



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:54 PM UTC

PDB ID : 5ENN / pdb_00005enn
Title : The crystal structure of Human VPS34 in complex with a selective and potent inhibitor
Authors : Kearney, E.P.
Deposited on : 2015-11-09
Resolution : 2.70 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

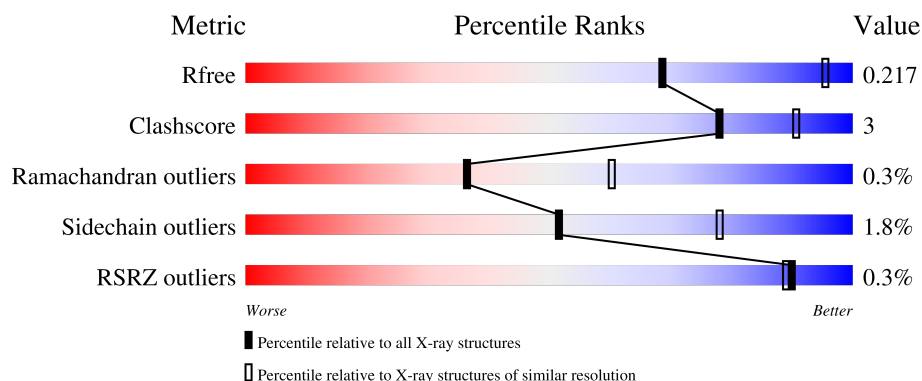
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	625	
1	B	625	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8831 atoms, of which 8 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 3-kinase catalytic subunit type 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	526	Total	C	N	O	S	0	2	0
			4231	2702	717	788	24			
1	B	531	Total	C	N	O	S	0	2	0
			4242	2711	715	792	24			

There are 60 discrepancies between the modelled and reference sequences:

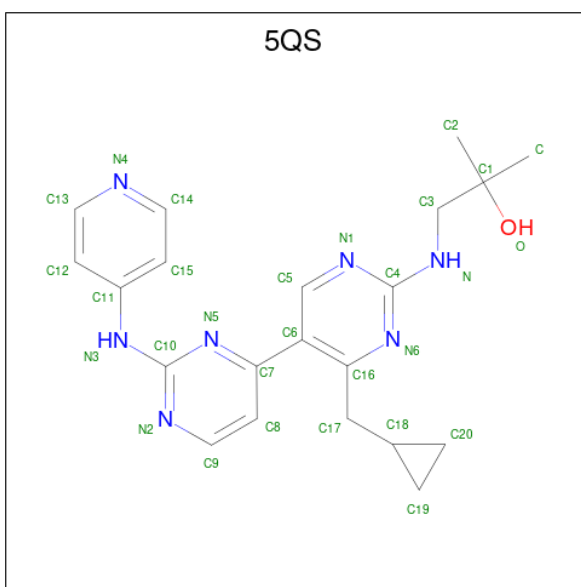
Chain	Residue	Modelled	Actual	Comment	Reference
A	263	MET	-	initiating methionine	UNP Q8NEB9
A	264	SER	-	expression tag	UNP Q8NEB9
A	265	TYR	-	expression tag	UNP Q8NEB9
A	266	TYR	-	expression tag	UNP Q8NEB9
A	267	HIS	-	expression tag	UNP Q8NEB9
A	268	HIS	-	expression tag	UNP Q8NEB9
A	269	HIS	-	expression tag	UNP Q8NEB9
A	270	HIS	-	expression tag	UNP Q8NEB9
A	271	HIS	-	expression tag	UNP Q8NEB9
A	272	HIS	-	expression tag	UNP Q8NEB9
A	273	ASP	-	expression tag	UNP Q8NEB9
A	274	TYR	-	expression tag	UNP Q8NEB9
A	275	ASP	-	expression tag	UNP Q8NEB9
A	276	ILE	-	expression tag	UNP Q8NEB9
A	277	PRO	-	expression tag	UNP Q8NEB9
A	278	THR	-	expression tag	UNP Q8NEB9
A	279	THR	-	expression tag	UNP Q8NEB9
A	280	GLU	-	expression tag	UNP Q8NEB9
A	281	ASN	-	expression tag	UNP Q8NEB9
A	282	LEU	-	expression tag	UNP Q8NEB9
A	283	TYR	-	expression tag	UNP Q8NEB9
A	284	PHE	-	expression tag	UNP Q8NEB9
A	285	GLN	-	expression tag	UNP Q8NEB9
A	286	GLY	-	expression tag	UNP Q8NEB9
A	287	ALA	-	expression tag	UNP Q8NEB9

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	288	MET	-	expression tag	UNP Q8NEB9
A	289	GLY	-	expression tag	UNP Q8NEB9
A	290	SER	-	expression tag	UNP Q8NEB9
A	291	GLY	-	expression tag	UNP Q8NEB9
A	292	ILE	-	expression tag	UNP Q8NEB9
B	263	MET	-	initiating methionine	UNP Q8NEB9
B	264	SER	-	expression tag	UNP Q8NEB9
B	265	TYR	-	expression tag	UNP Q8NEB9
B	266	TYR	-	expression tag	UNP Q8NEB9
B	267	HIS	-	expression tag	UNP Q8NEB9
B	268	HIS	-	expression tag	UNP Q8NEB9
B	269	HIS	-	expression tag	UNP Q8NEB9
B	270	HIS	-	expression tag	UNP Q8NEB9
B	271	HIS	-	expression tag	UNP Q8NEB9
B	272	HIS	-	expression tag	UNP Q8NEB9
B	273	ASP	-	expression tag	UNP Q8NEB9
B	274	TYR	-	expression tag	UNP Q8NEB9
B	275	ASP	-	expression tag	UNP Q8NEB9
B	276	ILE	-	expression tag	UNP Q8NEB9
B	277	PRO	-	expression tag	UNP Q8NEB9
B	278	THR	-	expression tag	UNP Q8NEB9
B	279	THR	-	expression tag	UNP Q8NEB9
B	280	GLU	-	expression tag	UNP Q8NEB9
B	281	ASN	-	expression tag	UNP Q8NEB9
B	282	LEU	-	expression tag	UNP Q8NEB9
B	283	TYR	-	expression tag	UNP Q8NEB9
B	284	PHE	-	expression tag	UNP Q8NEB9
B	285	GLN	-	expression tag	UNP Q8NEB9
B	286	GLY	-	expression tag	UNP Q8NEB9
B	287	ALA	-	expression tag	UNP Q8NEB9
B	288	MET	-	expression tag	UNP Q8NEB9
B	289	GLY	-	expression tag	UNP Q8NEB9
B	290	SER	-	expression tag	UNP Q8NEB9
B	291	GLY	-	expression tag	UNP Q8NEB9
B	292	ILE	-	expression tag	UNP Q8NEB9

- Molecule 2 is 1-[[4-(cyclopropylmethyl)-5-[2-(pyridin-4-ylamino)pyrimidin-4-yl]pyrimidin-2-yl]amino]-2-methyl-propan-2-ol (CCD ID: 5QS) (formula: C₂₁H₂₅N₇O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			29	21	7	1		
2	B	1	Total	C	N	O	0	0
			29	21	7	1		

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Na	0	0
			1	1		
3	B	1	Total	Na	0	0
			1	1		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

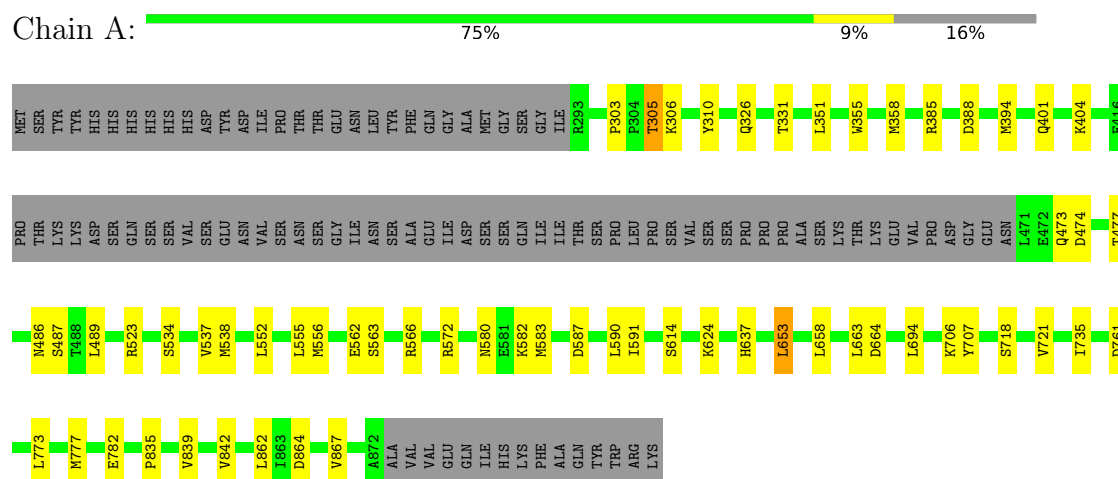
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	139	Total 139	O 139	0	0
6	B	140	Total 140	O 140	0	0

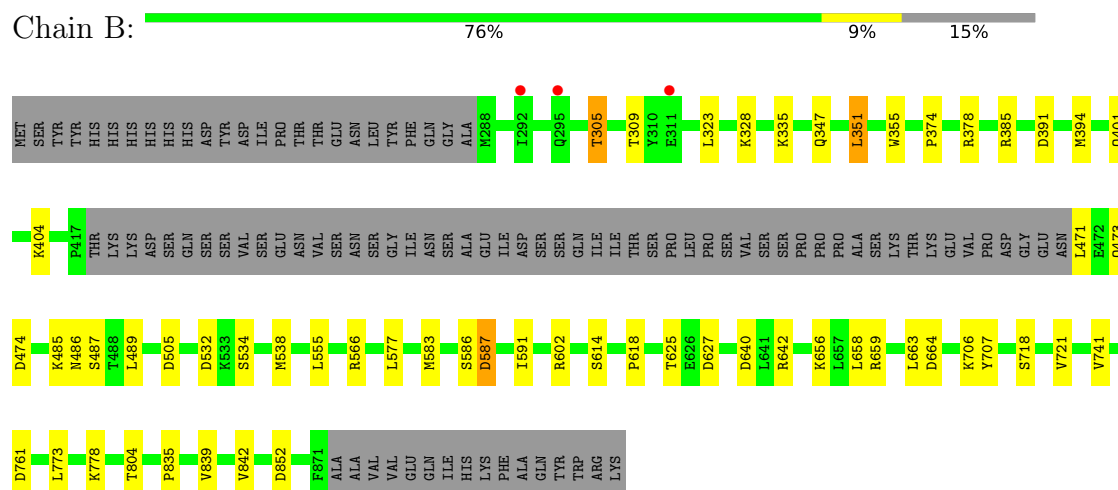
3 Residue-property plots [i](#)

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 3-kinase catalytic subunit type 3



- Molecule 1: Phosphatidylinositol 3-kinase catalytic subunit type 3



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	113.73Å 114.48Å 143.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.01 – 2.70 48.01 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.5 (48.01-2.70) 97.7 (48.01-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.23 (at 2.69Å)	Xtriage
Refinement program	BUSTER 2.11.6	Depositor
R, R_{free}	0.173 , 0.201 0.188 , 0.217	Depositor DCC
R_{free} test set	2615 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	58.4	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 53.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.298 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8831	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.14% 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: SO4, 5QS, GOL, NA

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	0.85	2/4313 (0.0%)	1.39	17/5833 (0.3%)
1	B	0.85	1/4323 (0.0%)	1.38	21/5848 (0.4%)
All	All	0.85	3/8636 (0.0%)	1.39	38/11681 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	773	LEU	CA-C	6.31	1.57	1.53
1	A	773	LEU	CA-C	5.94	1.57	1.53
1	A	735	ILE	CA-C	5.01	1.59	1.52

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	486	ASN	CA-C-N	6.70	129.56	120.38
1	B	486	ASN	C-N-CA	6.70	129.56	120.38
1	B	474	ASP	CA-CB-CG	6.69	119.29	112.60
1	A	486	ASN	CA-C-N	6.52	129.31	120.38
1	A	486	ASN	C-N-CA	6.52	129.31	120.38
1	A	474	ASP	CA-CB-CG	6.51	119.11	112.60
1	B	587	ASP	CA-C-N	6.20	129.49	122.48
1	B	587	ASP	C-N-CA	6.20	129.49	122.48
1	A	664	ASP	CA-CB-CG	6.06	118.66	112.60
1	B	664	ASP	CA-CB-CG	5.82	118.42	112.60
1	A	310	TYR	CA-C-N	5.81	128.07	120.28
1	A	310	TYR	C-N-CA	5.81	128.07	120.28
1	B	566	ARG	CA-C-N	5.74	127.90	120.44
1	B	566	ARG	C-N-CA	5.74	127.90	120.44
1	A	566	ARG	CA-C-N	5.51	127.60	120.44
1	A	566	ARG	C-N-CA	5.51	127.60	120.44

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	335	LYS	CA-C-N	5.45	128.12	120.28
1	B	335	LYS	C-N-CA	5.45	128.12	120.28
1	B	505	ASP	CA-CB-CG	5.43	118.03	112.60
1	A	563	SER	N-CA-C	-5.43	100.85	109.59
1	A	835	PRO	CA-C-N	5.42	129.27	120.60
1	A	835	PRO	C-N-CA	5.42	129.27	120.60
1	A	326	GLN	CA-C-N	5.36	127.78	120.54
1	A	326	GLN	C-N-CA	5.36	127.78	120.54
1	B	323	LEU	CA-C-N	5.29	128.10	120.38
1	B	323	LEU	C-N-CA	5.29	128.10	120.38
1	B	659	ARG	CA-C-N	5.16	127.15	120.44
1	B	659	ARG	C-N-CA	5.16	127.15	120.44
1	B	852	ASP	CA-C-N	5.10	127.11	120.28
1	B	852	ASP	C-N-CA	5.10	127.11	120.28
1	B	835	PRO	CA-C-N	5.06	128.69	120.60
1	B	835	PRO	C-N-CA	5.06	128.69	120.60
1	A	388	ASP	CA-CB-CG	5.04	117.64	112.60
1	B	842	VAL	CA-C-N	5.03	127.33	120.54
1	B	842	VAL	C-N-CA	5.03	127.33	120.54
1	A	842	VAL	CA-C-N	5.01	127.30	120.54
1	A	842	VAL	C-N-CA	5.01	127.30	120.54
1	A	864	ASP	CA-CB-CG	5.00	117.61	112.60

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	4231	0	4223	25	0
1	B	4242	0	4231	25	0
2	A	29	0	0	0	0
2	B	29	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	B	6	8	8	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	5	0	0	0	0
6	A	139	0	0	1	0
6	B	140	0	0	2	0
All	All	8823	8	8462	48	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 (48) 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:777:MET:HE2	1:A:862:LEU:HB3	1.64	0.78
1:A:537:VAL:HG11	6:B:1066:HOH:O	1.95	0.66
1:B:555:LEU:HB2	1:B:583:MET:HE2	1.78	0.66
1:B:555:LEU:HB2	1:B:583:MET:CE	2.26	0.66
1:B:485:LYS:HE3	6:B:1116:HOH:O	1.95	0.65
1:B:587:ASP:HA	1:B:602:ARG:O	2.00	0.61
1:A:580:ASN:ND2	1:A:587:ASP:HB3	2.16	0.60
1:A:556:MET:HE3	1:A:637:HIS:CE1	2.36	0.60
1:B:741:VAL:O	1:B:778[B]:LYS:HD2	2.02	0.60
1:A:614:SER:HB2	6:A:1063:HOH:O	2.02	0.59
1:A:394[A]:MET:HE2	1:A:653:LEU:HD12	1.86	0.57
1:A:303:PRO:HD2	1:A:306:LYS:HB2	1.86	0.56
1:A:582:LYS:HG3	1:A:583:MET:N	2.20	0.55
1:A:487:SER:HA	1:A:538:MET:HE3	1.91	0.53
1:A:477:THR:HG23	1:A:523:ARG:HH22	1.74	0.53
1:B:401:GLN:O	1:B:404:LYS:HG2	2.09	0.53
1:A:401:GLN:O	1:A:404:LYS:HG2	2.09	0.52
1:B:625:THR:HG22	1:B:627:ASP:H	1.75	0.51
1:B:305:THR:HG21	1:B:839:VAL:HG11	1.92	0.51
1:B:351:LEU:O	1:B:355:TRP:HB2	2.10	0.51
1:A:305:THR:HG21	1:A:839:VAL:HG11	1.94	0.50
1:B:487:SER:HA	1:B:538:MET:HE3	1.93	0.49
1:A:587:ASP:CG	1:B:532:ASP:HB2	2.38	0.49
1:B:351:LEU:HD23	1:B:355:TRP:CZ3	2.48	0.48
1:B:640:ASP:OD1	1:B:642:ARG:HD3	2.13	0.48
1:A:351:LEU:O	1:A:355:TRP:HB2	2.14	0.47
1:B:385:ARG:HH12	1:B:473:GLN:HB2	1.80	0.47
1:B:351:LEU:HD23	1:B:355:TRP:HZ3	1.79	0.47
1:A:562:GLU:OE1	1:A:572:ARG:HG3	2.16	0.46
1:B:555:LEU:HB2	1:B:583:MET:HE1	1.98	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:THR:HG23	1:A:358:MET:HE1	1.98	0.45
1:B:577:LEU:O	1:B:586:SER:HB2	2.16	0.45
1:B:374:PRO:O	1:B:378:ARG:HG2	2.17	0.45
1:A:706:LYS:HD3	1:A:707:TYR:CZ	2.52	0.44
1:B:706:LYS:HD3	1:B:707:TYR:CZ	2.52	0.44
1:A:477:THR:CG2	1:A:523:ARG:HH22	2.31	0.44
1:A:582:LYS:HG3	1:A:583:MET:H	1.81	0.44
1:B:391:ASP:O	1:B:394[B]:MET:HB2	2.17	0.44
1:A:587:ASP:OD2	1:B:532:ASP:HB2	2.19	0.43
1:A:385:ARG:HH12	1:A:473:GLN:HB2	1.83	0.43
1:B:718:SER:HB3	1:B:721:VAL:HB	2.01	0.42
1:B:658:LEU:HD22	1:B:663:LEU:HD23	2.02	0.42
1:A:658:LEU:HD22	1:A:663:LEU:HD23	2.01	0.42
1:B:614:SER:HB3	1:B:618:PRO:HD2	2.01	0.42
1:A:718:SER:HB3	1:A:721:VAL:HB	2.01	0.41
1:B:394[A]:MET:HE3	1:B:656:LYS:HE3	2.01	0.41
1:A:694:LEU:HD11	1:A:782:GLU:HG2	2.02	0.41
1:A:552:LEU:HD12	1:A:555:LEU:HD23	2.02	0.41

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	524/625 (84%)	509 (97%)	13 (2%)	2 (0%)	30	54
1	B	529/625 (85%)	514 (97%)	14 (3%)	1 (0%)	43	68
All	All	1053/1250 (84%)	1023 (97%)	27 (3%)	3 (0%)	36	60

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	590	LEU
1	A	761	ASP
1	B	761	ASP

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	463/565 (82%)	456 (98%)	7 (2%)	57	81
1	B	462/565 (82%)	452 (98%)	10 (2%)	45	74
All	All	925/1130 (82%)	908 (98%)	17 (2%)	51	78

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	305	THR
1	A	489	LEU
1	A	534	SER
1	A	591	ILE
1	A	624	LYS
1	A	653	LEU
1	A	867	VAL
1	B	305	THR
1	B	309	THR
1	B	328	LYS
1	B	347	GLN
1	B	351	LEU
1	B	471	LEU
1	B	489	LEU
1	B	534	SER
1	B	591	ILE
1	B	804	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	307	GLN
1	A	491	ASN
1	A	546	GLN
1	A	794	GLN
1	B	325	ASN
1	B	338	ASN
1	B	491	ASN
1	B	546	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](#)

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](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

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	$\# Z > 2$
5	SO4	B	904	-	4,4,4	0.27	0	6,6,6	0.24	0
4	GOL	B	903	-	5,5,5	0.10	0	5,5,5	0.17	0
2	5QS	B	901	-	32,32,32	1.82	6 (18%)	41,45,45	1.11	2 (4%)
2	5QS	A	901	-	32,32,32	1.75	8 (25%)	41,45,45	1.10	2 (4%)

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
4	GOL	B	903	-	-	0/4/4/4	-
2	5QS	B	901	-	-	0/18/20/20	0/4/4/4
2	5QS	A	901	-	-	0/18/20/20	0/4/4/4

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	901	5QS	C10-N2	5.46	1.42	1.34
2	A	901	5QS	C7-N5	4.44	1.41	1.34
2	B	901	5QS	C3-C1	3.43	1.58	1.52
2	B	901	5QS	C7-N5	3.28	1.39	1.34
2	B	901	5QS	C12-C11	3.28	1.44	1.39
2	B	901	5QS	C15-C11	2.69	1.43	1.39
2	A	901	5QS	C15-C11	2.65	1.43	1.39
2	A	901	5QS	C3-C1	2.57	1.57	1.52
2	A	901	5QS	C10-N2	2.55	1.38	1.34
2	B	901	5QS	C4-N	2.54	1.38	1.34
2	A	901	5QS	C4-N	2.30	1.38	1.34
2	A	901	5QS	C4-N1	2.28	1.37	1.34
2	A	901	5QS	C12-C11	2.28	1.43	1.39
2	A	901	5QS	C6-C16	2.13	1.45	1.41

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	901	5QS	N1-C4-N6	-3.07	123.46	126.42
2	A	901	5QS	C17-C16-N6	-3.00	111.42	117.08
2	B	901	5QS	N-C4-N1	2.38	120.41	117.21
2	A	901	5QS	N-C4-N1	2.19	120.15	117.21

There are no chirality outliers.

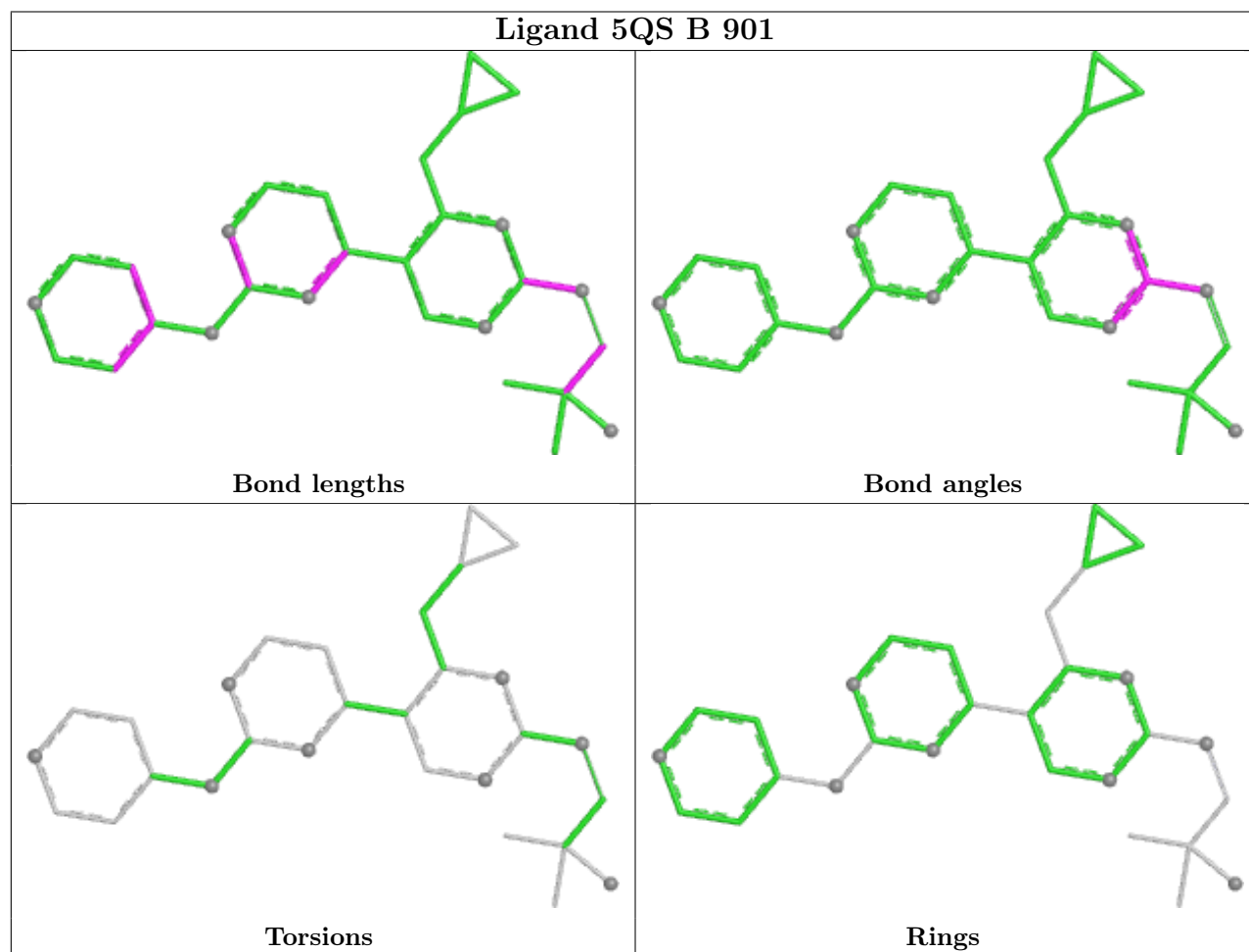
There are no torsion outliers.

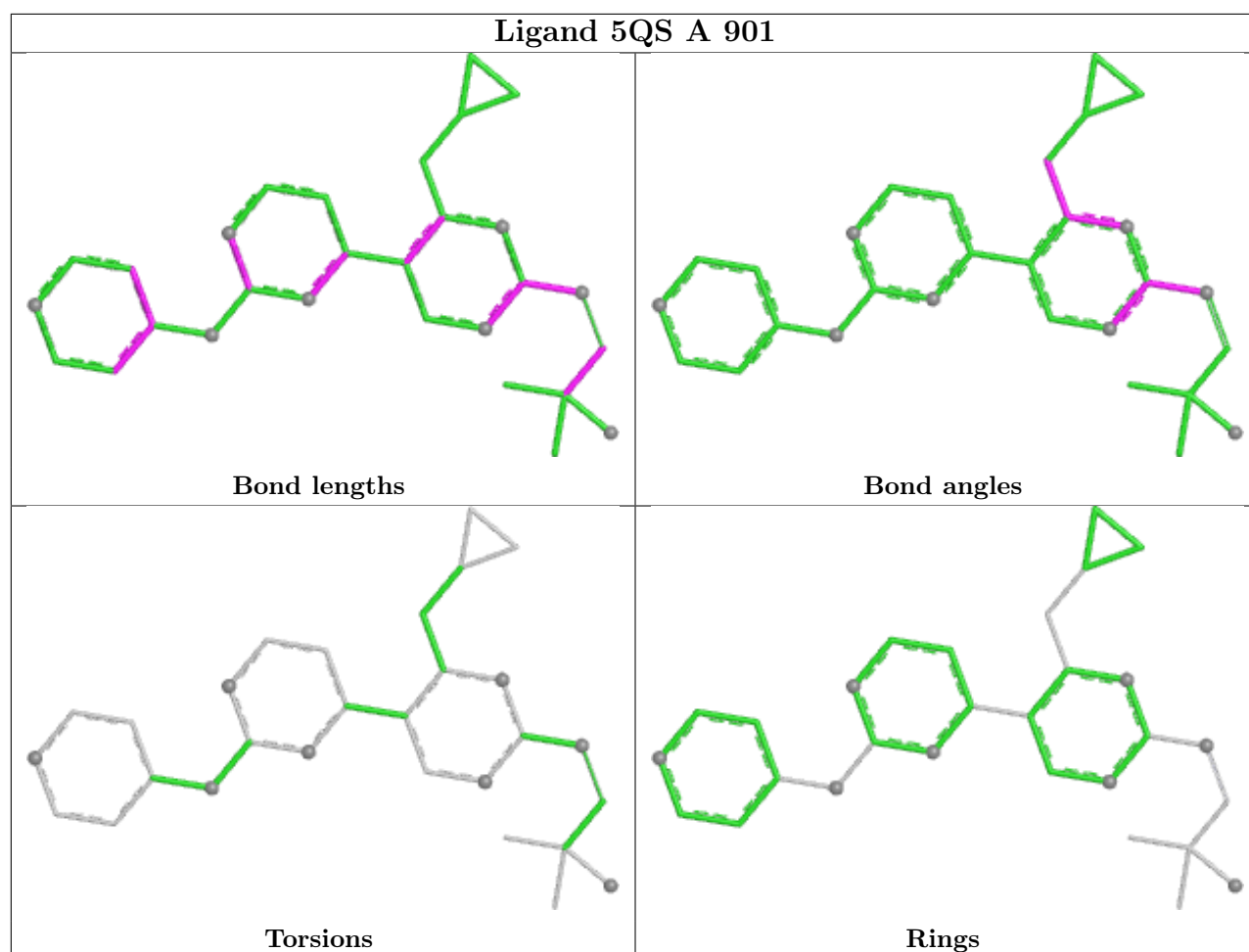
There are no ring outliers.

No monomer is involved in short contacts.

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	526/625 (84%)	-1.17	0	100 100	31, 69, 120, 165	2 (0%)
1	B	531/625 (84%)	-1.08	3 (0%)	85 85	26, 69, 125, 206	2 (0%)
All	All	1057/1250 (84%)	-1.13	3 (0%)	90 89	26, 69, 124, 206	4 (0%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	292	ILE	3.2
1	B	295	GLN	2.1
1	B	311	GLU	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(Å ²)	Q < 0.9
3	NA	A	902	1/1	0.98	0.10	62,62,62,62	0
3	NA	B	902	1/1	0.99	0.21	69,69,69,69	0

Continued on next page...

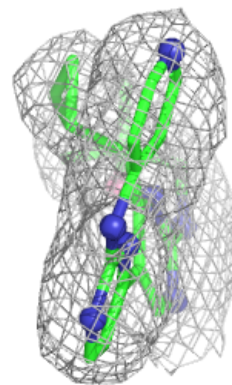
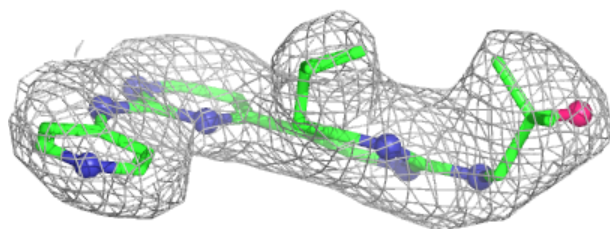
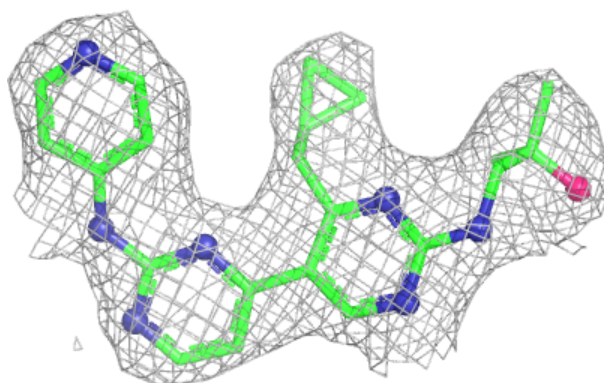
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	GOL	B	903	6/6	0.99	0.07	97,99,100,100	0
5	SO4	B	904	5/5	0.99	0.05	106,107,109,110	0
2	5QS	A	901	29/29	1.00	0.03	34,44,50,53	0
2	5QS	B	901	29/29	1.00	0.03	32,42,50,51	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.

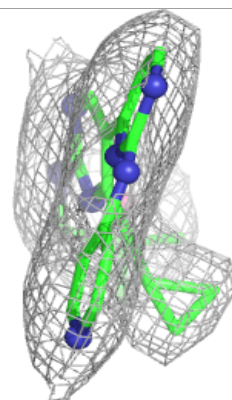
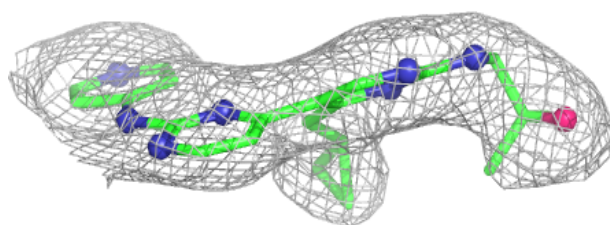
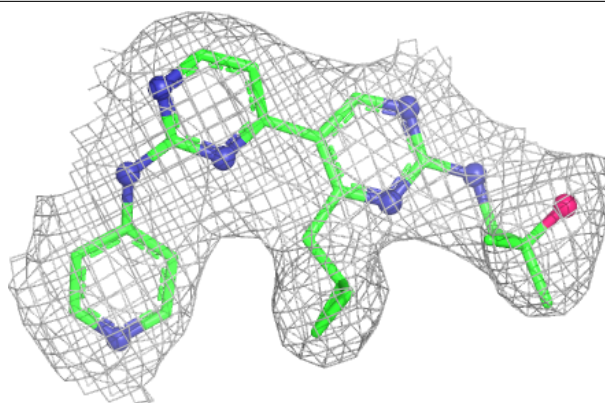
Electron density around 5QS A 901:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around 5QS B 901:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.