



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2021 – 02:02 PM EST

PDB ID : 7JWE
Title : Gedatolisib bound to the PI3Kg catalytic subunit p110 gamma
Authors : Burke, J.E.; Rathinaswamy, M.K.; Harris, N.J.
Deposited on : 2020-08-25
Resolution : 2.55 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.17.1
buster-report : 1.1.7 (2018)
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) : 2.17.1

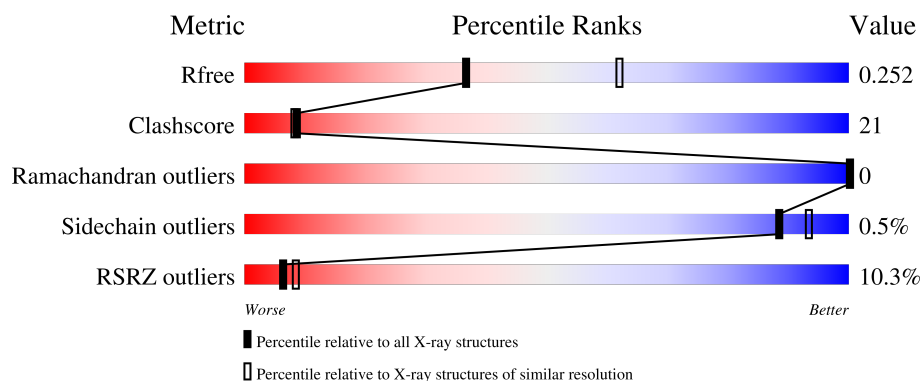
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1284 (2.56-2.52)
Clashscore	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)
RSRZ outliers	127900	1272 (2.56-2.52)

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	<div> <div>9%</div> <div>53%</div> <div>31%</div> <div>16%</div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6707 atoms, of which 41 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.

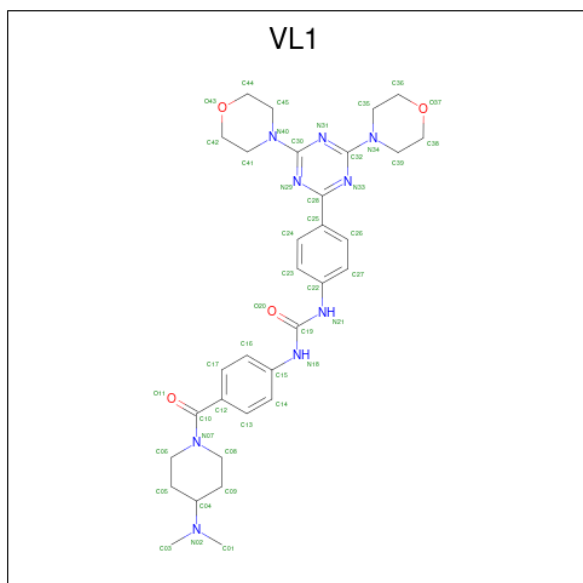
- Molecule 1 is a protein called Phosphatidylinositol 4,5-bisphosphate 3-kinase catalytic sub-unit gamma isoform.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	815	Total	C	N	O	S	0	0	0
			6612	4255	1122	1201	34			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	143	MET	-	initiating methionine	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

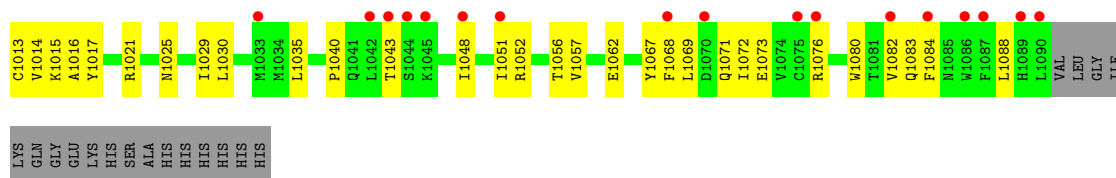
- Molecule 2 is Gedatolisib (three-letter code: VL1) (formula: C₃₂H₄₁N₉O₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	0	0
			86	32	41	9	4		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	9	Total	O	0	0
			9	9		



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	143.50Å 67.62Å 106.33Å 90.00° 95.43° 90.00°	Depositor
Resolution (Å)	44.59 – 2.55 44.59 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.4 (44.59-2.55) 90.4 (44.59-2.55)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.73 (at 2.54Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.209 , 0.253 0.209 , 0.252	Depositor DCC
R_{free} test set	1660 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	64.6	Xtriage
Anisotropy	0.296	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 64.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6707	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.07% 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: VL1

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	0.27	0/6752	0.44	0/9131

There are no bond length outliers.

There are no bond angle outliers.

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	6612	0	6641	274	0
2	A	45	41	0	3	0
3	A	9	0	0	1	0
All	All	6666	41	6641	274	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 21.

All (274) 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:272:LEU:HB3	1:A:305:VAL:HG11	1.48	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:804:MET:HG3	1:A:810:PRO:HG2	1.49	0.93
1:A:755:GLU:HG3	1:A:756:LYS:HG3	1.51	0.92
1:A:176:THR:HG23	1:A:674:ASP:HB2	1.52	0.91
1:A:1073:GLU:HA	1:A:1076:ARG:HE	1.37	0.87
1:A:983:VAL:HB	1:A:1082:VAL:HG11	1.56	0.87
1:A:992:LEU:O	1:A:996:GLY:N	2.09	0.85
1:A:218:ASN:OD1	1:A:237:PRO:HD2	1.81	0.81
1:A:606:LYS:O	1:A:609:GLN:HG2	1.83	0.79
1:A:396:GLN:O	1:A:397:ARG:NH1	2.16	0.77
1:A:1052:ARG:HG3	1:A:1057:VAL:HG21	1.66	0.76
1:A:233:ILE:HD11	1:A:248:PHE:HD1	1.50	0.76
1:A:804:MET:CG	1:A:810:PRO:HG2	2.17	0.75
1:A:215:ILE:HD11	1:A:237:PRO:HG3	1.69	0.75
1:A:1035:LEU:HD12	1:A:1048:ILE:HG12	1.66	0.74
1:A:855:TRP:CE3	1:A:862:LEU:HD13	2.22	0.74
1:A:887:THR:HG22	1:A:889:ALA:N	2.02	0.74
1:A:983:VAL:HG22	1:A:984:PRO:HD2	1.69	0.74
1:A:625:GLY:O	1:A:629:GLN:HG3	1.89	0.72
1:A:410:TRP:HB3	1:A:412:VAL:HG23	1.72	0.72
1:A:564:LEU:HD11	1:A:1048:ILE:HG22	1.72	0.72
1:A:198:MET:HE3	1:A:282:VAL:HG21	1.73	0.71
1:A:1013:CYS:HB3	1:A:1068:PHE:CE2	2.26	0.71
1:A:1003:SER:HB2	1:A:1004:PRO:HD2	1.72	0.70
1:A:435:CYS:HB3	1:A:461:LEU:HD11	1.73	0.70
1:A:272:LEU:CB	1:A:305:VAL:HG11	2.22	0.70
1:A:662:GLN:OE1	1:A:1030:LEU:HD22	1.92	0.69
1:A:1025:ASN:O	1:A:1029:ILE:HG22	1.92	0.68
1:A:1080:TRP:O	1:A:1084:PHE:N	2.16	0.68
1:A:185:MET:HE1	1:A:321:GLU:HB3	1.74	0.68
1:A:1073:GLU:HA	1:A:1076:ARG:NE	2.08	0.68
1:A:382:PHE:CE1	1:A:398:ARG:HD3	2.29	0.68
1:A:622:LEU:HD13	1:A:647:LYS:HB3	1.74	0.67
1:A:889:ALA:O	1:A:893:GLN:HG3	1.93	0.67
1:A:904:ASP:HB3	1:A:1080:TRP:HZ2	1.60	0.67
1:A:363:VAL:HG12	1:A:416:PHE:HE1	1.59	0.66
1:A:381:VAL:HA	1:A:434:TYR:O	1.96	0.66
1:A:833:LYS:HE3	1:A:836:ASP:OD2	1.95	0.66
1:A:364:LYS:HE2	1:A:411:ASN:OD1	1.96	0.65
1:A:921:PHE:O	1:A:925:VAL:HG23	1.95	0.65
1:A:185:MET:CE	1:A:321:GLU:HB3	2.28	0.65
1:A:240:THR:HG22	1:A:242:GLY:H	1.62	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:887:THR:HG22	1:A:889:ALA:H	1.61	0.64
1:A:1005:HIS:ND1	1:A:1008:LYS:HD3	2.13	0.64
1:A:370:ILE:HG12	1:A:372:VAL:H	1.62	0.64
1:A:804:MET:CB	1:A:810:PRO:HG2	2.28	0.64
1:A:308:ASP:OD1	1:A:309:THR:N	2.30	0.64
1:A:767:LEU:O	1:A:771:LEU:HG	1.96	0.64
1:A:887:THR:HG21	1:A:950:ASP:HA	1.80	0.63
1:A:202:VAL:HG13	1:A:285:THR:HG21	1.80	0.63
1:A:1008:LYS:O	1:A:1012:ILE:HG13	1.98	0.63
1:A:210:TYR:OH	1:A:856:GLU:HG3	1.98	0.62
1:A:597:LYS:HD3	1:A:600:GLN:NE2	2.14	0.62
1:A:245:LEU:HD13	1:A:272:LEU:HD11	1.81	0.62
1:A:561:THR:HB	1:A:565:ASN:HD22	1.65	0.62
1:A:966:GLY:HA3	2:A:1201:VL1:C15	2.30	0.62
1:A:862:LEU:O	1:A:931:SER:HA	2.00	0.62
1:A:1005:HIS:HA	1:A:1008:LYS:HD3	1.81	0.61
1:A:432:GLN:HB3	1:A:460:LEU:CD1	2.30	0.61
1:A:983:VAL:CG2	1:A:984:PRO:HD2	2.30	0.61
1:A:240:THR:O	1:A:244:ILE:HG13	1.99	0.61
1:A:154:LEU:O	1:A:158:ILE:HG13	2.01	0.61
1:A:561:THR:OG1	1:A:591:LYS:HE2	1.99	0.61
1:A:215:ILE:HG23	1:A:216:ALA:O	2.00	0.61
1:A:278:ASP:OD2	1:A:792:LYS:HE3	2.00	0.61
1:A:1040:PRO:O	1:A:1043:THR:HG23	2.00	0.61
1:A:215:ILE:CD1	1:A:237:PRO:HG3	2.31	0.60
1:A:586:PRO:HA	1:A:589:TYR:CD1	2.36	0.60
1:A:370:ILE:HD11	1:A:372:VAL:HB	1.82	0.60
1:A:991:PHE:O	1:A:995:MET:HG3	2.02	0.60
1:A:186:ALA:O	1:A:190:SER:HB2	2.02	0.59
1:A:988:THR:OG1	1:A:990:ASP:OD1	2.13	0.59
1:A:378:ASP:C	1:A:379:LEU:HD12	2.23	0.59
1:A:233:ILE:HD11	1:A:248:PHE:CD1	2.36	0.59
1:A:939:THR:HB	1:A:945:GLY:HA2	1.84	0.58
1:A:697:TRP:CH2	1:A:739:ILE:HD13	2.38	0.58
1:A:1052:ARG:HG3	1:A:1057:VAL:CG2	2.32	0.58
1:A:148:GLN:O	1:A:152:ARG:HG3	2.04	0.58
1:A:839:ARG:HG2	1:A:842:MET:HE2	1.86	0.58
1:A:741:MET:HB3	1:A:774:LEU:HD21	1.84	0.57
1:A:1021:ARG:NE	1:A:1056:THR:OG1	2.37	0.57
1:A:561:THR:CB	1:A:565:ASN:HD22	2.18	0.57
1:A:359:ARG:HG2	1:A:360:LYS:O	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:285:THR:HG22	1:A:289:ASN:HB2	1.87	0.57
1:A:968:ILE:O	1:A:968:ILE:HD12	2.05	0.57
1:A:1084:PHE:CE2	1:A:1088:LEU:HD11	2.39	0.57
1:A:896:VAL:O	1:A:899:THR:HG22	2.04	0.57
1:A:945:GLY:O	1:A:985:PHE:HA	2.05	0.57
1:A:564:LEU:CD1	1:A:1048:ILE:HG22	2.35	0.57
1:A:1069:LEU:O	1:A:1073:GLU:HG3	2.04	0.57
1:A:304:HIS:CG	1:A:823:LEU:HD11	2.40	0.56
1:A:895:THR:HG21	1:A:906:VAL:HG13	1.87	0.56
1:A:202:VAL:HG12	1:A:203:THR:N	2.20	0.56
1:A:422:ASP:HB3	1:A:599:GLY:O	2.05	0.56
1:A:804:MET:HB2	1:A:810:PRO:HG2	1.87	0.56
1:A:838:LEU:HD12	1:A:877:GLY:HA3	1.87	0.56
1:A:983:VAL:CB	1:A:1082:VAL:HG11	2.33	0.56
1:A:180:LEU:O	1:A:183:PRO:HD2	2.06	0.56
1:A:480:TYR:HB2	1:A:518:ILE:HD11	1.87	0.56
1:A:1067:TYR:O	1:A:1071:GLN:HG2	2.06	0.56
1:A:370:ILE:CD1	1:A:372:VAL:HB	2.37	0.55
1:A:753:SER:OG	1:A:809:LYS:HE2	2.07	0.55
1:A:887:THR:CG2	1:A:950:ASP:HA	2.35	0.55
1:A:888:ILE:HB	1:A:949:ASN:OD1	2.07	0.55
1:A:176:THR:O	1:A:180:LEU:HG	2.07	0.55
1:A:303:ILE:HD12	1:A:303:ILE:H	1.71	0.55
1:A:749:ILE:CD1	1:A:767:LEU:HD13	2.36	0.55
1:A:697:TRP:HH2	1:A:739:ILE:HD13	1.71	0.55
1:A:758:ASP:O	1:A:759:VAL:HG23	2.07	0.55
1:A:855:TRP:CD2	1:A:862:LEU:HD13	2.40	0.55
1:A:888:ILE:HG23	1:A:907:LEU:HD21	1.89	0.55
1:A:210:TYR:CD1	1:A:211:LEU:HD13	2.42	0.55
1:A:240:THR:HG22	1:A:242:GLY:N	2.21	0.55
1:A:996:GLY:O	1:A:1003:SER:HB3	2.07	0.54
1:A:224:ILE:O	1:A:230:SER:HA	2.08	0.54
1:A:750:LYS:HE3	1:A:808:LYS:HB3	1.90	0.54
1:A:271:VAL:HG23	1:A:282:VAL:HG12	1.89	0.54
1:A:434:TYR:CE2	1:A:460:LEU:HB2	2.43	0.54
1:A:706:SER:O	1:A:710:GLN:HB3	2.08	0.53
1:A:939:THR:HG21	1:A:962:HIS:CE1	2.44	0.53
1:A:361:PHE:HA	1:A:420:ILE:HD11	1.91	0.53
1:A:362:ARG:HG2	1:A:415:GLU:OE1	2.09	0.53
1:A:432:GLN:HB3	1:A:460:LEU:HD11	1.89	0.53
1:A:435:CYS:HB3	1:A:461:LEU:CD1	2.39	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:736:VAL:HA	1:A:739:ILE:HG12	1.91	0.53
1:A:767:LEU:HD22	1:A:803:VAL:HG23	1.90	0.53
1:A:787:TYR:CE1	1:A:880:GLU:HB2	2.44	0.53
1:A:804:MET:O	1:A:805:ALA:HB3	2.08	0.53
1:A:198:MET:SD	1:A:282:VAL:HG11	2.50	0.52
1:A:804:MET:HG3	1:A:810:PRO:CG	2.31	0.52
1:A:555:LEU:HD11	1:A:575:LEU:HD23	1.92	0.52
1:A:589:TYR:CD2	1:A:593:PHE:HE2	2.28	0.52
1:A:896:VAL:HG11	1:A:903:LYS:HG3	1.91	0.52
1:A:964:ASP:HA	2:A:1201:VL1:C22	2.40	0.52
1:A:905:GLU:HG2	1:A:993:PHE:CE2	2.44	0.52
1:A:384:GLU:OE2	1:A:398:ARG:NH1	2.43	0.52
1:A:546:GLU:OE1	1:A:546:GLU:HA	2.11	0.51
1:A:1072:ILE:O	1:A:1076:ARG:HG3	2.09	0.51
1:A:410:TRP:HB3	1:A:412:VAL:CG2	2.38	0.51
1:A:549:ASN:OD1	1:A:553:LYS:HE3	2.09	0.51
1:A:635:PHE:O	1:A:641:ARG:HD2	2.11	0.51
1:A:552:ARG:HG2	1:A:556:GLU:OE2	2.11	0.51
1:A:946:ASP:CG	1:A:983:VAL:HG12	2.31	0.51
1:A:240:THR:HG23	1:A:241:PRO:HD2	1.93	0.51
1:A:372:VAL:HG12	1:A:373:LEU:N	2.26	0.51
1:A:562:ASP:HB2	1:A:563:PRO:CD	2.41	0.50
1:A:1013:CYS:HB3	1:A:1068:PHE:HE2	1.76	0.50
1:A:466:LEU:HD11	1:A:476:ARG:HD3	1.93	0.50
1:A:1004:PRO:O	1:A:1008:LYS:HG3	2.11	0.50
1:A:917:THR:CG2	1:A:919:GLU:HG2	2.42	0.49
1:A:1011:ASP:OD1	1:A:1015:LYS:NZ	2.39	0.49
1:A:968:ILE:O	1:A:969:LEU:HD12	2.12	0.49
1:A:432:GLN:NE2	1:A:434:TYR:OH	2.45	0.49
1:A:749:ILE:HD11	1:A:767:LEU:HA	1.95	0.49
1:A:888:ILE:CG2	1:A:907:LEU:HD21	2.42	0.49
1:A:586:PRO:O	1:A:626:LEU:HD21	2.12	0.49
1:A:760:SER:OG	1:A:763:VAL:HG23	2.12	0.49
1:A:180:LEU:C	1:A:183:PRO:HD2	2.33	0.49
1:A:862:LEU:HD21	1:A:1016:ALA:HB2	1.93	0.49
1:A:302:GLU:HB2	1:A:304:HIS:CD2	2.48	0.49
1:A:1003:SER:HB2	1:A:1004:PRO:CD	2.43	0.49
1:A:806:SER:HB3	1:A:810:PRO:HD2	1.95	0.48
1:A:199:HIS:O	1:A:199:HIS:ND1	2.47	0.48
1:A:842:MET:HE1	3:A:1305:HOH:O	2.13	0.48
1:A:410:TRP:C	1:A:412:VAL:HG23	2.34	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:561:THR:CG2	1:A:565:ASN:HD22	2.26	0.48
1:A:1011:ASP:O	1:A:1014:VAL:HG22	2.13	0.48
1:A:301:GLU:HG2	1:A:302:GLU:N	2.28	0.48
1:A:731:ASP:OD2	1:A:784:ARG:NE	2.40	0.48
1:A:964:ASP:HA	2:A:1201:VL1:N21	2.28	0.48
1:A:1017:TYR:O	1:A:1021:ARG:HG3	2.12	0.48
1:A:939:THR:HB	1:A:945:GLY:CA	2.43	0.48
1:A:163:THR:HG22	1:A:177:ARG:HH12	1.80	0.47
1:A:428:LEU:HD23	1:A:467:LEU:HD23	1.95	0.47
1:A:609:GLN:O	1:A:612:ALA:HB3	2.15	0.47
1:A:853:SER:O	1:A:857:THR:HG23	2.14	0.47
1:A:202:VAL:CG1	1:A:203:THR:N	2.77	0.47
1:A:307:LEU:O	1:A:307:LEU:HD23	2.15	0.47
1:A:364:LYS:HB3	1:A:519:LEU:HB3	1.97	0.47
1:A:804:MET:SD	1:A:812:TRP:HE3	2.38	0.47
1:A:865:LEU:CD2	1:A:882:VAL:HG11	2.45	0.47
1:A:935:TYR:CE2	1:A:961:PHE:HA	2.50	0.47
1:A:989:PRO:HB2	1:A:1080:TRP:CZ2	2.50	0.47
1:A:177:ARG:HG2	1:A:715:VAL:HG13	1.96	0.47
1:A:739:ILE:O	1:A:743:GLN:HG3	2.15	0.46
1:A:767:LEU:HD11	1:A:771:LEU:HD11	1.96	0.46
1:A:220:ILE:HD12	1:A:237:PRO:HA	1.96	0.46
1:A:246:GLN:O	1:A:250:THR:HG23	2.16	0.46
1:A:410:TRP:O	1:A:412:VAL:HG23	2.15	0.46
1:A:385:ALA:CB	1:A:416:PHE:HZ	2.29	0.46
1:A:808:LYS:HD3	1:A:835:GLY:HA3	1.97	0.46
1:A:304:HIS:ND1	1:A:823:LEU:HD11	2.31	0.46
1:A:743:GLN:O	1:A:747:LEU:HG	2.16	0.46
1:A:983:VAL:HB	1:A:1082:VAL:CG1	2.37	0.45
1:A:1082:VAL:HG23	1:A:1083:GLN:N	2.32	0.45
1:A:364:LYS:HB2	1:A:413:TRP:CZ3	2.51	0.45
1:A:1062:GLU:OE1	1:A:1062:GLU:N	2.36	0.45
1:A:865:LEU:HD21	1:A:882:VAL:HG11	1.99	0.45
1:A:302:GLU:O	1:A:304:HIS:HD2	2.00	0.45
1:A:364:LYS:HD2	1:A:413:TRP:CZ2	2.52	0.45
1:A:862:LEU:HD12	1:A:862:LEU:N	2.30	0.45
1:A:469:ILE:HG22	1:A:473:PHE:HA	1.99	0.45
1:A:800:LYS:HD3	1:A:814:GLU:OE1	2.17	0.45
1:A:806:SER:OG	1:A:807:LYS:N	2.49	0.45
1:A:224:ILE:HD12	1:A:233:ILE:HD12	1.99	0.45
1:A:834:HIS:HA	1:A:875:LYS:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:864:LEU:C	1:A:866:PRO:HD3	2.38	0.45
1:A:463:TYR:CD1	1:A:487:ILE:HD11	2.52	0.44
1:A:907:LEU:HD13	1:A:990:ASP:HB2	1.99	0.44
1:A:245:LEU:HD11	1:A:307:LEU:HD12	1.99	0.44
1:A:372:VAL:O	1:A:373:LEU:HB2	2.17	0.44
1:A:640:VAL:O	1:A:643:ILE:HG12	2.18	0.44
1:A:202:VAL:CG1	1:A:285:THR:HG21	2.47	0.44
1:A:271:VAL:CG2	1:A:282:VAL:HG12	2.48	0.44
1:A:303:ILE:HD12	1:A:303:ILE:N	2.32	0.44
1:A:597:LYS:HB2	1:A:603:ILE:CD1	2.47	0.44
1:A:904:ASP:CB	1:A:1080:TRP:HZ2	2.28	0.44
1:A:689:LYS:HG2	1:A:728:MET:SD	2.57	0.44
1:A:280:TYR:HB3	1:A:282:VAL:HG13	2.00	0.44
1:A:157:LEU:HD23	1:A:157:LEU:HA	1.79	0.44
1:A:359:ARG:O	1:A:420:ILE:HG12	2.18	0.44
1:A:361:PHE:HD2	1:A:416:PHE:CD1	2.35	0.44
1:A:480:TYR:HB2	1:A:518:ILE:CD1	2.47	0.44
1:A:835:GLY:O	1:A:875:LYS:HD2	2.18	0.44
1:A:196:TYR:O	1:A:689:LYS:HE3	2.19	0.43
1:A:315:LEU:O	1:A:727:ALA:HB2	2.18	0.43
1:A:425:LYS:HD2	1:A:473:PHE:CE2	2.53	0.43
1:A:1082:VAL:HG23	1:A:1083:GLN:H	1.83	0.43
1:A:273:ARG:O	1:A:305:VAL:HG13	2.19	0.43
1:A:245:LEU:HD13	1:A:272:LEU:CD1	2.47	0.43
1:A:287:ILE:HA	1:A:290:PHE:HD1	1.84	0.43
1:A:370:ILE:HG12	1:A:372:VAL:N	2.30	0.43
1:A:627:THR:HG21	1:A:648:LEU:HG	2.00	0.43
1:A:317:GLU:O	1:A:726:THR:HG23	2.19	0.43
1:A:564:LEU:HD11	1:A:1048:ILE:CG2	2.46	0.43
1:A:699:LEU:O	1:A:703:ILE:HG13	2.18	0.43
1:A:949:ASN:HB2	1:A:1083:GLN:OE1	2.19	0.43
1:A:547:MET:HE3	1:A:578:PHE:CE2	2.54	0.43
1:A:385:ALA:HB1	1:A:416:PHE:HZ	1.84	0.42
1:A:743:GLN:HB3	1:A:876:ILE:CD1	2.49	0.42
1:A:887:THR:HG21	1:A:889:ALA:HB3	2.01	0.42
1:A:1002:THR:HG22	1:A:1003:SER:N	2.34	0.42
1:A:597:LYS:HD2	1:A:603:ILE:CD1	2.49	0.42
1:A:1073:GLU:HB3	1:A:1076:ARG:HH21	1.84	0.42
1:A:210:TYR:CE1	1:A:211:LEU:HD13	2.55	0.42
1:A:862:LEU:HB3	1:A:934:GLY:HA3	2.01	0.42
1:A:584:LYS:O	1:A:586:PRO:HD3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:703:ILE:CD1	1:A:714:ALA:HA	2.50	0.42
1:A:630:LEU:HB2	1:A:644:ALA:HB2	2.02	0.42
1:A:231:GLN:HG3	1:A:232:THR:O	2.20	0.42
1:A:291:GLN:HA	1:A:291:GLN:OE1	2.20	0.42
1:A:806:SER:HB3	1:A:810:PRO:CD	2.50	0.42
1:A:1051:ILE:HG23	1:A:1052:ARG:N	2.35	0.42
1:A:741:MET:SD	1:A:774:LEU:HD11	2.59	0.41
1:A:905:GLU:HG2	1:A:993:PHE:CZ	2.55	0.41
1:A:483:HIS:HB3	1:A:513:SER:OG	2.21	0.41
1:A:567:LEU:HD11	1:A:591:LYS:HB3	2.03	0.41
1:A:593:PHE:CZ	1:A:611:LEU:HD21	2.55	0.41
1:A:561:THR:HG21	1:A:565:ASN:HD22	1.85	0.41
1:A:487:ILE:HG22	1:A:488:SER:N	2.36	0.41
1:A:547:MET:HG2	1:A:578:PHE:CD1	2.56	0.41
1:A:927:ARG:HH11	1:A:959:ASN:ND2	2.19	0.41
1:A:146:GLU:OE1	1:A:146:GLU:HA	2.20	0.41
1:A:988:THR:HG21	1:A:1083:GLN:HG3	2.02	0.41
1:A:207:LEU:HD22	1:A:288:LYS:HG3	2.02	0.41
1:A:887:THR:CG2	1:A:889:ALA:HB3	2.51	0.41
1:A:470:ASP:HB3	1:A:476:ARG:NH2	2.36	0.41
1:A:888:ILE:HG12	1:A:952:ILE:O	2.21	0.40
1:A:285:THR:CG2	1:A:289:ASN:HB2	2.50	0.40
1:A:501:LYS:HE3	1:A:501:LYS:HB3	1.82	0.40
1:A:547:MET:HG2	1:A:578:PHE:CE1	2.57	0.40
1:A:231:GLN:HG3	1:A:232:THR:N	2.36	0.40
1:A:561:THR:HG21	1:A:565:ASN:ND2	2.37	0.40
1:A:819:ASP:OD1	1:A:820:PRO:HD2	2.22	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	793/966 (82%)	755 (95%)	38 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	731/864 (85%)	727 (100%)	4 (0%)	88	93

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	190	SER
1	A	282	VAL
1	A	415	GLU
1	A	731	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	432	GLN
1	A	565	ASN
1	A	959	ASN

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 [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	$\# Z > 2$
2	VL1	A	1201	-	49,50,50	2.95	14 (28%)	67,69,69	2.61	18 (26%)

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	VL1	A	1201	-	-	3/32/58/58	0/6/6/6

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1201	VL1	C06-N07	7.65	1.60	1.47
2	A	1201	VL1	C10-N07	7.58	1.51	1.34
2	A	1201	VL1	C30-N40	7.49	1.49	1.35
2	A	1201	VL1	C08-N07	6.90	1.59	1.47
2	A	1201	VL1	C32-N34	5.67	1.46	1.35
2	A	1201	VL1	C19-N21	5.48	1.48	1.37
2	A	1201	VL1	C05-C04	-5.28	1.38	1.52
2	A	1201	VL1	C09-C04	-4.91	1.39	1.52
2	A	1201	VL1	C19-N18	4.49	1.46	1.37
2	A	1201	VL1	C12-C10	4.30	1.57	1.50
2	A	1201	VL1	C15-N18	3.01	1.47	1.41
2	A	1201	VL1	C22-N21	2.99	1.47	1.41
2	A	1201	VL1	C25-C28	2.75	1.55	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1201	VL1	O20-C19	-2.37	1.18	1.23

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1201	VL1	C08-N07-C06	-10.92	91.60	112.62
2	A	1201	VL1	C32-N33-C28	7.34	119.76	114.43
2	A	1201	VL1	C30-N29-C28	7.28	119.72	114.43
2	A	1201	VL1	C05-C06-N07	5.72	119.52	110.82
2	A	1201	VL1	C42-C41-N40	4.13	117.63	110.02
2	A	1201	VL1	C32-N31-C30	4.09	119.68	113.55
2	A	1201	VL1	C06-N07-C10	3.79	134.77	122.78
2	A	1201	VL1	N31-C30-N29	-3.76	120.11	126.31
2	A	1201	VL1	N33-C32-N31	-3.60	120.36	126.31
2	A	1201	VL1	C08-N07-C10	3.43	133.63	122.78
2	A	1201	VL1	N33-C28-N29	-3.01	120.36	125.23
2	A	1201	VL1	N31-C32-N34	2.99	120.80	117.11
2	A	1201	VL1	O43-C42-C41	2.71	117.76	111.80
2	A	1201	VL1	N31-C30-N40	2.45	120.13	117.11
2	A	1201	VL1	C41-N40-C30	-2.32	117.85	121.69
2	A	1201	VL1	C03-N02-C04	-2.29	108.45	112.39
2	A	1201	VL1	N29-C30-N40	2.15	119.77	117.11
2	A	1201	VL1	C25-C28-N29	2.07	120.82	117.33

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1201	VL1	N31-C30-N40-C41
2	A	1201	VL1	N29-C30-N40-C41
2	A	1201	VL1	C09-C04-N02-C03

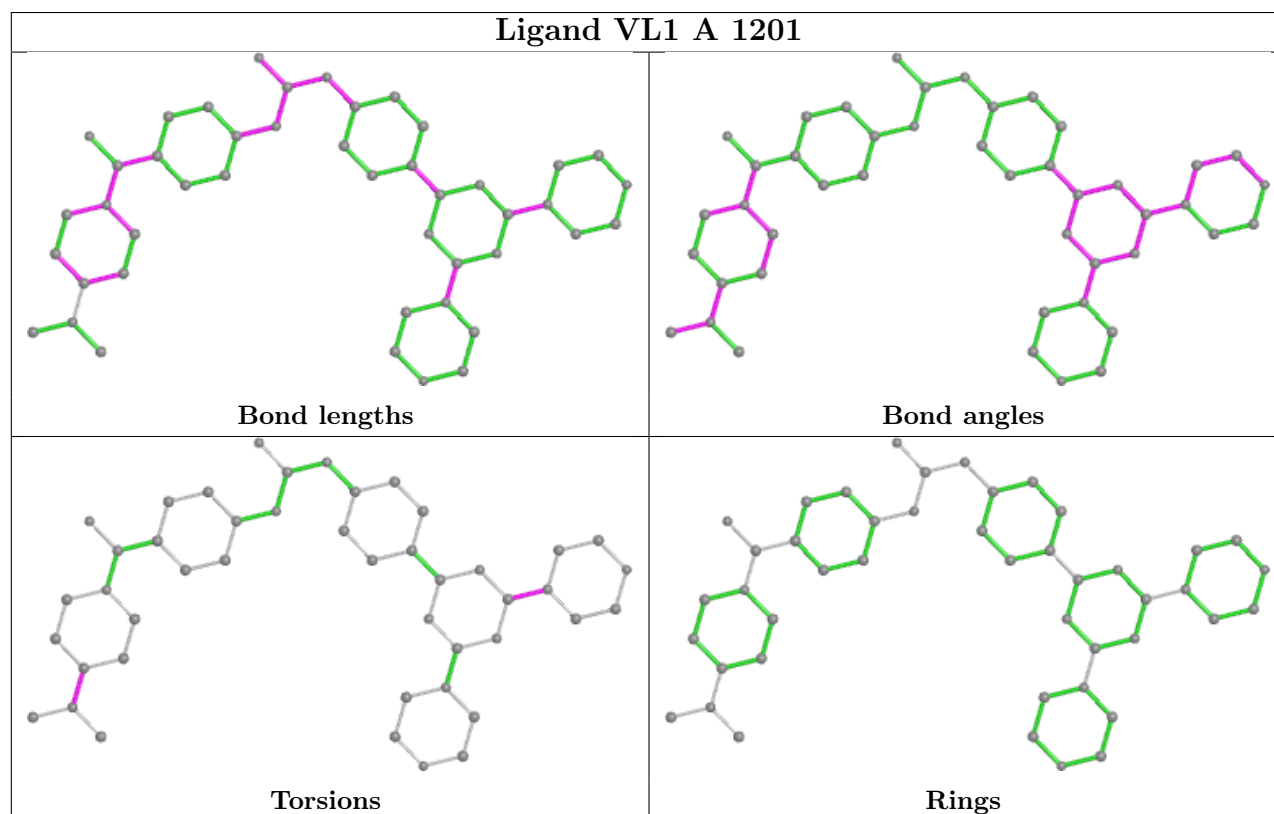
There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1201	VL1	3	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 ⓘ

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	815/966 (84%)	0.72	84 (10%) 6 8	50, 85, 133, 174	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	235	VAL	6.9
1	A	995	MET	5.5
1	A	270	PHE	5.1
1	A	217	ASN	4.8
1	A	759	VAL	4.6
1	A	911	LEU	4.3
1	A	1082	VAL	4.2
1	A	307	LEU	4.2
1	A	404	PHE	4.1
1	A	373	LEU	4.1
1	A	757	TYR	4.0
1	A	378	ASP	4.0
1	A	212	TRP	3.9
1	A	218	ASN	3.9
1	A	991	PHE	3.9
1	A	1042	LEU	3.9
1	A	907	LEU	3.9
1	A	1075	CYS	3.9
1	A	1084	PHE	3.8
1	A	221	PHE	3.6
1	A	1087	PHE	3.6
1	A	1089	HIS	3.6
1	A	220	ILE	3.5
1	A	902	PHE	3.5
1	A	489	GLY	3.5
1	A	1090	LEU	3.5
1	A	869	CYS	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	754	ALA	3.4
1	A	372	VAL	3.4
1	A	237	PRO	3.3
1	A	233	ILE	3.2
1	A	219	CYS	3.2
1	A	1068	PHE	3.2
1	A	245	LEU	3.1
1	A	216	ALA	3.0
1	A	231	GLN	3.0
1	A	1086	TRP	2.9
1	A	234	LYS	2.9
1	A	188	VAL	2.8
1	A	1076	ARG	2.8
1	A	870	ILE	2.8
1	A	321	GLU	2.8
1	A	895	THR	2.7
1	A	303	ILE	2.7
1	A	249	PHE	2.7
1	A	774	LEU	2.7
1	A	269	ASP	2.6
1	A	772	GLU	2.6
1	A	662	GLN	2.6
1	A	320	LYS	2.6
1	A	969	LEU	2.6
1	A	741	MET	2.6
1	A	1070	ASP	2.6
1	A	842	MET	2.6
1	A	146	GLU	2.5
1	A	211	LEU	2.5
1	A	248	PHE	2.5
1	A	994	VAL	2.5
1	A	823	LEU	2.4
1	A	982	ARG	2.4
1	A	1045	LYS	2.4
1	A	660	LEU	2.4
1	A	550	GLN	2.4
1	A	379	LEU	2.4
1	A	1048	ILE	2.3
1	A	624	VAL	2.3
1	A	900	GLY	2.3
1	A	401	PRO	2.3
1	A	244	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	993	PHE	2.3
1	A	317	GLU	2.2
1	A	1033	MET	2.2
1	A	301	GLU	2.2
1	A	236	SER	2.2
1	A	545	ALA	2.2
1	A	989	PRO	2.2
1	A	1044	SER	2.2
1	A	756	LYS	2.1
1	A	1043	THR	2.1
1	A	909	HIS	2.1
1	A	659	TYR	2.1
1	A	1051	ILE	2.1
1	A	713	PHE	2.1
1	A	771	LEU	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](#)

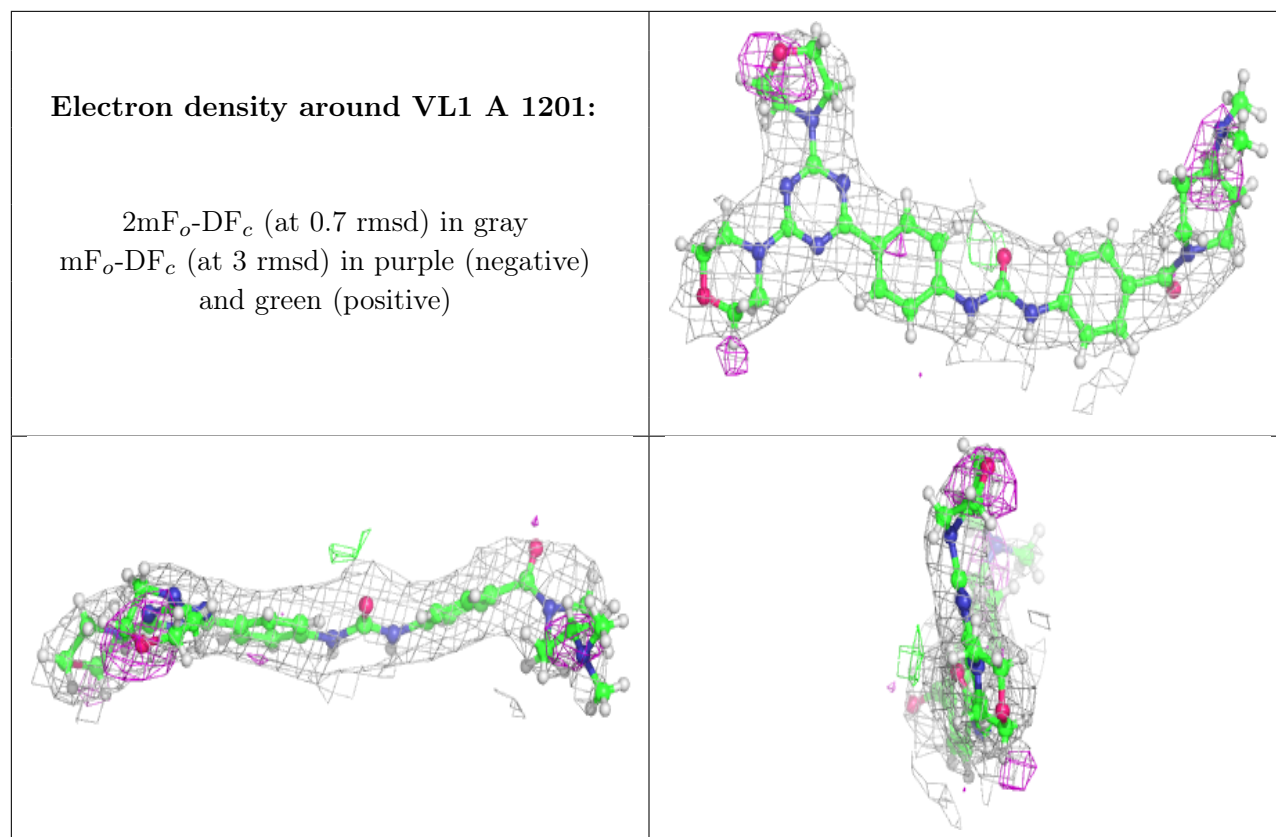
There are no monosaccharides in this entry.

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(\AA^2)	Q<0.9
2	VL1	A	1201	45/45	0.85	0.23	60,81,132,138	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.