



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:24 AM UTC

PDB ID : 4AJW / pdb_00004ajw
Title : Discovery and Optimization of New Benzimidazole- and Benzoxazole-Pyrimidone Selective PI3KBeta Inhibitors for the Treatment of Phosphatase and TENSin homologue (PTEN)-Deficient Cancers
Authors : Certal, V.; Halley, F.; Virone-Oddos, A.; Delorme, C.; Karlsson, A.; Rak, A.; Thompson, F.; Filoche-Romme, B.; El-Ahmad, Y.; Carry, J.C.; Abecassis, P.Y.; Lejeune, P.; Bonnevaux, H.; Nicolas, J.P.; Bertrand, T.; Marquette, J.P.; Michot, N.; Benard, T.; Below, P.; Vade, I.; Chatreaux, F.; Lebourg, G.; Pilorge, F.; Angouillant-Boniface, O.; Louboutin, A.; Lengauer, C.; Schio, L.
Deposited on : 2012-02-20
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)

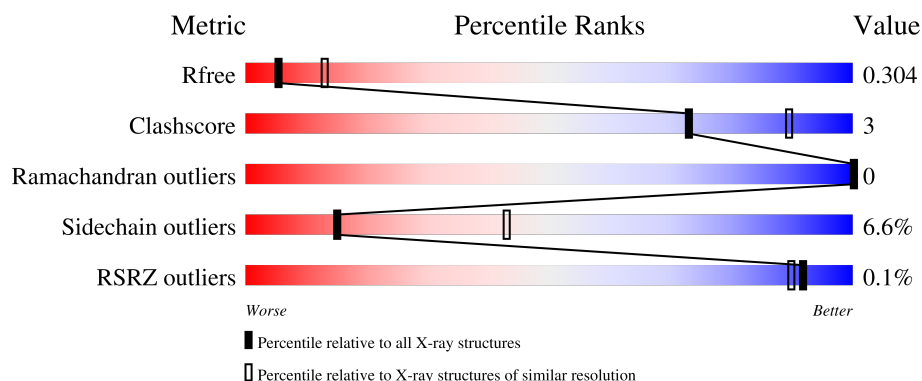
1 Overall quality at a glance

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	934	 73% 8% • 18%
1	B	934	 75% 8% • 15%

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

2 Entry composition [i](#)

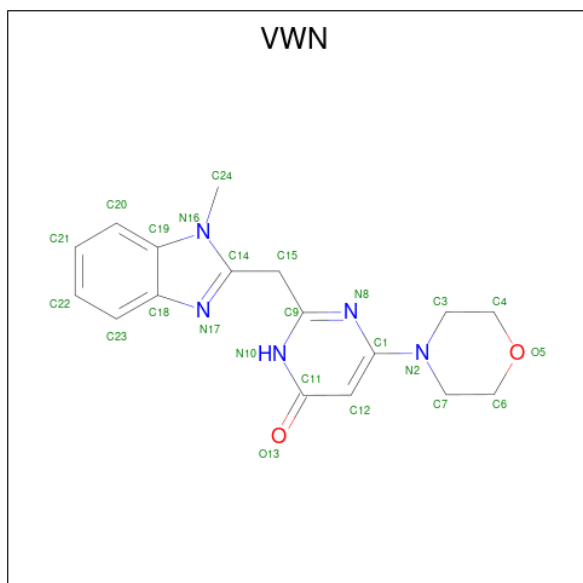
There are 3 unique types of molecules in this entry. The entry contains 12596 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 DELTA ISOFORM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	765	Total	C	N	O	S	0	0	0
			6151	3942	1042	1116	51			
1	B	792	Total	C	N	O	S	0	0	0
			6363	4076	1076	1158	53			

- Molecule 2 is 2-[(1-methyl-1H-benzimidazol-2-yl)methyl]-6-morpholin-4-ylpyrimidin-4(3H)-one (CCD ID: VWN) (formula: $C_{17}H_{19}N_5O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			24	17	5	2		
2	B	1	Total	C	N	O	0	0
			24	17	5	2		

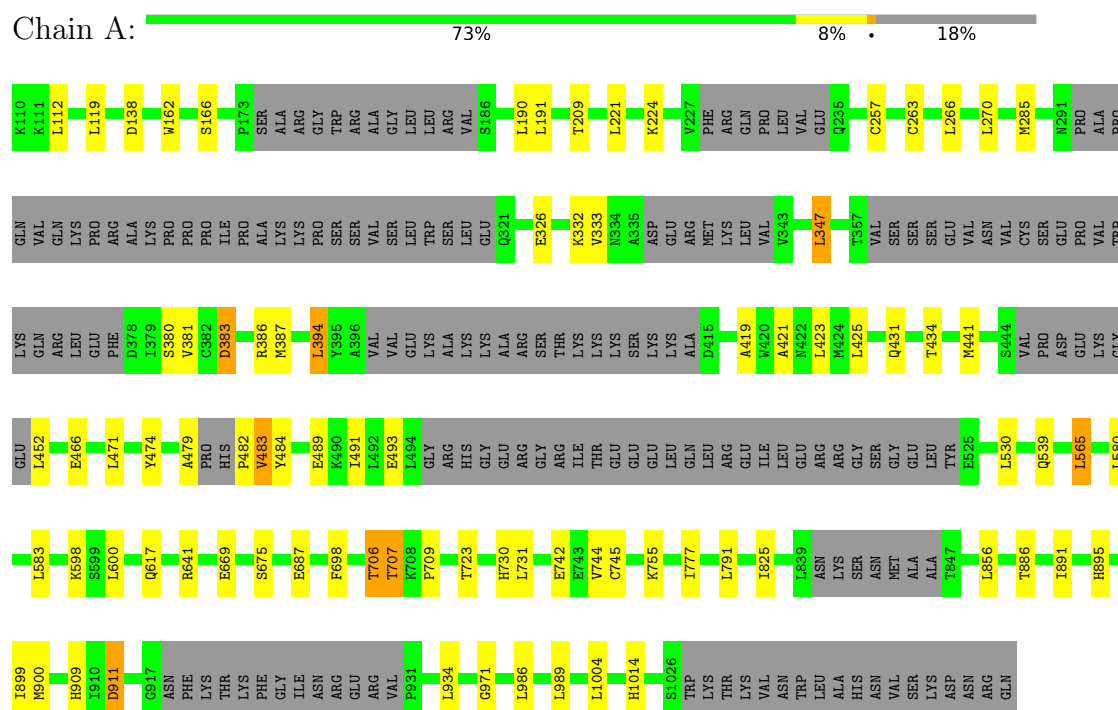
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	19	Total	O	0	0
			19	19		
3	B	15	Total	O	0	0
			15	15		

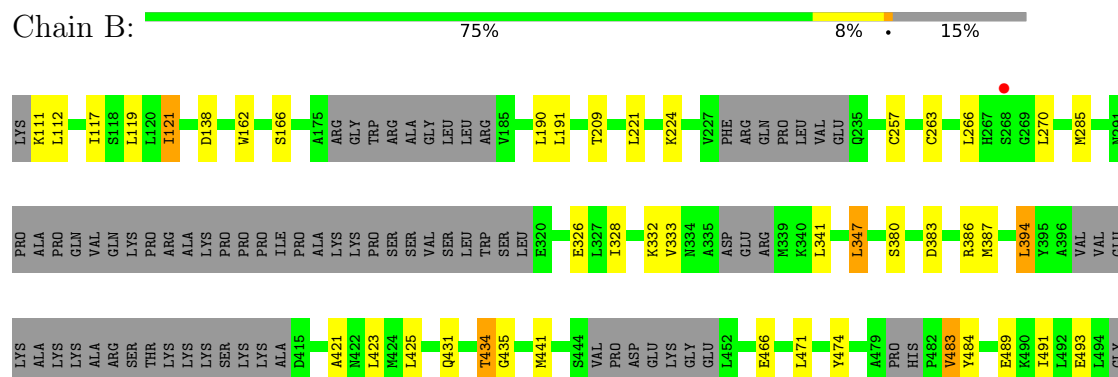
3 Residue-property plots

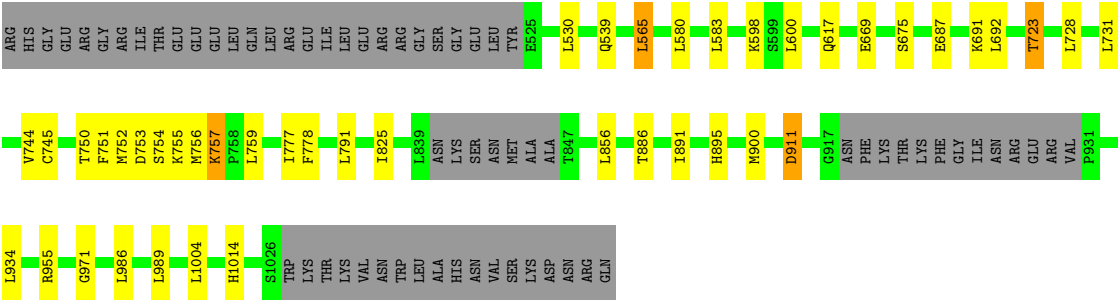
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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 DELTA ISOFORM



• Molecule 1: PHOSPHATIDYLINOSITOL-4,5-BISPHOSPHATE 3-KINASE CATALYTIC SUBUNIT DELTA ISOFORM





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.40Å 220.47Å 78.26Å 90.00° 113.87° 90.00°	Depositor
Resolution (Å)	71.57 – 2.80 71.57 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.4 (71.57-2.80) 98.2 (71.57-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.13 (at 2.82Å)	Xtriage
Refinement program	BUSTER-TNT 2.9.7	Depositor
R, R_{free}	0.252 , 0.277 0.281 , 0.304	Depositor DCC
R_{free} test set	2422 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	68.5	Xtriage
Anisotropy	0.236	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 81.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.457 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12596	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.11% 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: VWN

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	0.74	0/6282	1.29	3/8470 (0.0%)
1	B	0.76	2/6499 (0.0%)	1.29	3/8767 (0.0%)
All	All	0.75	2/12781 (0.0%)	1.29	6/17237 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	757	LYS	C-O	-5.32	1.21	1.23
1	B	745	CYS	CA-C	-5.25	1.46	1.53

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	754	SER	N-CA-C	-7.81	97.84	108.38
1	B	911	ASP	N-CA-C	7.03	119.14	110.91
1	A	911	ASP	N-CA-C	6.89	118.97	110.91
1	B	756	MET	N-CA-C	6.67	121.42	113.28
1	A	745	CYS	CA-C-N	5.75	128.01	120.60
1	A	745	CYS	C-N-CA	5.75	128.01	120.60

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	6151	0	6129	35	0
1	B	6363	0	6340	46	0
2	A	24	0	19	1	0
2	B	24	0	19	1	0
3	A	19	0	0	0	0
3	B	15	0	0	0	0
All	All	12596	0	12507	81	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 (81) 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:B:753:ASP:CG	1:B:757:LYS:NZ	2.08	1.12
1:B:753:ASP:CG	1:B:757:LYS:HZ1	1.79	0.91
1:B:753:ASP:HA	1:B:757:LYS:HZ2	1.32	0.91
1:B:753:ASP:HA	1:B:757:LYS:NZ	1.85	0.90
1:A:707:THR:OG1	1:A:709:PRO:HD2	1.74	0.86
1:A:989:LEU:O	1:A:989:LEU:HG	1.75	0.84
1:B:431:GLN:HG3	1:B:484:TYR:CE1	2.16	0.81
1:A:483:VAL:CG2	1:A:483:VAL:O	2.30	0.80
1:A:483:VAL:O	1:A:483:VAL:HG22	1.82	0.78
1:B:989:LEU:O	1:B:989:LEU:HG	1.87	0.75
1:A:431:GLN:HG3	1:A:484:TYR:CE1	2.21	0.75
1:A:730:HIS:HD2	1:A:742:GLU:OE1	1.70	0.74
1:B:483:VAL:O	1:B:483:VAL:CG2	2.36	0.73
1:B:753:ASP:CA	1:B:757:LYS:HZ2	2.03	0.72
1:B:753:ASP:OD2	1:B:757:LYS:NZ	2.18	0.71
1:B:753:ASP:CA	1:B:757:LYS:NZ	2.54	0.71
1:B:117:ILE:O	1:B:121:ILE:HG23	1.89	0.71
1:B:753:ASP:CG	1:B:757:LYS:HZ2	2.05	0.64
1:B:431:GLN:HG3	1:B:484:TYR:CD1	2.34	0.62
1:A:431:GLN:HG3	1:A:484:TYR:CD1	2.34	0.62
1:A:479:ALA:HB3	1:A:482:PRO:HD2	1.81	0.61
1:B:753:ASP:CG	1:B:757:LYS:HZ3	2.07	0.61
1:A:706:THR:OG1	1:A:707:THR:N	2.29	0.60
1:B:483:VAL:O	1:B:483:VAL:HG22	2.02	0.60
1:B:989:LEU:O	1:B:989:LEU:CG	2.45	0.59
1:A:479:ALA:CB	1:A:482:PRO:HD2	2.32	0.59
1:B:421:ALA:HB2	1:B:441:MET:HG2	1.84	0.59
1:B:989:LEU:C	1:B:989:LEU:HD12	2.30	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:394:LEU:HD13	1:B:441:MET:HE1	1.88	0.55
1:B:753:ASP:CB	1:B:757:LYS:NZ	2.70	0.54
1:B:753:ASP:CB	1:B:757:LYS:HZ2	2.21	0.54
1:A:421:ALA:HB2	1:A:441:MET:HG2	1.91	0.53
1:A:479:ALA:HB3	1:A:482:PRO:CD	2.39	0.53
1:A:989:LEU:O	1:A:989:LEU:CG	2.45	0.53
1:A:479:ALA:C	1:A:482:PRO:HD3	2.34	0.53
1:A:431:GLN:HA	1:A:484:TYR:HA	1.92	0.52
1:B:347:LEU:HD21	1:B:425:LEU:HD11	1.91	0.52
1:B:483:VAL:O	1:B:483:VAL:HG23	2.09	0.51
1:A:380:SER:HB3	1:A:383:ASP:HB2	1.92	0.51
1:A:347:LEU:HD21	1:A:425:LEU:HD11	1.91	0.51
1:A:162:TRP:HE1	1:A:285:MET:HE2	1.76	0.51
1:B:162:TRP:HE1	1:B:285:MET:HE2	1.75	0.50
1:A:899:ILE:HG12	1:A:909:HIS:CE1	2.47	0.50
1:A:491:ILE:HG21	1:A:565:LEU:HD22	1.94	0.50
1:B:491:ILE:HG21	1:B:565:LEU:HD22	1.94	0.49
1:B:911:ASP:O	1:B:911:ASP:OD1	2.30	0.49
1:B:753:ASP:HA	1:B:757:LYS:HZ3	1.76	0.48
1:B:886:THR:HA	1:B:891:ILE:HD12	1.96	0.47
1:A:209:THR:HG22	1:A:257:CYS:HB3	1.96	0.47
1:A:886:THR:HA	1:A:891:ILE:HD12	1.96	0.47
1:A:263:CYS:HA	1:A:266:LEU:HD12	1.98	0.47
1:B:326:GLU:HB3	1:B:474:TYR:HB3	1.98	0.46
1:A:394:LEU:HD13	1:A:441:MET:HE1	1.97	0.46
1:A:326:GLU:HB3	1:A:474:TYR:HB3	1.98	0.45
1:B:692:LEU:HG	1:B:728:LEU:HD21	1.97	0.45
1:B:263:CYS:HA	1:B:266:LEU:HD12	1.98	0.45
1:B:209:THR:HG22	1:B:257:CYS:HB3	1.99	0.45
1:A:730:HIS:CD2	1:A:742:GLU:OE1	2.61	0.44
1:A:707:THR:OG1	1:A:709:PRO:CD	2.56	0.44
1:B:434:THR:HA	1:B:435:GLY:HA2	1.77	0.44
1:B:751:PHE:CD1	1:B:751:PHE:C	2.95	0.44
1:A:777:ILE:HB	1:A:825:ILE:HB	1.99	0.43
1:A:580:LEU:HA	1:A:583:LEU:HD12	2.01	0.43
1:A:419:ALA:HB1	1:A:441:MET:HB3	2.01	0.43
1:A:900:MET:SD	2:A:2027:VWN:H72C	2.59	0.42
1:B:691:LYS:HE3	1:B:723:THR:HB	2.01	0.42
1:B:752:MET:O	1:B:757:LYS:HA	2.18	0.42
1:B:753:ASP:CA	1:B:757:LYS:HZ3	2.28	0.42
1:B:580:LEU:HA	1:B:583:LEU:HD12	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:386:ARG:HG3	1:A:387:MET:HE2	2.01	0.42
1:B:759:LEU:HB2	1:B:778:PHE:HB3	2.02	0.42
1:B:777:ILE:HB	1:B:825:ILE:HB	2.02	0.42
1:B:900:MET:SD	2:B:2027:VWN:H72C	2.60	0.42
1:B:971:GLY:HA3	1:B:1004:LEU:HD11	2.02	0.42
1:B:386:ARG:HG3	1:B:387:MET:HE2	2.01	0.41
1:B:583:LEU:HD11	1:B:600:LEU:HD11	2.03	0.41
1:B:328:ILE:HD11	1:B:474:TYR:HB2	2.02	0.41
1:A:730:HIS:H	1:A:742:GLU:HA	1.86	0.41
1:A:971:GLY:HA3	1:A:1004:LEU:HD11	2.02	0.41
1:B:750:THR:OG1	1:B:751:PHE:N	2.54	0.41
1:A:583:LEU:HD11	1:A:600:LEU:HD11	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	741/934 (79%)	718 (97%)	23 (3%)	0	100	100
1	B	770/934 (82%)	746 (97%)	24 (3%)	0	100	100
All	All	1511/1868 (81%)	1464 (97%)	47 (3%)	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	674/823 (82%)	628 (93%)	46 (7%)	14	42
1	B	700/823 (85%)	656 (94%)	44 (6%)	16	45
All	All	1374/1646 (84%)	1284 (93%)	90 (7%)	15	43

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	112	LEU
1	A	119	LEU
1	A	138	ASP
1	A	166	SER
1	A	190	LEU
1	A	191	LEU
1	A	221	LEU
1	A	224	LYS
1	A	270	LEU
1	A	332	LYS
1	A	333	VAL
1	A	347	LEU
1	A	381	VAL
1	A	383	ASP
1	A	394	LEU
1	A	423	LEU
1	A	434	THR
1	A	452	LEU
1	A	466	GLU
1	A	471	LEU
1	A	483	VAL
1	A	489	GLU
1	A	493	GLU
1	A	530	LEU
1	A	539	GLN
1	A	565	LEU
1	A	598	LYS
1	A	617	GLN
1	A	641	ARG
1	A	669	GLU
1	A	675	SER
1	A	687	GLU
1	A	698	PHE
1	A	706	THR

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Mol	Chain	Res	Type
1	A	707	THR
1	A	723	THR
1	A	731	LEU
1	A	744	VAL
1	A	755	LYS
1	A	791	LEU
1	A	856	LEU
1	A	895	HIS
1	A	911	ASP
1	A	934	LEU
1	A	986	LEU
1	A	1014	HIS
1	B	111	LYS
1	B	112	LEU
1	B	119	LEU
1	B	121	ILE
1	B	138	ASP
1	B	166	SER
1	B	190	LEU
1	B	191	LEU
1	B	221	LEU
1	B	224	LYS
1	B	270	LEU
1	B	332	LYS
1	B	333	VAL
1	B	341	LEU
1	B	347	LEU
1	B	380	SER
1	B	383	ASP
1	B	394	LEU
1	B	423	LEU
1	B	434	THR
1	B	466	GLU
1	B	471	LEU
1	B	483	VAL
1	B	489	GLU
1	B	493	GLU
1	B	530	LEU
1	B	539	GLN
1	B	565	LEU
1	B	598	LYS
1	B	617	GLN

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Mol	Chain	Res	Type
1	B	669	GLU
1	B	675	SER
1	B	687	GLU
1	B	723	THR
1	B	731	LEU
1	B	744	VAL
1	B	755	LYS
1	B	791	LEU
1	B	856	LEU
1	B	895	HIS
1	B	934	LEU
1	B	955	ARG
1	B	986	LEU
1	B	1014	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	116	GLN
1	A	126	HIS
1	A	193	ASN
1	A	242	GLN
1	A	258	HIS
1	A	422	ASN
1	A	464	ASN
1	A	536	HIS
1	A	563	GLN
1	A	610	GLN
1	A	730	HIS
1	A	732	GLN
1	A	773	ASN
1	A	786	GLN
1	A	795	GLN
1	A	857	ASN
1	A	906	GLN
1	B	126	HIS
1	B	193	ASN
1	B	242	GLN
1	B	244	ASN
1	B	258	HIS
1	B	422	ASN
1	B	464	ASN

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Mol	Chain	Res	Type
1	B	536	HIS
1	B	563	GLN
1	B	610	GLN
1	B	773	ASN
1	B	786	GLN
1	B	795	GLN
1	B	906	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](#)

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	$\# Z > 2$
2	VWN	B	2027	-	25,27,27	0.83	2 (8%)	32,38,38	1.12	4 (12%)
2	VWN	A	2027	-	25,27,27	0.82	2 (8%)	32,38,38	1.12	4 (12%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.

'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	VWN	B	2027	-	-	0/6/16/16	0/4/4/4
2	VWN	A	2027	-	-	0/6/16/16	0/4/4/4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2027	VWN	C1-N2	2.51	1.42	1.37
2	A	2027	VWN	C1-N2	2.49	1.42	1.37
2	B	2027	VWN	C14-N16	2.21	1.40	1.36
2	A	2027	VWN	C14-N16	2.17	1.40	1.36

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2027	VWN	C11-C12-C1	3.44	121.44	119.28
2	B	2027	VWN	C11-C12-C1	3.38	121.40	119.28
2	A	2027	VWN	C12-C1-N8	-2.91	120.90	122.86
2	B	2027	VWN	C12-C1-N8	-2.89	120.91	122.86
2	B	2027	VWN	C11-N10-C9	-2.75	121.16	123.42
2	A	2027	VWN	C11-N10-C9	-2.72	121.19	123.42
2	B	2027	VWN	C24-N16-C14	2.67	130.60	127.26
2	A	2027	VWN	C24-N16-C14	2.64	130.56	127.26

There are no chirality outliers.

There are no torsion outliers.

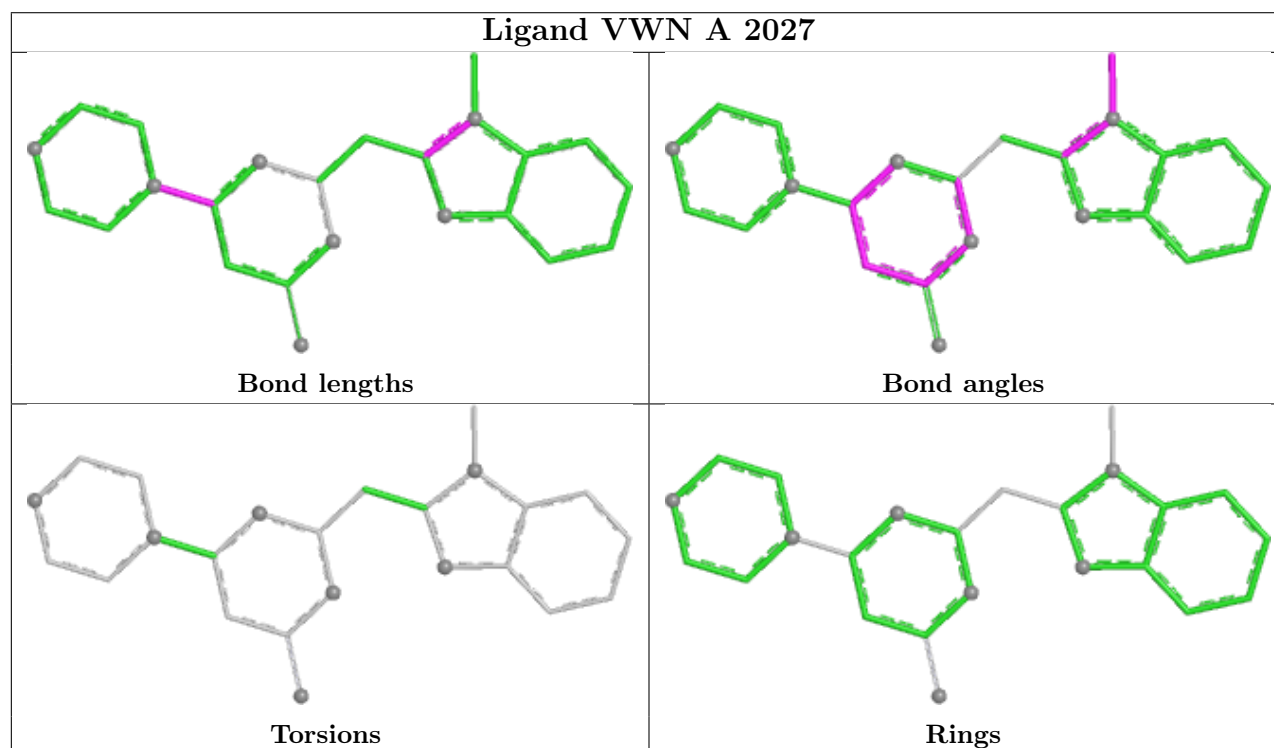
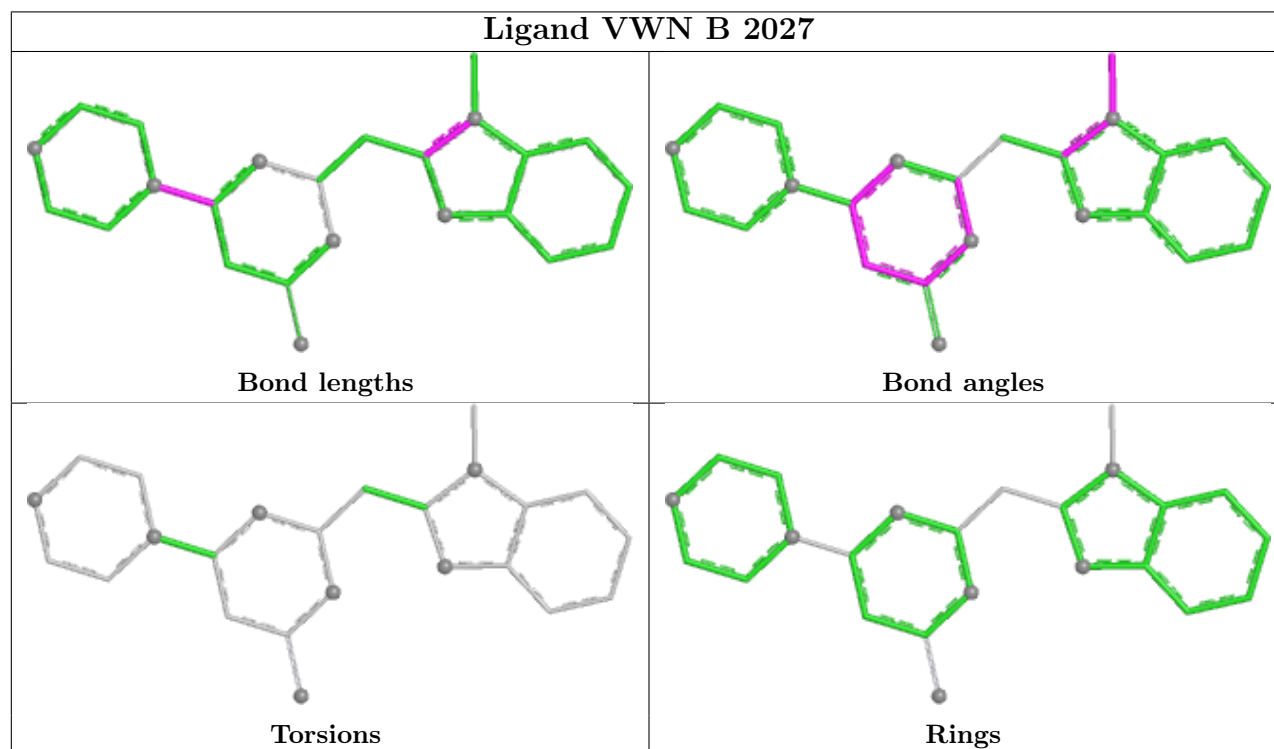
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2027	VWN	1	0
2	A	2027	VWN	1	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	765/934 (81%)	-0.84	0	100 100	29, 71, 103, 124	5 (0%)
1	B	792/934 (84%)	-0.87	1 (0%)	92 90	28, 73, 102, 129	5 (0%)
All	All	1557/1868 (83%)	-0.86	1 (0%)	92 90	28, 72, 103, 129	10 (0%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	268	SER	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.

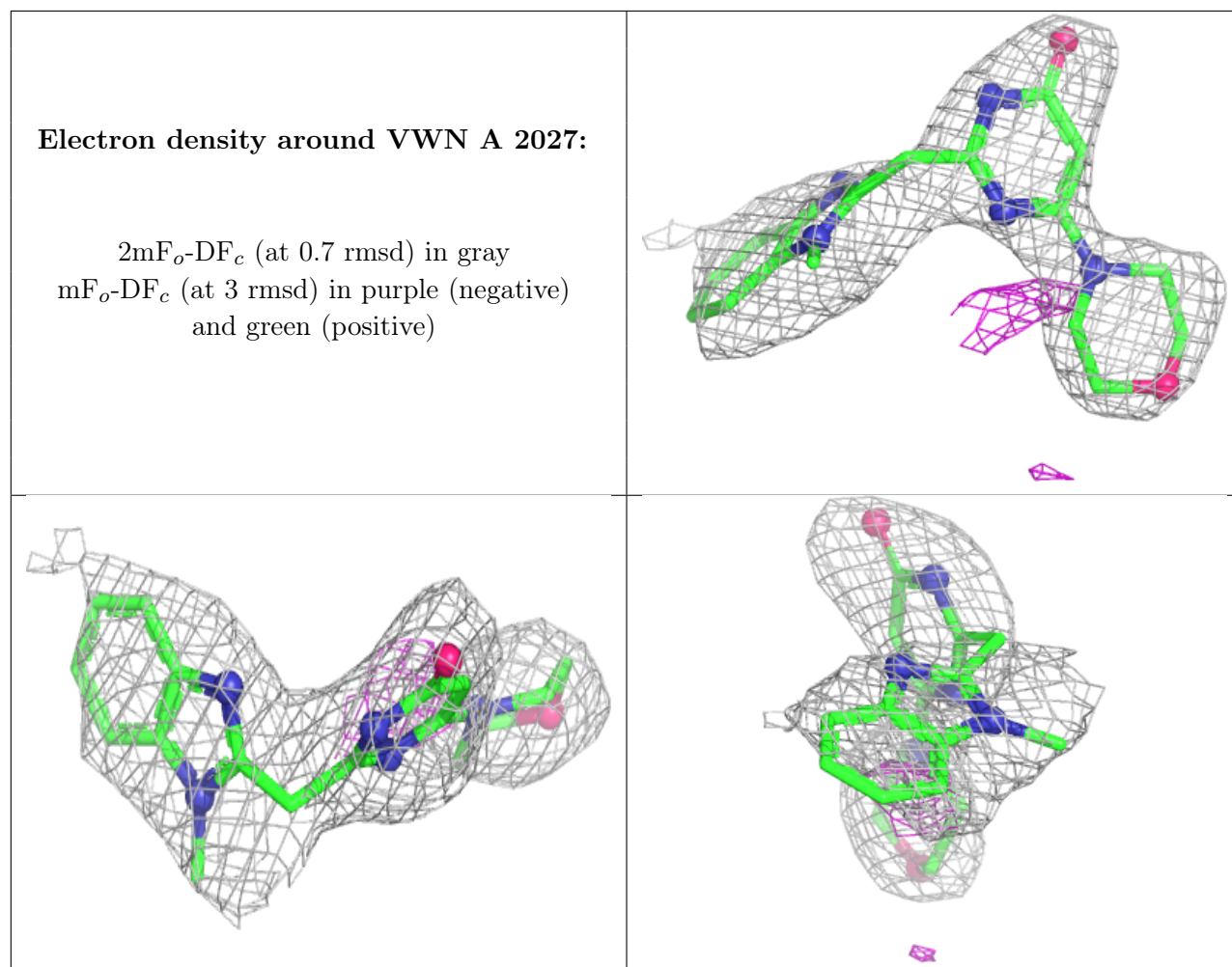
6.4 Ligands [i](#)

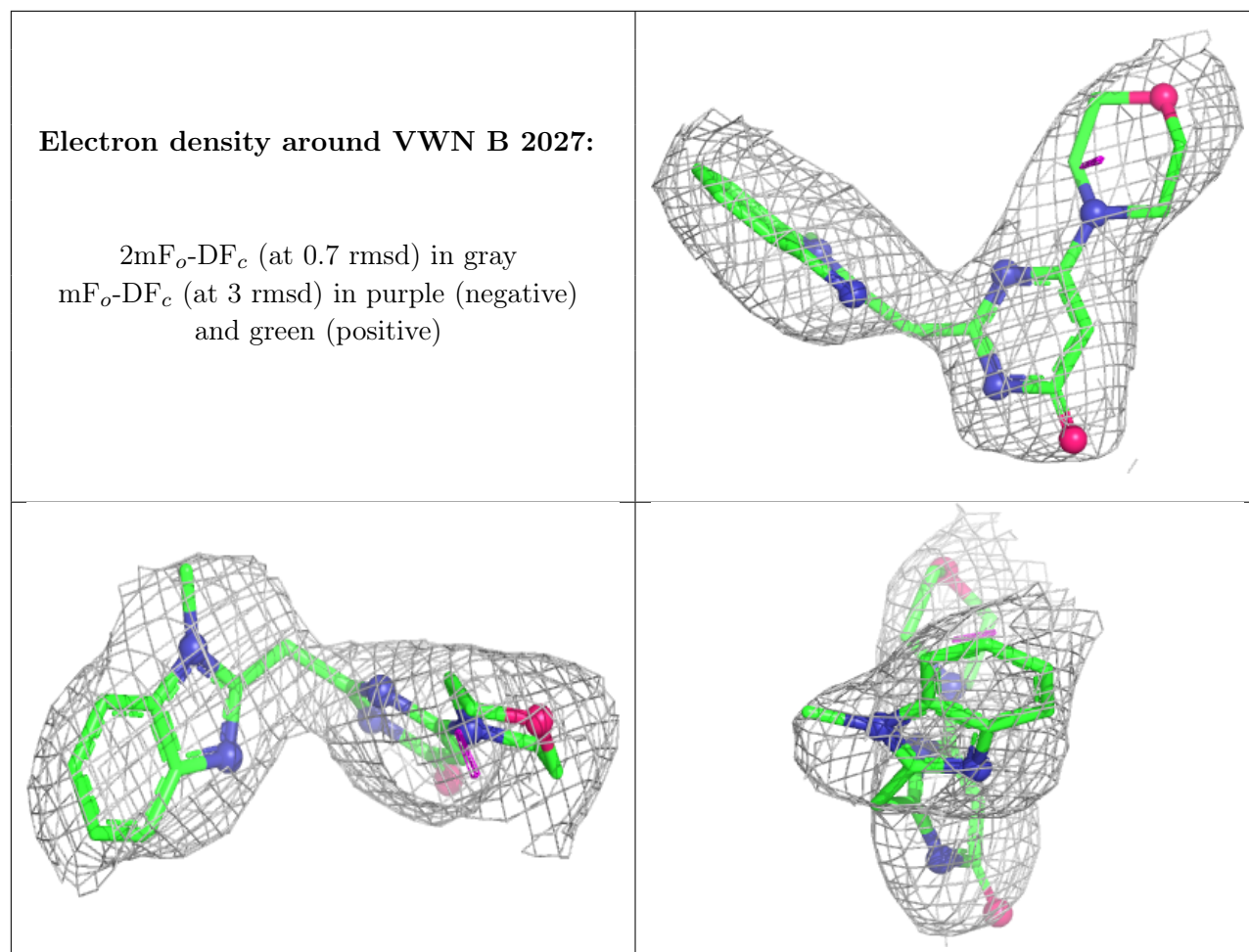
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q < 0.9
2	VWN	A	2027	24/24	0.98	0.07	80,81,81,81	0
2	VWN	B	2027	24/24	0.99	0.06	72,73,73,74	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.