



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2022 – 12:11 PM JST

PDB ID : 7VDQ
Title : The structure of cyclin-dependent kinase 5 (CDK5) in complex with p25 and Compound 7
Authors : Malojcic, G.; Clugston, S.L.; Daniels, M.; Harmange, J.C.; Ledeborer, M.
Deposited on : 2021-09-07
Resolution : 2.91 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 : 2.27
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) : 2.27

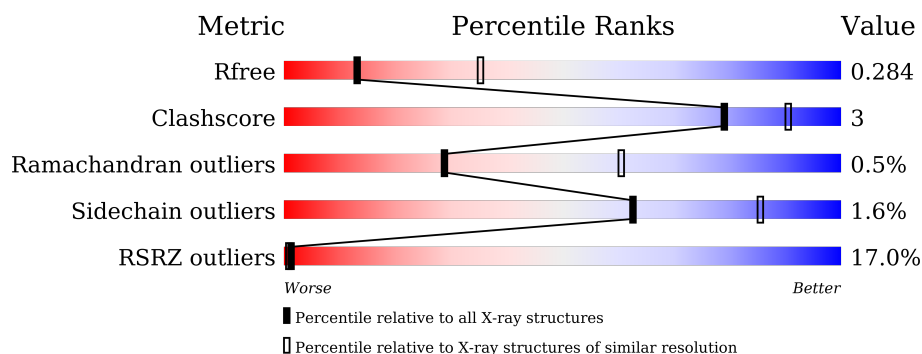
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.91 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	2307 (2.94-2.90)
Clashscore	141614	2531 (2.94-2.90)
Ramachandran outliers	138981	2462 (2.94-2.90)
Sidechain outliers	138945	2464 (2.94-2.90)
RSRZ outliers	127900	2248 (2.94-2.90)

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	292	<div> <div>2%</div> <div>92%</div> <div>7%</div> </div>
1	B	292	<div> <div>26%</div> <div>85%</div> <div>10%</div> <div>5%</div> </div>
2	C	209	<div> <div>64%</div> <div>7%</div> <div>29%</div> </div>
2	D	209	<div> <div>31%</div> <div>57%</div> <div>8%</div> <div>35%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6944 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 Cyclin-dependent-like kinase 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	290	Total	C	N	O	S	0	0	0
			2330	1495	400	424	11			
1	B	277	Total	C	N	O	S	0	0	0
			2233	1433	385	405	10			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	SER	-	expression tag	UNP Q00535
B	1	SER	-	expression tag	UNP Q00535

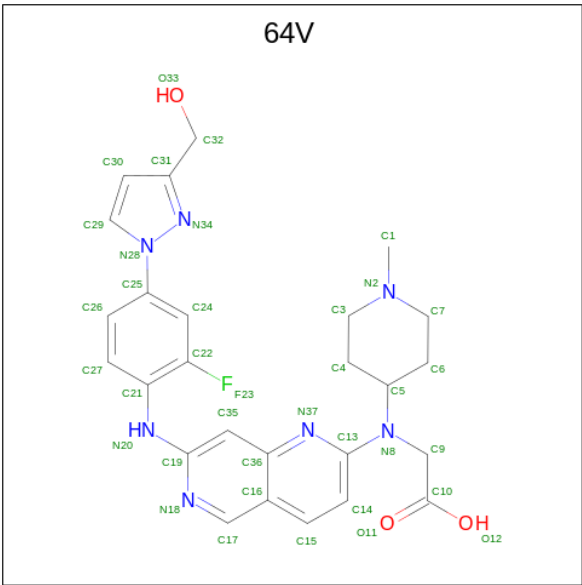
- Molecule 2 is a protein called Cyclin-dependent kinase 5 activator 1, p25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	148	Total	C	N	O	S	0	0	0
			1196	768	197	220	11			
2	D	136	Total	C	N	O	S	0	0	0
			1104	711	179	204	10			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	99	MET	-	initiating methionine	UNP Q15078
D	99	MET	-	initiating methionine	UNP Q15078

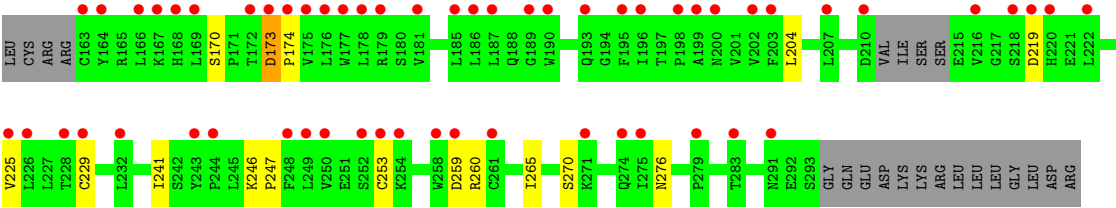
- Molecule 3 is 2-[[7-[[2-fluoranyl-4-[3-(hydroxymethyl)pyrazol-1-yl]phenyl]amino]-1,6-naphthylidin-2-yl]-(1-methylpiperidin-4-yl)amino]ethanoic acid (three-letter code: 64V) (formula: C₂₆H₂₈FN₇O₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			37	26	1	7	3		
3	B	1	Total	C	F	N	O	0	0
			37	26	1	7	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	4	Total	O	0	0
			4	4		
4	B	1	Total	O	0	0
			1	1		
4	C	2	Total	O	0	0
			2	2		



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	117.61Å 117.61Å 154.01Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.50 – 2.91 38.50 – 2.91	Depositor EDS
% Data completeness (in resolution range)	99.8 (38.50-2.91) 99.9 (38.50-2.91)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.98 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.252 , 0.284 0.253 , 0.284	Depositor DCC
R_{free} test set	1392 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	78.1	Xtriage
Anisotropy	0.177	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 52.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.028 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6944	wwPDB-VP
Average B, all atoms (Å ²)	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 64V

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	0.65	0/2385	0.78	0/3229
1	B	0.67	0/2284	0.77	0/3091
2	C	0.65	0/1224	0.74	0/1660
2	D	0.67	0/1129	0.70	0/1530
All	All	0.66	0/7022	0.76	0/9510

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 [i](#)

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	2330	0	2344	13	0
1	B	2233	0	2252	18	0
2	C	1196	0	1185	7	0
2	D	1104	0	1088	8	0
3	A	37	0	0	0	0
3	B	37	0	0	0	0
4	A	4	0	0	0	0
4	B	1	0	0	0	0
4	C	2	0	0	0	0
All	All	6944	0	6869	38	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 (38) 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:104:LYS:NZ	1:A:285:TYR:O	2.21	0.73
1:A:152:GLY:HA3	2:C:276:ASN:O	1.90	0.71
1:A:50:ARG:NH2	2:C:235:SER:O	2.25	0.64
1:B:247:SER:O	1:B:248:LEU:HB2	1.97	0.64
1:B:50:ARG:NH2	2:D:241:ILE:O	2.32	0.62
2:D:204:LEU:HD12	2:D:229:CYS:SG	2.40	0.62
1:A:43:GLY:O	1:A:46:SER:HB3	2.02	0.60
1:B:109:GLN:OE1	1:B:140:LEU:N	2.35	0.56
1:A:34:ARG:NH2	1:B:41:ASP:OD2	2.37	0.55
1:B:44:VAL:N	1:B:45:PRO:HD2	2.24	0.53
2:C:246:LYS:HB3	2:C:247:PRO:HD3	1.90	0.53
2:C:178:LEU:HD22	2:C:230:LEU:HD13	1.91	0.52
1:B:38:ASP:OD1	1:B:40:ASP:HB2	2.10	0.51
2:C:173:ASP:HB2	2:C:174:PRO:HD3	1.93	0.51
1:B:157:CYS:HB2	1:B:177:LYS:HB3	1.93	0.50
1:B:152:GLY:HA3	2:D:276:ASN:O	2.12	0.48
1:A:53:CYS:SG	2:C:265:ILE:HG12	2.54	0.48
1:A:107:LEU:HD22	1:A:192:PHE:CD1	2.49	0.47
1:B:55:LEU:HD21	1:B:145:PHE:CD1	2.49	0.47
1:B:157:CYS:HB2	1:B:177:LYS:CB	2.45	0.47
1:B:222:PRO:HG3	1:B:269:CYS:SG	2.54	0.47
2:D:246:LYS:HB3	2:D:247:PRO:HD3	1.97	0.47
1:A:44:VAL:HB	1:A:45:PRO:HD3	1.96	0.46
1:B:166:TRP:CD1	1:B:204:PRO:HA	2.51	0.46
1:B:107:LEU:HD22	1:B:192:PHE:CD1	2.52	0.45
1:A:166:TRP:CD1	1:A:204:PRO:HA	2.51	0.45
2:C:186:LEU:C	2:C:186:LEU:HD13	2.37	0.45
1:A:166:TRP:CZ3	1:B:92:ASP:HB3	2.53	0.44
1:B:36:ARG:C	1:B:37:LEU:HD12	2.38	0.44
1:B:49:LEU:HA	1:B:52:ILE:HD12	2.01	0.43
2:D:253:CYS:O	2:D:253:CYS:SG	2.76	0.43
1:A:174:PHE:CD2	1:A:236:TYR:HA	2.54	0.43
1:B:50:ARG:HD2	1:B:148:ALA:O	2.20	0.42
1:A:55:LEU:HD21	1:A:145:PHE:CD1	2.55	0.42
1:A:78:LEU:HD12	1:A:78:LEU:N	2.36	0.41
2:D:173:ASP:H	2:D:174:PRO:HD2	1.86	0.41
2:D:225:VAL:HG22	2:D:260:ARG:NH1	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:49:LEU:HB2	2:D:265:ILE:HD11	2.03	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	288/292 (99%)	267 (93%)	20 (7%)	1 (0%)	41	70
1	B	273/292 (94%)	249 (91%)	22 (8%)	2 (1%)	22	53
2	C	146/209 (70%)	144 (99%)	2 (1%)	0	100	100
2	D	130/209 (62%)	124 (95%)	5 (4%)	1 (1%)	19	49
All	All	837/1002 (84%)	784 (94%)	49 (6%)	4 (0%)	29	60

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	163	VAL
1	B	163	VAL
1	B	248	LEU
2	D	173	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	258/260 (99%)	257 (100%)	1 (0%)	91	97
1	B	248/260 (95%)	247 (100%)	1 (0%)	91	97
2	C	138/187 (74%)	133 (96%)	5 (4%)	35	67
2	D	128/187 (68%)	123 (96%)	5 (4%)	32	64
All	All	772/894 (86%)	760 (98%)	12 (2%)	62	85

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	251	VAL
1	B	207	ASP
2	C	147	SER
2	C	149	SER
2	C	242	SER
2	C	254	LYS
2	C	270	SER
2	D	154	CYS
2	D	170	SER
2	D	219	ASP
2	D	259	ASP
2	D	270	SER

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	273	GLN
1	A	282	GLN
1	B	95	ASN
1	B	265	ASN
2	C	188	GLN
2	C	193	GLN
2	D	188	GLN
2	D	239	ASN

5.3.3 RNA ⓘ

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	64V	B	301	-	36,41,41	0.80	1 (2%)	47,58,58	2.23	16 (34%)
3	64V	A	301	-	36,41,41	0.72	1 (2%)	47,58,58	2.13	17 (36%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	64V	B	301	-	-	7/18/32/32	0/5/5/5
3	64V	A	301	-	-	6/18/32/32	0/5/5/5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	301	64V	C31-N34	-2.34	1.30	1.34
3	A	301	64V	C31-N34	-2.09	1.31	1.34

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	301	64V	C17-N18-C19	6.14	124.03	117.81
3	A	301	64V	C29-N28-N34	-5.86	106.75	112.72
3	B	301	64V	C29-N28-N34	-5.76	106.84	112.72
3	B	301	64V	C17-N18-C19	5.28	123.16	117.81
3	B	301	64V	C3-N2-C7	4.65	116.02	109.52
3	B	301	64V	C13-N37-C36	4.02	121.96	117.49
3	A	301	64V	C10-C9-N8	-3.86	108.46	114.82
3	B	301	64V	C24-C22-C21	-3.67	120.00	123.50
3	A	301	64V	C13-N37-C36	3.44	121.31	117.49
3	A	301	64V	C9-N8-C5	3.20	121.92	118.20
3	B	301	64V	N37-C13-N8	3.07	119.98	116.57
3	B	301	64V	C10-C9-N8	-3.06	109.77	114.82
3	B	301	64V	C31-N34-N28	3.00	109.14	105.66
3	B	301	64V	C29-C30-C31	2.96	107.88	105.54
3	B	301	64V	C35-C36-N37	2.92	123.16	118.72
3	A	301	64V	C31-N34-N28	2.80	108.91	105.66
3	A	301	64V	O33-C32-C31	-2.64	106.28	112.10
3	A	301	64V	C24-C22-C21	-2.57	121.05	123.50
3	A	301	64V	C35-C36-N37	2.57	122.62	118.72
3	A	301	64V	C3-N2-C7	2.48	112.99	109.52
3	B	301	64V	C27-C21-C22	2.42	119.61	117.17
3	A	301	64V	C26-C25-C24	-2.41	118.89	121.74
3	A	301	64V	C29-C30-C31	2.37	107.41	105.54
3	B	301	64V	C30-C31-N34	-2.34	107.98	110.44
3	B	301	64V	C26-C25-C24	-2.28	119.04	121.74
3	B	301	64V	C4-C5-N8	-2.25	109.78	112.80
3	B	301	64V	C9-N8-C5	2.17	120.72	118.20
3	A	301	64V	C6-C5-C4	-2.15	106.24	111.19
3	B	301	64V	C16-C17-N18	-2.15	120.04	124.08
3	A	301	64V	C4-C5-N8	-2.12	109.96	112.80
3	A	301	64V	C35-C19-N18	-2.08	119.83	122.75
3	A	301	64V	C16-C17-N18	-2.06	120.20	124.08
3	A	301	64V	C16-C36-N37	-2.00	119.28	122.26

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	301	64V	C14-C13-N8-C5
3	A	301	64V	N37-C13-N8-C5
3	A	301	64V	C4-C5-N8-C9
3	A	301	64V	C6-C5-N8-C9
3	B	301	64V	C6-C5-N8-C9

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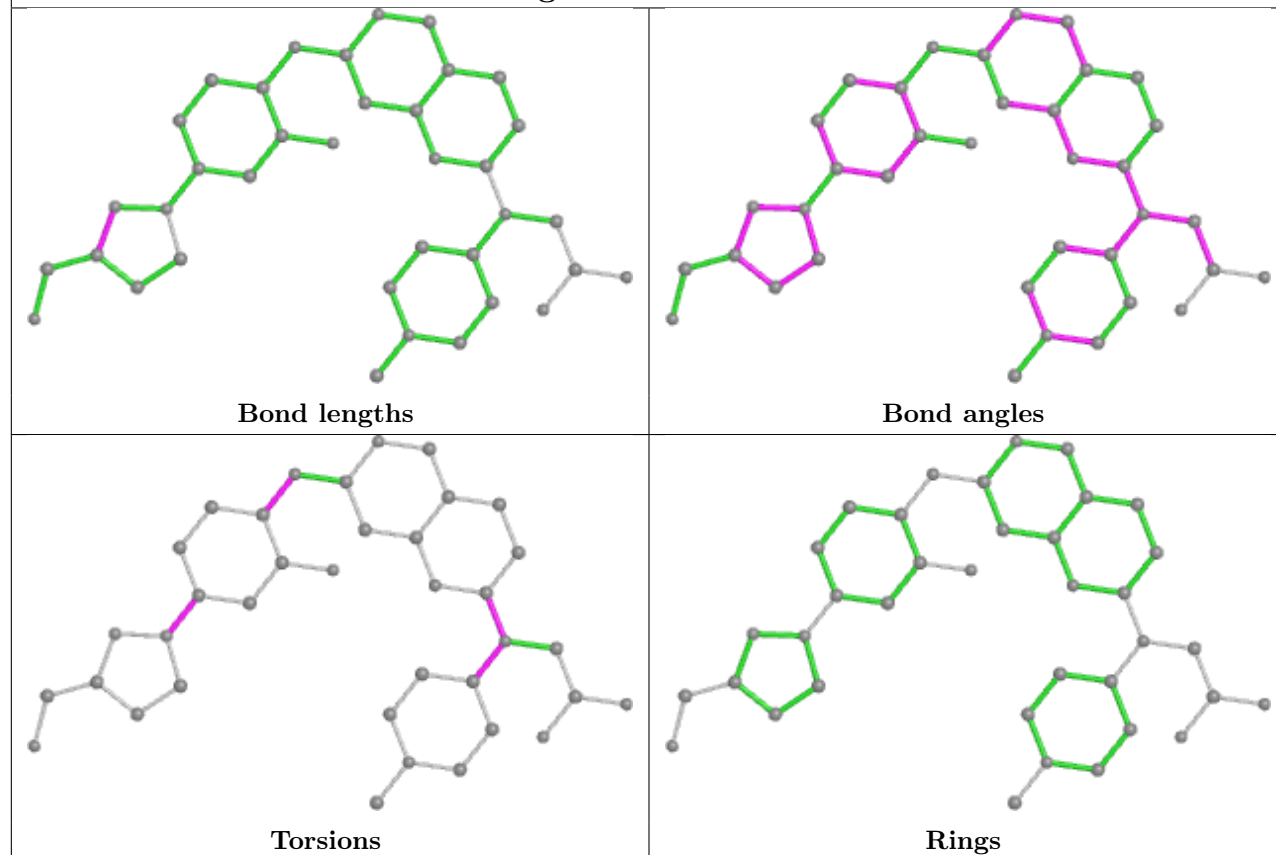
Mol	Chain	Res	Type	Atoms
3	B	301	64V	C22-C21-N20-C19
3	A	301	64V	C24-C25-N28-C29
3	B	301	64V	C27-C21-N20-C19
3	A	301	64V	C26-C25-N28-C29
3	B	301	64V	C26-C25-N28-C29
3	B	301	64V	C14-C13-N8-C5
3	B	301	64V	C4-C5-N8-C9
3	B	301	64V	C24-C25-N28-C29

There are no ring outliers.

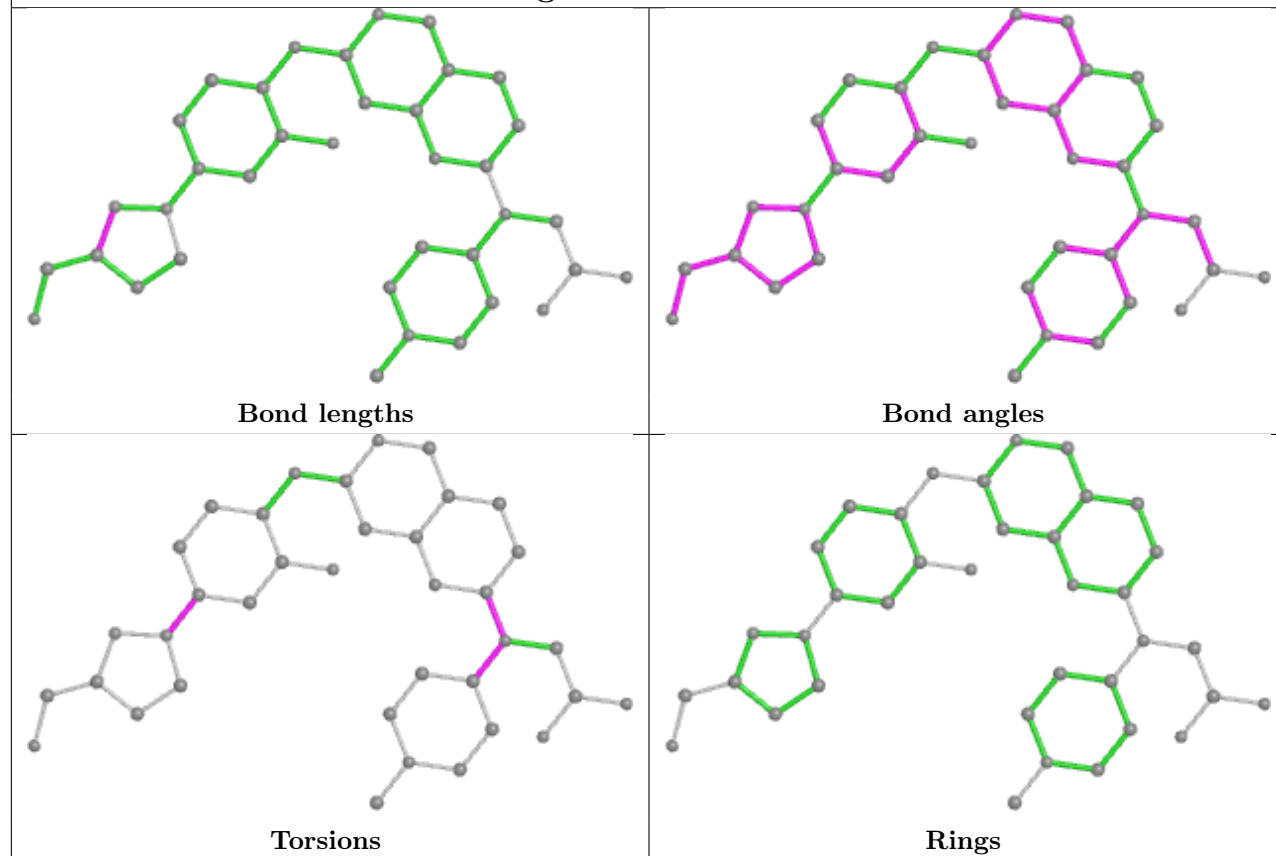
No monomer is involved in short contacts.

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.

Ligand 64V B 301



Ligand 64V A 301



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	290/292 (99%)	0.13	6 (2%) 63 62	50, 65, 102, 143	0
1	B	277/292 (94%)	1.43	75 (27%) 0 0	72, 105, 155, 180	0
2	C	148/209 (70%)	0.05	0 100 100	51, 67, 90, 99	0
2	D	136/209 (65%)	2.57	64 (47%) 0 0	106, 151, 186, 216	0
All	All	851/1002 (84%)	0.93	145 (17%) 1 1	50, 88, 165, 216	0

All (145) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	175	VAL	11.5
2	D	164	TYR	10.0
2	D	196	ILE	9.8
2	D	200	ASN	8.7
2	D	152	LEU	8.1
2	D	229	CYS	7.8
1	B	106	PHE	7.4
2	D	252	SER	7.3
2	D	220	HIS	7.0
1	B	22	LYS	6.8
2	D	151	LEU	6.7
1	B	245	THR	6.7
1	B	28	GLU	6.6
2	D	226	LEU	6.6
1	B	156	ARG	6.4
2	D	177	TRP	6.4
2	D	166	LEU	6.2
1	A	40	ASP	6.0
2	D	258	TRP	5.9
2	D	259	ASP	5.9
1	B	29	ILE	5.6

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Mol	Chain	Res	Type	RSRZ
1	B	287	SER	5.5
2	D	210	ASP	5.5
1	B	64	VAL	5.3
1	B	263	LEU	5.3
1	B	247	SER	5.3
1	B	110	LEU	5.2
2	D	261	CYS	5.2
2	D	163	CYS	5.1
1	B	70	LEU	5.0
2	D	202	VAL	4.9
1	B	31	ALA	4.9
2	D	203	PHE	4.9
1	A	41	ASP	4.9
1	B	30	VAL	4.7
2	D	222	LEU	4.7
1	B	23	ASN	4.7
1	B	286	PHE	4.6
2	D	173	ASP	4.5
1	B	24	ARG	4.5
1	B	140	LEU	4.5
1	B	103	VAL	4.5
2	D	174	PRO	4.4
1	B	20	LYS	4.4
1	B	78	LEU	4.4
2	D	248	PHE	4.3
2	D	216	VAL	4.2
1	B	32	LEU	4.2
1	B	111	LEU	4.2
1	B	21	ALA	4.1
2	D	185	LEU	4.0
1	B	27	HIS	4.0
1	B	55	LEU	4.0
1	B	246	THR	3.9
1	B	280	ALA	3.9
2	D	225	VAL	3.8
2	D	186	LEU	3.8
1	A	291	PRO	3.8
1	B	66	LEU	3.7
2	D	167	LYS	3.7
2	D	181	VAL	3.7
1	B	104	LYS	3.7
2	D	172	THR	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	65	ARG	3.6
2	D	155	LEU	3.5
2	D	179	ARG	3.5
2	D	249	LEU	3.5
1	B	284	PRO	3.4
2	D	254	LYS	3.4
2	D	178	LEU	3.4
1	B	71	HIS	3.4
1	B	108	PHE	3.4
2	D	228	THR	3.3
2	D	195	PHE	3.3
1	B	249	VAL	3.3
2	D	253	CYS	3.2
1	B	102	ILE	3.2
1	B	206	ASN	3.2
1	B	257	ALA	3.2
2	D	153	ARG	3.2
1	B	53	CYS	3.1
1	B	80	PHE	3.1
1	B	250	ASN	3.1
1	B	251	VAL	3.0
2	D	199	ALA	3.0
2	D	154	CYS	3.0
1	B	68	ASP	3.0
1	B	77	THR	2.8
1	A	224	GLU	2.8
1	B	221	THR	2.8
1	B	19	PHE	2.8
2	D	150	GLU	2.8
2	D	190	TRP	2.8
2	D	274	GLN	2.8
1	B	99	ASP	2.7
2	D	218	SER	2.7
2	D	244	PRO	2.7
2	D	207	LEU	2.7
2	D	219	ASP	2.7
1	B	42	GLU	2.7
1	B	265	ASN	2.6
1	B	56	LYS	2.6
1	B	133	LEU	2.5
2	D	291	ASN	2.5
1	B	216	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	101	GLU	2.5
1	B	281	LEU	2.4
1	B	190	CYS	2.4
2	D	189	GLY	2.4
1	B	217	ARG	2.4
2	D	232	LEU	2.3
2	D	283	THR	2.3
1	B	139	GLU	2.3
2	D	243	TYR	2.3
2	D	169	LEU	2.3
1	B	18	VAL	2.3
1	B	238	PRO	2.3
1	B	193	ALA	2.3
1	B	267	LEU	2.3
2	D	275	ILE	2.3
1	B	164	THR	2.3
1	B	244	ALA	2.3
2	D	176	LEU	2.3
1	B	248	LEU	2.2
1	B	135	ASN	2.2
2	D	271	LYS	2.2
1	B	4	TYR	2.2
2	D	148	THR	2.2
1	B	58	LEU	2.2
2	D	279	PRO	2.2
1	B	142	LEU	2.2
1	B	255	LEU	2.1
1	B	152	GLY	2.1
1	B	212	LEU	2.1
2	D	187	LEU	2.1
1	A	152	GLY	2.1
2	D	250	VAL	2.1
1	B	213	LYS	2.1
1	B	134	ILE	2.1
1	A	156	ARG	2.1
2	D	198	PRO	2.1
2	D	193	GLN	2.1
2	D	168	HIS	2.0
1	B	26	THR	2.0
1	B	222	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 monosaccharides in this entry.

6.4 Ligands [i](#)

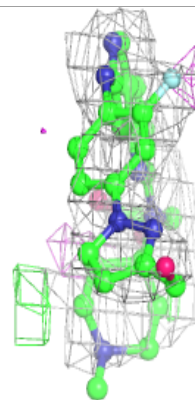
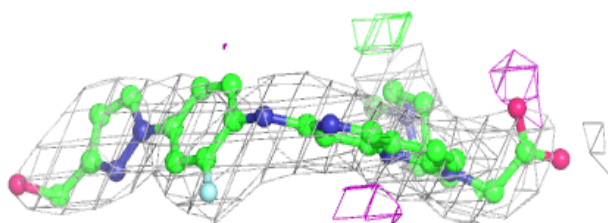
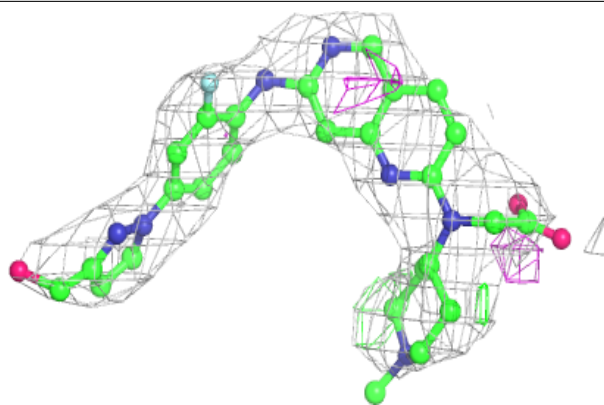
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	64V	B	301	37/37	0.81	0.29	86,96,106,112	0
3	64V	A	301	37/37	0.93	0.21	55,59,76,84	0

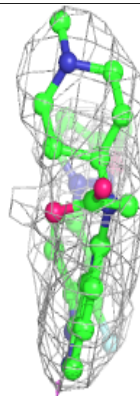
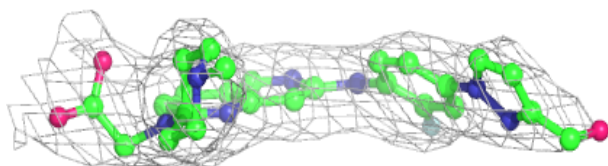
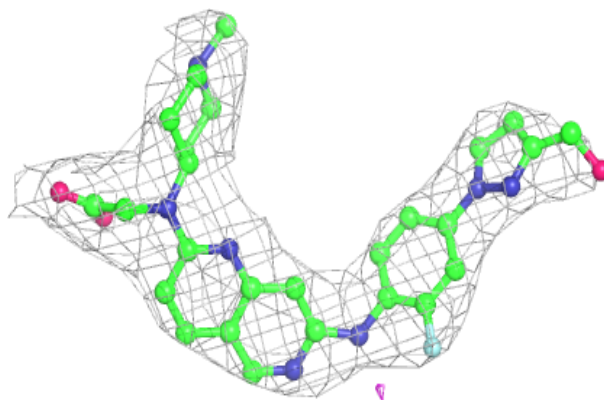
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 64V B 301:

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

**Electron density around 64V A 301:**

$2mF_o-DF_c$ (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.