:HZ3	1.83	0.42
1:A:470:ASP:CB	1:A:476:ARG:NH1	2.83	0.42
1:A:586:PRO:O	1:A:626:LEU:HD21	2.20	0.42
1:A:589:TYR:HB2	1:A:590:PRO:HD3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:905:GLU:HG2	1:A:993:PHE:CE1	2.54	0.42
3:A:1109:L64:H29A	3:A:1109:L64:H14	2.01	0.42
1:A:241:PRO:O	1:A:245:LEU:HD12	2.20	0.41
1:A:640:VAL:HG12	1:A:641:ARG:N	2.35	0.41
1:A:817:CYS:O	1:A:819:ASP:N	2.53	0.41
1:A:864:LEU:C	1:A:866:PRO:HD3	2.45	0.41
1:A:1089:HIS:O	1:A:1091:VAL:N	2.53	0.41
1:A:706:SER:O	1:A:710:GLN:HB3	2.20	0.41
1:A:829:GLY:C	1:A:881:ILE:HD12	2.45	0.41
1:A:886:THR:HG23	1:A:887:THR:O	2.20	0.41
1:A:1029:ILE:HD12	1:A:1033:MET:HE3	2.02	0.41
1:A:478:GLY:N	1:A:520:LEU:HB2	2.34	0.41
1:A:632:ASP:OD1	1:A:632:ASP:C	2.64	0.41
1:A:850:ILE:O	1:A:854:ILE:HG13	2.21	0.41
1:A:879:ILE:CG2	1:A:880:GLU:N	2.83	0.41
1:A:798:ILE:H	1:A:798:ILE:HG12	1.65	0.41
1:A:937:VAL:O	1:A:938:ALA:C	2.63	0.41
1:A:964:ASP:HA	3:A:1109:L64:C20	2.51	0.41
1:A:420:ILE:O	1:A:420:ILE:HG22	2.21	0.41
1:A:470:ASP:OD1	1:A:470:ASP:C	2.64	0.41
1:A:1060:ASN:C	1:A:1062:GLU:H	2.29	0.41
1:A:311:PRO:O	1:A:313:PRO:HD3	2.21	0.41
1:A:785:VAL:HA	1:A:786:PRO:HD3	1.91	0.41
1:A:817:CYS:C	1:A:819:ASP:H	2.28	0.41
1:A:917:THR:HG23	1:A:920:LYS:HD3	2.03	0.41
1:A:364:LYS:O	1:A:518:ILE:HA	2.21	0.41
1:A:354:LEU:O	1:A:527:ILE:CD1	2.67	0.41
1:A:838:LEU:HA	1:A:838:LEU:HD23	1.88	0.40
1:A:817:CYS:SG	1:A:819:ASP:HB3	2.62	0.40
1:A:819:ASP:HA	1:A:820:PRO:HD3	1.90	0.40
1:A:1020:LEU:O	1:A:1021:ARG:C	2.64	0.40
1:A:883:LYS:HD3	1:A:883:LYS:HA	1.86	0.40
1:A:1055:LEU:C	1:A:1056:THR:HG22	2.47	0.40
1:A:177:ARG:HG2	1:A:715:VAL:HG13	2.02	0.40
1:A:194:LYS:O	1:A:198:MET:HE3	2.21	0.40
1:A:529:LEU:HA	1:A:530:PRO:HD3	1.85	0.40
1:A:560:ALA:HB2	4:A:86:HOH:O	2.22	0.40
1:A:564:LEU:HD13	1:A:1028:ILE:CG2	2.52	0.40
1:A:621:ALA:O	1:A:622:LEU:C	2.64	0.40
1:A:622:LEU:HD21	1:A:651:LEU:HD21	2.04	0.40
1:A:1008:LYS:O	1:A:1011:ASP:CB	2.70	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1009:PHE:HE2	1:A:1072:ILE:HD13	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	828/966 (86%)	714 (86%)	94 (11%)	20 (2%)	4 17

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	303	ILE
1	A	371	PRO
1	A	216	ALA
1	A	372	VAL
1	A	638	GLU
1	A	691	ILE
1	A	964	ASP
1	A	1049	GLU
1	A	379	LEU
1	A	532	HIS
1	A	806	SER
1	A	818	ALA
1	A	1090	LEU
1	A	411	ASN
1	A	938	ALA
1	A	144	SER
1	A	162	VAL
1	A	989	PRO
1	A	215	ILE
1	A	1038	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	671/864 (78%)	604 (90%)	67 (10%)	7 24

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	162	VAL
1	A	168	VAL
1	A	203	THR
1	A	225	HIS
1	A	229	THR
1	A	235	VAL
1	A	247	SER
1	A	281	LEU
1	A	282	VAL
1	A	285	THR
1	A	306	VAL
1	A	309	THR
1	A	322	GLU
1	A	355	TRP
1	A	368	ILE
1	A	369	ASP
1	A	370	ILE
1	A	372	VAL
1	A	382	PHE
1	A	383	VAL
1	A	395	CYS
1	A	400	SER
1	A	415	GLU
1	A	418	ILE
1	A	435	CYS
1	A	469	ILE
1	A	487	ILE
1	A	488	SER
1	A	512	ASN
1	A	518	ILE

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Mol	Chain	Res	Type
1	A	519	LEU
1	A	564	LEU
1	A	585	HIS
1	A	591	LYS
1	A	601	GLN
1	A	639	ASN
1	A	647	LYS
1	A	701	SER
1	A	705	GLN
1	A	731	ASP
1	A	740	GLU
1	A	763	VAL
1	A	764	ILE
1	A	798	ILE
1	A	814	GLU
1	A	847	ILE
1	A	870	ILE
1	A	884	ASP
1	A	894	SER
1	A	913	GLU
1	A	917	THR
1	A	920	LYS
1	A	949	ASN
1	A	952	ILE
1	A	954	ILE
1	A	955	THR
1	A	989	PRO
1	A	1002	THR
1	A	1011	ASP
1	A	1023	HIS
1	A	1029	ILE
1	A	1039	MET
1	A	1049	GLU
1	A	1060	ASN
1	A	1072	ILE
1	A	1087	PHE
1	A	1091	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	218	ASN

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Mol	Chain	Res	Type
1	A	246	GLN
1	A	295	HIS
1	A	299	ASN
1	A	665	GLN
1	A	825	ASN

5.3.3 RNA [i](#)

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](#)

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	L64	A	1109	-	51,51,51	0.78	3 (5%)	69,71,71	1.93	19 (27%)
2	SO4	A	1	-	4,4,4	0.24	0	6,6,6	0.23	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	L64	A	1109	-	-	1/28/46/46	0/7/7/7

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1109	L64	C3-C6	-2.24	1.46	1.50
3	A	1109	L64	C14-N4	2.21	1.34	1.32
3	A	1109	L64	C6-N2	2.00	1.38	1.34

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1109	L64	C14-N4-N3	-5.57	102.15	105.95
3	A	1109	L64	C12-N3-N4	5.34	114.15	110.94
3	A	1109	L64	C8-C9-N3	4.91	116.53	111.88
3	A	1109	L64	C3-C6-N2	-4.68	112.82	118.66
3	A	1109	L64	C1-C2-C3	-3.21	117.20	120.36
3	A	1109	L64	C13-C15-N10	3.16	127.80	122.81
3	A	1109	L64	C13-C14-N4	3.14	113.76	111.38
3	A	1109	L64	O1-C6-N2	2.92	126.96	122.35
3	A	1109	L64	C10-C9-N3	-2.83	109.20	111.88
3	A	1109	L64	N6-C12-N3	-2.76	125.04	128.68
3	A	1109	L64	C12-N3-C9	-2.75	122.14	128.97
3	A	1109	L64	N8-C23-N7	-2.68	107.25	112.44
3	A	1109	L64	C32-N10-C29	2.65	117.54	111.57
3	A	1109	L64	C13-C15-N5	-2.33	117.01	122.71
3	A	1109	L64	C2-C3-C4	2.31	120.19	117.61
3	A	1109	L64	C13-C12-N6	2.21	127.97	126.56
3	A	1109	L64	C11-C10-C9	-2.19	106.78	110.78
3	A	1109	L64	C24-N8-C23	2.17	131.04	126.61
3	A	1109	L64	C8-C7-N2	-2.14	106.47	110.66

There are no chirality outliers.

All (1) torsion outliers are listed below:

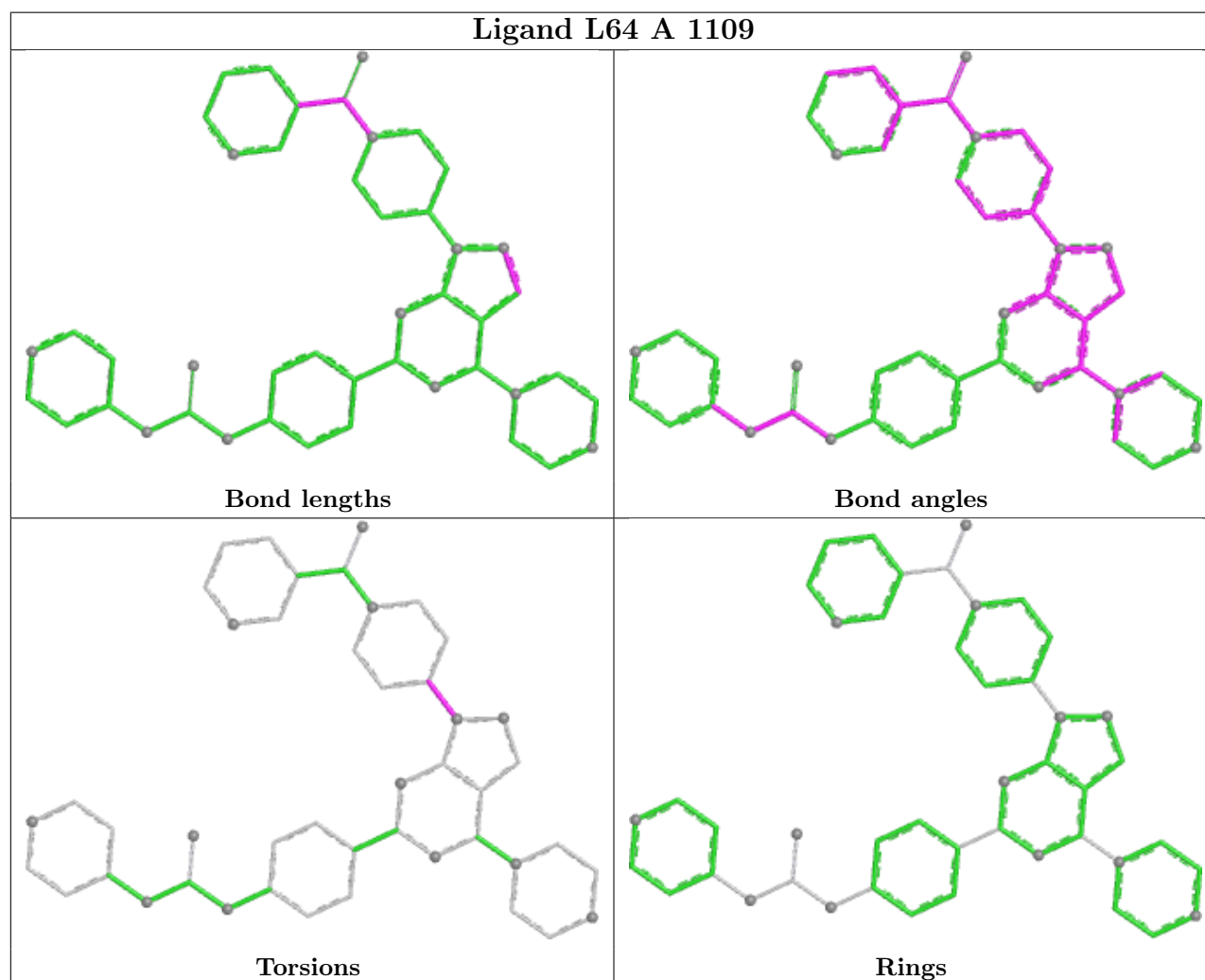
Mol	Chain	Res	Type	Atoms
3	A	1109	L64	C10-C9-N3-C12

There are no ring outliers.

1 monomer is involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1109	L64	9	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	850/966 (87%)	-0.22	9 (1%)	78 70	34, 68, 115, 148	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	543	VAL	2.8
1	A	268	GLN	2.5
1	A	1040	PRO	2.5
1	A	1092	LEU	2.4
1	A	394	LEU	2.3
1	A	967	HIS	2.2
1	A	361	PHE	2.1
1	A	901	ALA	2.1
1	A	436	GLY	2.1

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](#)

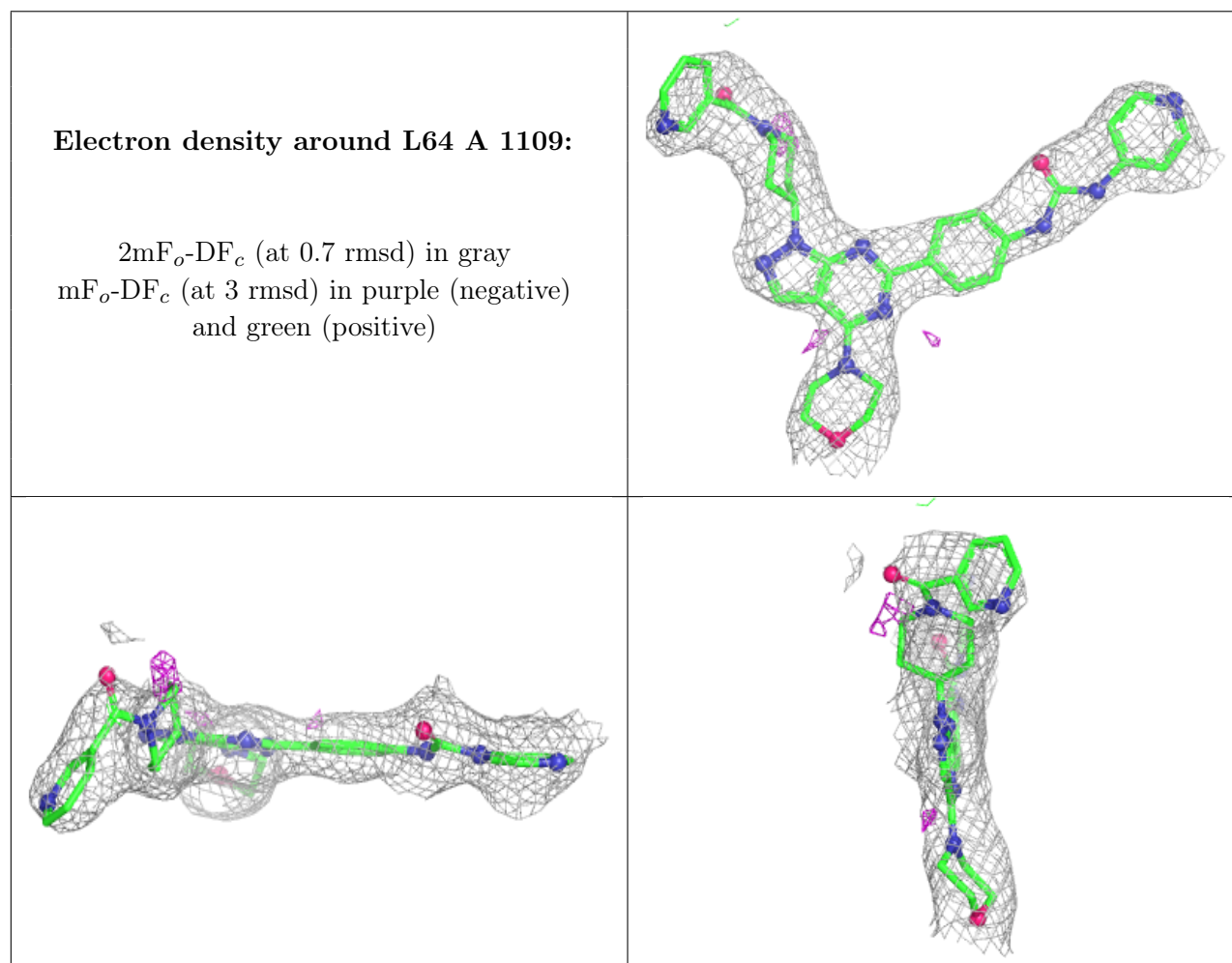
There are no oligosaccharides in this entry.

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
3	L64	A	1109	45/45	0.86	0.11	52,75,98,110	0
2	SO4	A	1	5/5	0.94	0.09	62,75,89,94	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.