



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:19 AM UTC

PDB ID : 3EFL / pdb_00003efl
Title : Crystal structure of the VEGFR2 kinase domain in complex with motesanib
Authors : Kim, J.L.; Whittington, D.A.; Long, A.M.; Rose, P.; Gu, Y.; Zhao, H.
Deposited on : 2008-09-09
Resolution : 2.20 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

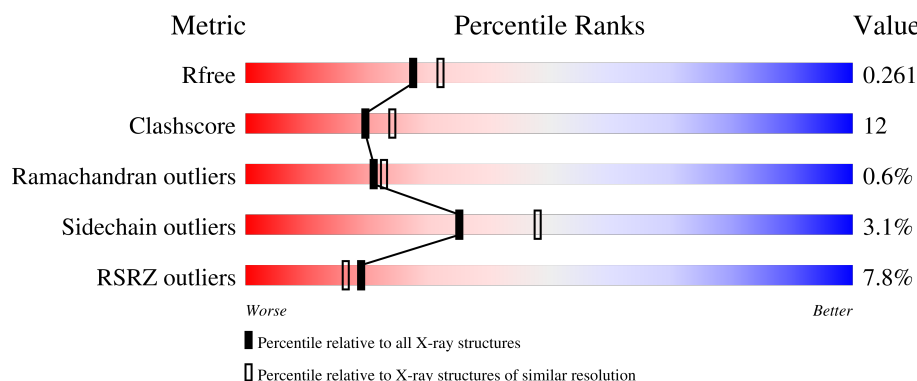
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	314	4% 71% 18% • 9%
1	B	314	9% 61% 19% • 16%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4623 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.

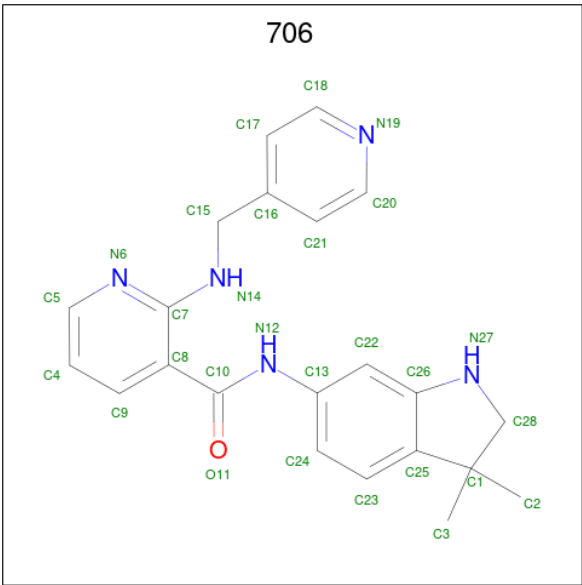
- Molecule 1 is a protein called Vascular endothelial growth factor receptor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	286	Total	C	N	O	S	0	0	0
			2315	1482	409	408	16			
1	B	263	Total	C	N	O	S	0	0	0
			2110	1360	364	370	16			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	817	ALA	CYS	engineered mutation	UNP P35968
A	990	VAL	GLU	engineered mutation	UNP P35968
A	1172	ARG	-	expression tag	UNP P35968
A	1173	HIS	-	expression tag	UNP P35968
A	1174	HIS	-	expression tag	UNP P35968
A	1175	HIS	-	expression tag	UNP P35968
A	1176	HIS	-	expression tag	UNP P35968
A	1177	HIS	-	expression tag	UNP P35968
A	1178	HIS	-	expression tag	UNP P35968
B	817	ALA	CYS	engineered mutation	UNP P35968
B	990	VAL	GLU	engineered mutation	UNP P35968
B	1172	ARG	-	expression tag	UNP P35968
B	1173	HIS	-	expression tag	UNP P35968
B	1174	HIS	-	expression tag	UNP P35968
B	1175	HIS	-	expression tag	UNP P35968
B	1176	HIS	-	expression tag	UNP P35968
B	1177	HIS	-	expression tag	UNP P35968
B	1178	HIS	-	expression tag	UNP P35968

- Molecule 2 is N-(3,3-dimethyl-2,3-dihydro-1H-indol-6-yl)-2-[(pyridin-4-ylmethyl)amino]pyridine-3-carboxamide (CCD ID: 706) (formula: C₂₂H₂₃N₅O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			28	22	5	1		
2	B	1	Total	C	N	O	0	0
			28	22	5	1		

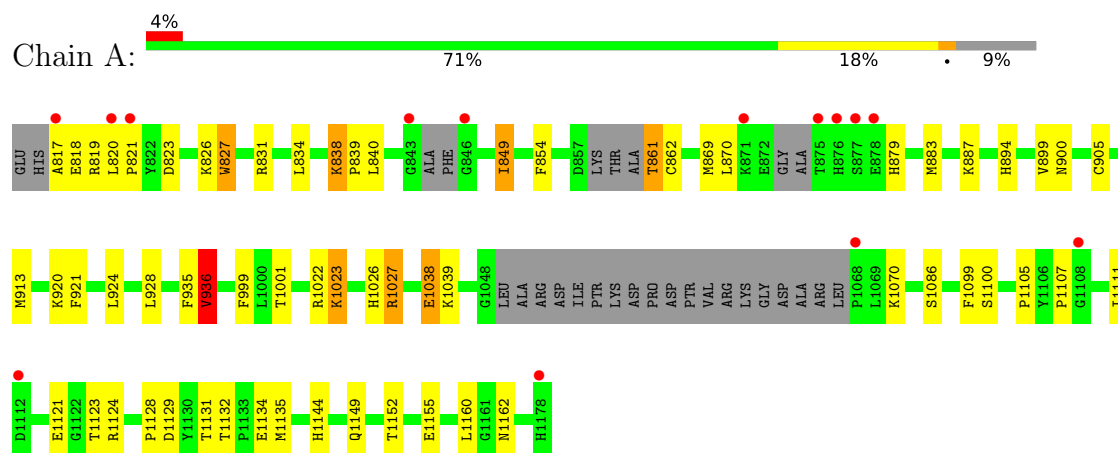
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	100	Total	O	0	0
			100	100		
3	B	42	Total	O	0	0
			42	42		

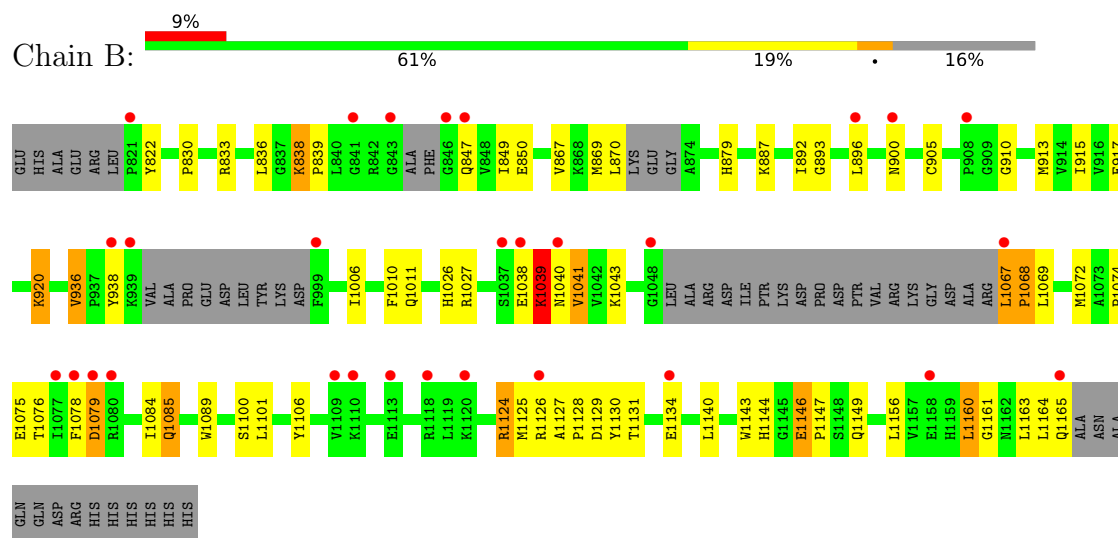
3 Residue-property plots [i](#)

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: Vascular endothelial growth factor receptor 2



- Molecule 1: Vascular endothelial growth factor receptor 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.59Å 66.81Å 91.42Å 90.00° 93.27° 90.00°	Depositor
Resolution (Å)	38.04 – 2.20 38.04 – 2.20	Depositor EDS
% Data completeness (in resolution range)	87.0 (38.04-2.20) 87.0 (38.04-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.62 (at 2.10Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.230 , 0.264 0.229 , 0.261	Depositor DCC
R_{free} test set	1528 reflections (4.41%)	wwPDB-VP
Wilson B-factor (Å ²)	29.8	Xtriage
Anisotropy	0.608	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 37.2	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	0.93	EDS
Total number of atoms	4623	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.90% 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:
706

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	0.43	0/2373	0.96	12/3201 (0.4%)
1	B	0.41	0/2160	1.01	16/2913 (0.5%)
All	All	0.42	0/4533	0.99	28/6114 (0.5%)

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	936	VAL	N-CA-C	-10.80	100.33	109.19
1	B	1079	ASP	N-CA-C	8.84	124.37	113.50
1	A	827	TRP	N-CA-C	8.20	123.42	113.41
1	A	1100	SER	N-CA-C	-8.16	101.43	112.94
1	B	1100	SER	N-CA-C	-8.07	101.50	113.61
1	B	1067	LEU	CA-C-N	7.74	129.52	119.84
1	B	1067	LEU	C-N-CA	7.74	129.52	119.84
1	A	1132	THR	N-CA-C	-7.56	101.23	110.31
1	A	840	LEU	N-CA-C	-6.43	105.74	113.97
1	B	920	LYS	N-CA-C	6.28	119.79	111.75
1	A	838	LYS	CA-C-N	6.22	126.17	119.76
1	A	838	LYS	C-N-CA	6.22	126.17	119.76
1	B	1129	ASP	N-CA-C	6.10	118.73	111.71
1	B	838	LYS	CA-C-N	6.08	126.02	119.76
1	B	838	LYS	C-N-CA	6.08	126.02	119.76
1	B	910	GLY	CA-C-N	6.07	125.98	119.85
1	B	910	GLY	C-N-CA	6.07	125.98	119.85
1	B	867	VAL	N-CA-C	5.99	116.49	107.75
1	A	1086	SER	N-CA-C	-5.54	105.24	112.23
1	A	1038	GLU	N-CA-C	5.53	122.58	110.80
1	A	900	ASN	N-CA-C	5.49	117.68	108.96
1	B	1041	VAL	N-CA-C	-5.27	100.88	108.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	938	TYR	N-CA-C	5.24	117.15	109.24
1	A	1129	ASP	N-CA-C	5.24	118.93	112.54
1	B	915	ILE	N-CA-C	5.21	115.47	108.17
1	B	936	VAL	CA-C-N	5.13	126.26	119.84
1	B	936	VAL	C-N-CA	5.13	126.26	119.84
1	A	894	HIS	N-CA-C	5.02	118.25	110.17

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	2315	0	2296	47	0
1	B	2110	0	2124	58	0
2	A	28	0	23	1	0
2	B	28	0	23	0	0
3	A	100	0	0	3	0
3	B	42	0	0	2	0
All	All	4623	0	4466	103	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 12.

All (103) 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:1128:PRO:HG2	1:B:1131:THR:HB	1.63	0.81
1:A:870:LEU:HD11	1:A:879:HIS:HA	1.62	0.80
1:A:1026:HIS:O	1:A:1027:ARG:HB2	1.85	0.76
1:B:822:TYR:HB2	1:B:887:LYS:NZ	2.06	0.69
1:A:854:PHE:CE1	1:A:861:THR:HG23	2.27	0.69
1:B:870:LEU:HD11	1:B:879:HIS:HA	1.75	0.69
1:A:1023:LYS:HE3	1:A:1023:LYS:HA	1.74	0.68
1:B:1068:PRO:O	1:B:1072:MET:HG3	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1070:LYS:HE3	1:A:1111:ILE:HD11	1.75	0.67
1:B:1069:LEU:HD23	1:B:1072:MET:HE3	1.77	0.66
1:B:1069:LEU:HD23	1:B:1072:MET:CE	2.25	0.66
1:B:905:CYS:SG	1:B:913:MET:HE3	2.40	0.62
1:B:1038:GLU:C	1:B:1040:ASN:N	2.53	0.60
1:B:1010:PHE:CG	1:B:1164:LEU:HD11	2.38	0.59
1:A:905:CYS:SG	1:A:913:MET:HE3	2.44	0.58
1:A:861:THR:HG22	1:A:862:CYS:N	2.17	0.58
1:A:1162:ASN:ND2	3:A:178:HOH:O	2.37	0.58
1:B:822:TYR:HB2	1:B:887:LYS:HZ3	1.68	0.57
1:A:820:LEU:HD12	1:A:821:PRO:HD2	1.87	0.57
1:B:1161:GLY:HA2	1:B:1164:LEU:HD12	1.87	0.56
1:A:936:VAL:HG22	1:A:999:PHE:HD2	1.70	0.56
1:B:1134:GLU:HG2	1:B:1163:LEU:HD21	1.88	0.56
1:B:1160:LEU:HD12	1:B:1163:LEU:HD12	1.87	0.55
1:B:1038:GLU:C	1:B:1040:ASN:H	2.13	0.55
1:B:900:ASN:HB3	1:B:917:GLU:OE1	2.07	0.55
1:B:1085:GLN:HE21	1:B:1085:GLN:H	1.54	0.55
1:B:836:LEU:HD13	1:B:869:MET:HE1	1.88	0.54
1:B:1040:ASN:H	1:B:1040:ASN:HD22	1.56	0.54
1:B:1069:LEU:HA	1:B:1072:MET:CE	2.37	0.54
1:A:1022:ARG:HH21	1:A:1022:ARG:HG3	1.73	0.54
1:A:1152:THR:OG1	1:A:1155:GLU:HG3	2.08	0.54
1:B:1026:HIS:O	1:B:1027:ARG:HB2	2.07	0.53
1:A:1128:PRO:HG2	1:A:1131:THR:HB	1.89	0.53
1:A:869:MET:HB3	1:A:913:MET:HG2	1.89	0.53
1:A:849:ILE:H	1:A:849:ILE:HD13	1.74	0.53
1:A:899:VAL:CG1	2:A:501:706:H24	2.40	0.52
1:A:936:VAL:O	1:A:1001:THR:HA	2.10	0.51
1:B:1074:PRO:O	1:B:1078:PHE:HD1	1.94	0.51
1:B:1075:GLU:O	1:B:1079:ASP:HB2	2.10	0.51
1:A:1039:LYS:N	1:A:1039:LYS:HD2	2.26	0.51
1:B:1038:GLU:C	1:B:1039:LYS:HD2	2.37	0.50
1:A:838:LYS:HD3	1:B:1038:GLU:CD	2.37	0.50
1:B:830:PRO:HB2	1:B:833:ARG:HG3	1.93	0.49
1:A:854:PHE:HE1	1:A:861:THR:HG23	1.73	0.49
1:A:1038:GLU:HB3	1:A:1039:LYS:HD2	1.94	0.49
1:A:826:LYS:HD3	1:A:827:TRP:NE1	2.27	0.49
1:B:892:ILE:HG13	1:B:893:GLY:N	2.27	0.49
1:B:920:LYS:HG2	3:B:41:HOH:O	2.12	0.49
1:A:1070:LYS:HE3	1:A:1111:ILE:CD1	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1069:LEU:HA	1:B:1072:MET:HE2	1.96	0.48
1:A:1070:LYS:HD3	1:A:1105:PRO:HG2	1.95	0.47
1:A:1022:ARG:HG3	1:A:1022:ARG:NH2	2.29	0.47
1:B:1006:ILE:HG22	1:B:1164:LEU:CD2	2.44	0.47
1:A:1070:LYS:CE	1:A:1111:ILE:HD11	2.44	0.47
1:A:1099:PHE:CE2	1:A:1135:MET:HG2	2.49	0.47
1:A:839:PRO:HD2	1:B:1039:LYS:HE2	1.97	0.47
1:B:1038:GLU:O	1:B:1040:ASN:N	2.48	0.47
1:B:1011:GLN:OE1	1:B:1041:VAL:HG23	2.15	0.47
1:B:1126:ARG:HG3	3:B:112:HOH:O	2.15	0.46
1:A:849:ILE:HD13	1:A:849:ILE:N	2.31	0.46
1:B:1144:HIS:HE2	1:B:1149:GLN:NE2	2.14	0.46
1:A:935:PHE:CD1	1:A:935:PHE:C	2.93	0.46
1:B:1126:ARG:HG2	1:B:1127:ALA:N	2.30	0.46
1:B:1040:ASN:N	1:B:1040:ASN:HD22	2.14	0.46
1:B:1085:GLN:H	1:B:1085:GLN:NE2	2.14	0.46
1:B:1101:LEU:HD12	1:B:1130:TYR:CZ	2.51	0.46
1:A:817:ALA:C	1:A:819:ARG:N	2.73	0.46
1:A:818:GLU:HA	1:A:887:LYS:HD2	1.98	0.46
1:A:1134:GLU:H	1:A:1134:GLU:CD	2.24	0.45
1:A:883:MET:CE	1:A:887:LYS:HE3	2.46	0.45
1:A:920:LYS:HE3	1:A:921:PHE:CZ	2.52	0.45
1:A:831:ARG:HG2	1:A:913:MET:CE	2.47	0.45
1:B:838:LYS:HA	1:B:839:PRO:HD3	1.81	0.45
1:B:1160:LEU:O	1:B:1164:LEU:HG	2.17	0.44
1:B:917:GLU:OE2	1:B:1043:LYS:HD3	2.17	0.44
1:B:1069:LEU:HA	1:B:1072:MET:HE3	1.99	0.44
1:B:1156:LEU:O	1:B:1160:LEU:HB2	2.17	0.44
1:A:817:ALA:C	1:A:819:ARG:H	2.26	0.44
1:A:839:PRO:HA	1:A:849:ILE:HG22	2.00	0.44
1:A:924:LEU:O	1:A:928:LEU:HG	2.18	0.44
1:A:817:ALA:O	1:A:820:LEU:N	2.51	0.43
1:B:1067:LEU:HA	1:B:1068:PRO:HD2	1.75	0.43
1:A:1123:THR:O	1:A:1124:ARG:HD3	2.18	0.43
1:B:1011:GLN:OE1	1:B:1041:VAL:HA	2.18	0.43
1:A:1038:GLU:HA	1:A:1038:GLU:OE2	2.19	0.43
1:A:1107:PRO:HB3	3:A:159:HOH:O	2.18	0.43
1:B:1144:HIS:NE2	1:B:1149:GLN:NE2	2.66	0.43
1:B:1165:GLN:O	1:B:1165:GLN:HG2	2.19	0.42
1:B:822:TYR:HB2	1:B:887:LYS:HZ1	1.84	0.42
1:B:849:ILE:HG22	1:B:850:GLU:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1084:ILE:HG13	1:B:1085:GLN:N	2.33	0.41
1:B:1106:TYR:CE1	1:B:1125:MET:HG3	2.56	0.41
1:B:1130:TYR:CD1	1:B:1130:TYR:N	2.88	0.41
1:A:920:LYS:HG2	3:A:52:HOH:O	2.20	0.41
1:A:1144:HIS:CE1	1:A:1149:GLN:HG3	2.56	0.41
1:B:1124:ARG:HG2	1:B:1143:TRP:HE3	1.86	0.41
1:A:823:ASP:OD2	1:A:826:LYS:HB2	2.20	0.41
1:B:896:LEU:HD12	1:B:896:LEU:O	2.21	0.41
1:B:1075:GLU:HG2	1:B:1076:THR:N	2.35	0.41
1:B:1146:GLU:HA	1:B:1147:PRO:HD3	1.89	0.41
1:A:1026:HIS:O	1:A:1027:ARG:CB	2.62	0.40
1:B:1089:TRP:CD1	1:B:1089:TRP:C	2.99	0.40
1:B:836:LEU:CD1	1:B:869:MET:HE1	2.51	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	276/314 (88%)	266 (96%)	9 (3%)	1 (0%)	30	34
1	B	253/314 (81%)	242 (96%)	9 (4%)	2 (1%)	16	16
All	All	529/628 (84%)	508 (96%)	18 (3%)	3 (1%)	21	23

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1039	LYS
1	A	1027	ARG
1	B	1068	PRO

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	251/270 (93%)	244 (97%)	7 (3%)	38	52
1	B	230/270 (85%)	222 (96%)	8 (4%)	32	43
All	All	481/540 (89%)	466 (97%)	15 (3%)	35	48

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	834	LEU
1	A	849	ILE
1	A	861	THR
1	A	936	VAL
1	A	1023	LYS
1	A	1121	GLU
1	A	1160	LEU
1	B	847	GLN
1	B	936	VAL
1	B	1039	LYS
1	B	1085	GLN
1	B	1124	ARG
1	B	1140	LEU
1	B	1146	GLU
1	B	1160	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	933	ASN
1	A	1149	GLN
1	A	1159	HIS
1	A	1162	ASN
1	A	1165	GLN
1	A	1169	GLN
1	B	847	GLN

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Mol	Chain	Res	Type
1	B	894	HIS
1	B	1040	ASN
1	B	1085	GLN
1	B	1149	GLN
1	B	1165	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	706	B	502	-	30,31,31	1.09	3 (10%)	41,44,44	1.27	6 (14%)
2	706	A	501	-	30,31,31	1.11	3 (10%)	41,44,44	1.29	6 (14%)

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	706	B	502	-	-	0/13/25/25	0/4/4/4
2	706	A	501	-	-	1/13/25/25	0/4/4/4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	706	C13-N12	-4.02	1.33	1.41
2	B	502	706	C13-N12	-3.83	1.33	1.41
2	A	501	706	C8-C7	-2.33	1.39	1.42
2	B	502	706	C8-C7	-2.28	1.39	1.42
2	A	501	706	C26-N27	-2.26	1.33	1.38
2	B	502	706	C26-N27	-2.24	1.33	1.38

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	A	501	706	C8-C7-N14	-3.35	119.81	121.81
2	A	501	706	C5-N6-C7	2.94	122.45	116.74
2	B	502	706	C5-N6-C7	2.93	122.45	116.74
2	B	502	706	C8-C7-N14	-2.85	120.11	121.81
2	B	502	706	C23-C25-C1	2.64	131.85	125.15
2	A	501	706	C20-N19-C18	2.48	122.54	116.86
2	A	501	706	C23-C25-C1	2.46	131.39	125.15
2	B	502	706	C20-N19-C18	2.42	122.39	116.86
2	B	502	706	C3-C1-C2	2.12	111.90	109.17
2	B	502	706	C4-C5-N6	-2.09	120.11	123.42
2	A	501	706	C9-C8-C7	2.03	118.89	117.47
2	A	501	706	C3-C1-C2	2.01	111.76	109.17

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	706	C24-C13-N12-C10

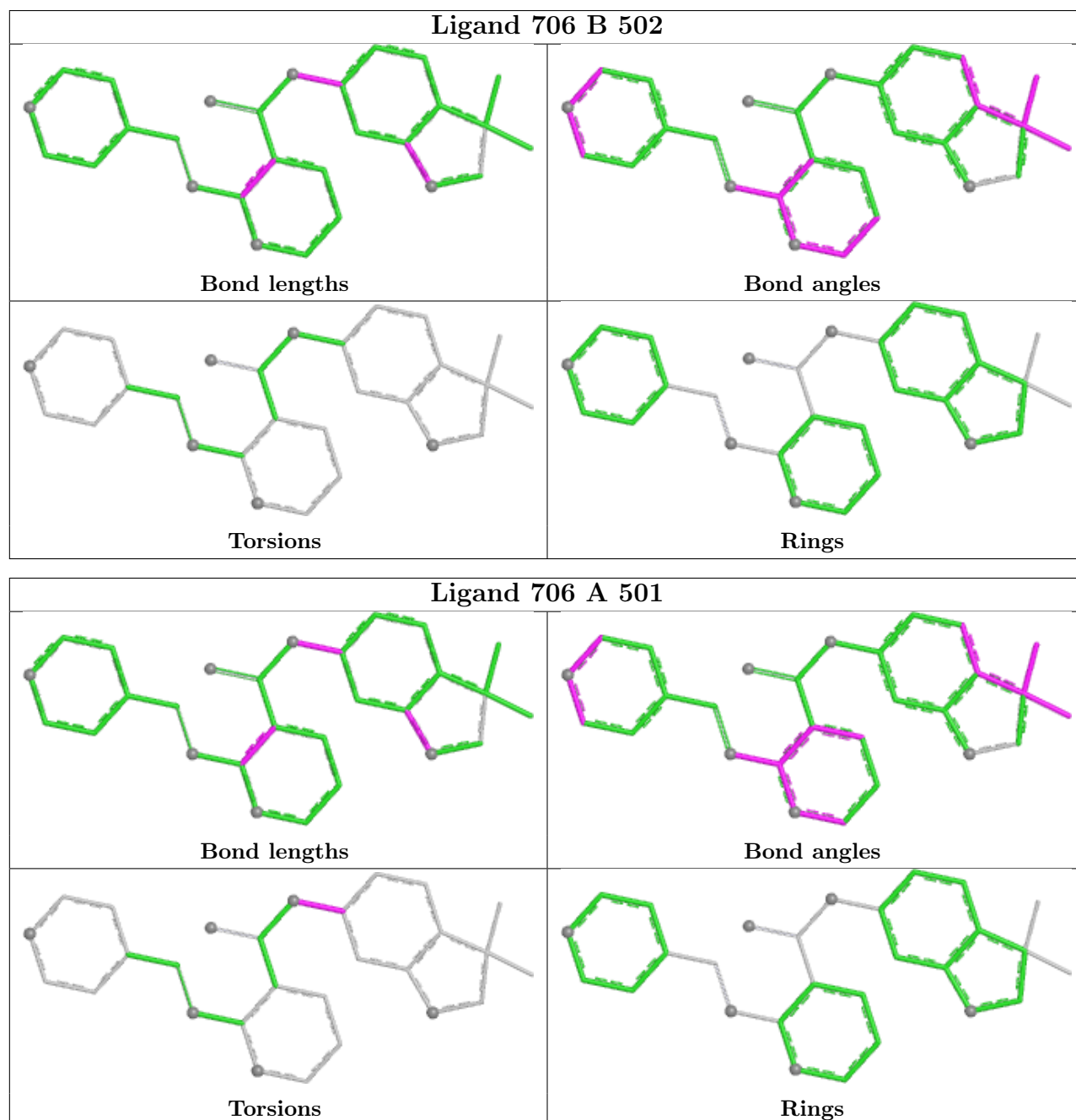
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	706	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

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	286/314 (91%)	0.31	14 (4%) 35 31	17, 33, 66, 87	0
1	B	263/314 (83%)	0.76	29 (11%) 10 8	24, 45, 73, 85	0
All	All	549/628 (87%)	0.53	43 (7%) 19 16	17, 38, 70, 87	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1079	ASP	4.6
1	B	999	PHE	3.9
1	A	876	HIS	3.7
1	A	1178	HIS	3.5
1	B	841	GLY	3.5
1	B	1078	PHE	3.4
1	B	1134	GLU	3.4
1	B	1048	GLY	3.4
1	A	871	LYS	3.2
1	B	1165	GLN	3.1
1	A	875	THR	3.0
1	A	846	GLY	3.0
1	B	1067	LEU	3.0
1	B	1037	SER	3.0
1	B	846	GLY	3.0
1	A	1068	PRO	2.9
1	A	820	LEU	2.8
1	A	817	ALA	2.8
1	B	1118	ARG	2.7
1	A	843	GLY	2.6
1	B	847	GLN	2.6
1	A	1108	GLY	2.6
1	B	843	GLY	2.6
1	B	900	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	938	TYR	2.5
1	B	1126	ARG	2.5
1	A	1112	ASP	2.5
1	A	878	GLU	2.4
1	A	821	PRO	2.4
1	A	877	SER	2.4
1	B	1158	GLU	2.4
1	B	1040	ASN	2.3
1	B	1120	LYS	2.3
1	B	1038	GLU	2.2
1	B	939	LYS	2.2
1	B	1113	GLU	2.2
1	B	896	LEU	2.1
1	B	1110	LYS	2.1
1	B	1077	ILE	2.1
1	B	1109	VAL	2.1
1	B	1080	ARG	2.1
1	B	908	PRO	2.1
1	B	821	PRO	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

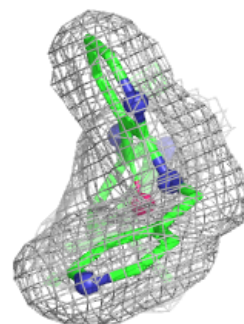
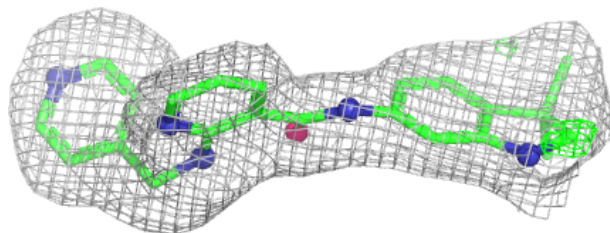
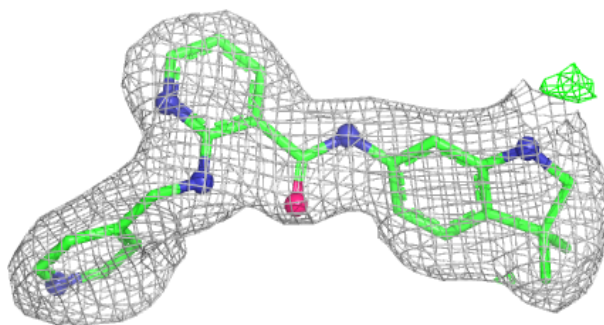
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	706	A	501	28/28	0.94	0.06	20,25,30,30	0
2	706	B	502	28/28	0.96	0.06	20,23,33,34	0

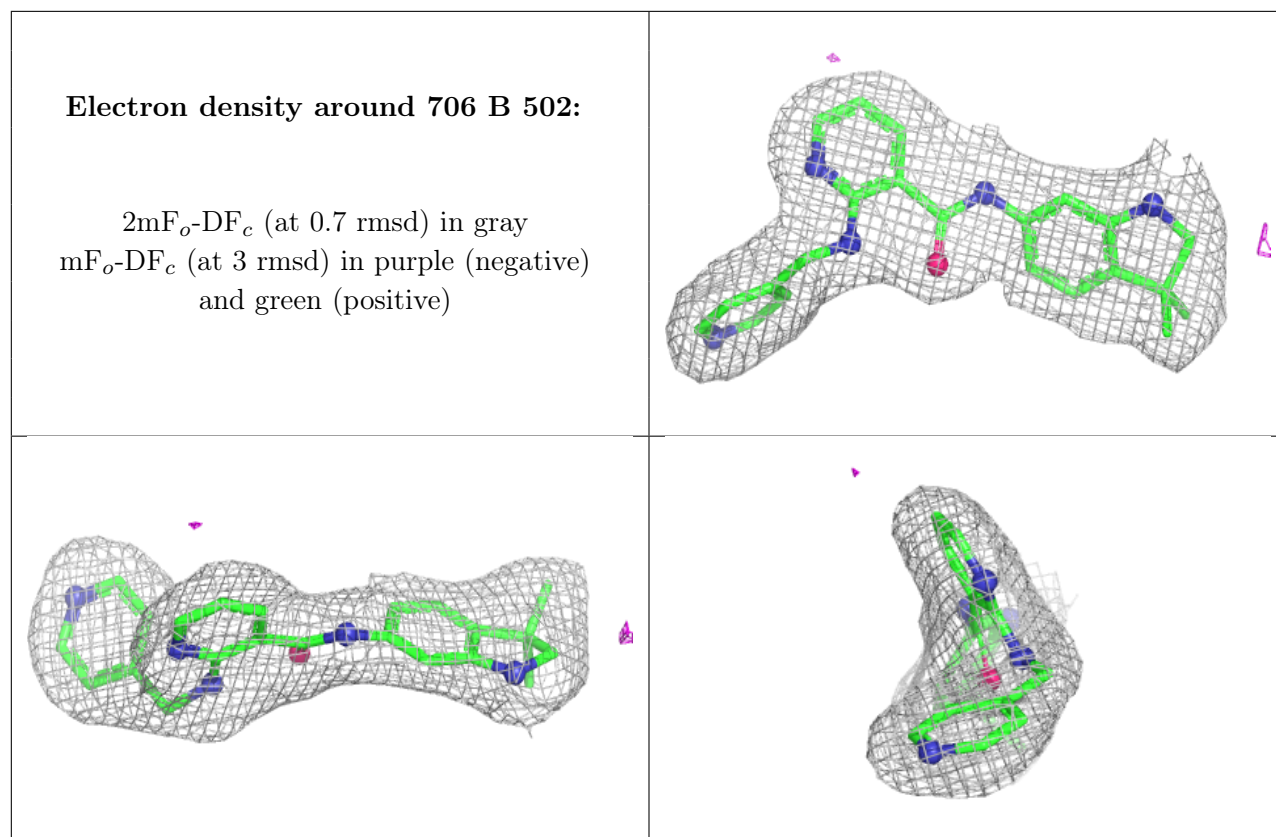
The following is a graphical depiction of the model fit to experimental electron density of all

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 706 A 501:

$2mF_o-DF_c$ (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.