



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:44 AM UTC

PDB ID : 6YVS / pdb_00006yvs
Title : FOCAL ADHESION KINASE CATALYTIC DOMAIN IN COMPLEX WITH
5-{4-[(Pyridin-3-ylmethyl)-amino]-5-trifluoromethyl-pyrimidin-2-ylamino}-1,3
-dihydro-indol-2-one
Authors : Musil, D.; Heinrich, T.; Amaral, M.
Deposited on : 2020-04-28
Resolution : 1.81 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

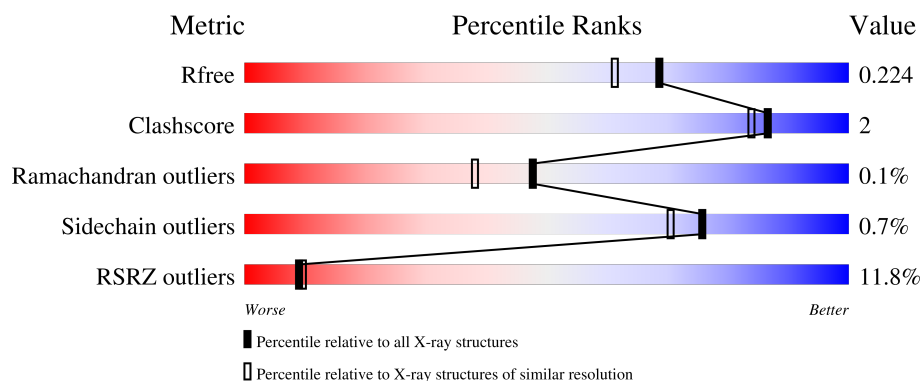
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.81 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	1112 (1.82-1.82)
Clashscore	190562	1148 (1.82-1.82)
Ramachandran outliers	187476	1140 (1.82-1.82)
Sidechain outliers	187428	1140 (1.82-1.82)
RSRZ outliers	180081	1112 (1.82-1.82)

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	282	10% 88% 5% 8%
1	B	282	10% 85% 5% 10%
1	C	282	14% 85% 5% 10%
1	D	282	9% 87% 5% 9%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9136 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.

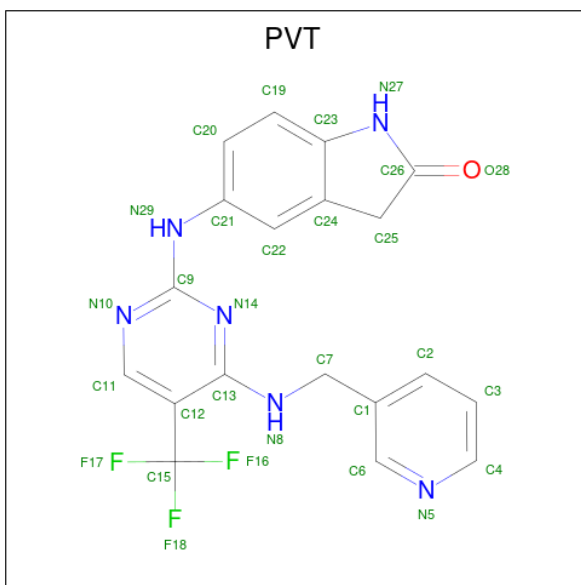
- Molecule 1 is a protein called Focal adhesion kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	260	Total	C	N	O	S	0	4	0
			2114	1346	367	384	17			
1	B	254	Total	C	N	O	S	0	4	0
			2068	1321	357	373	17			
1	C	254	Total	C	N	O	S	0	2	0
			2055	1309	355	374	17			
1	D	257	Total	C	N	O	S	0	0	0
			2069	1320	359	373	17			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	408	GLY	-	expression tag	UNP Q05397
A	409	SER	-	expression tag	UNP Q05397
A	410	GLY	-	expression tag	UNP Q05397
B	408	GLY	-	expression tag	UNP Q05397
B	409	SER	-	expression tag	UNP Q05397
B	410	GLY	-	expression tag	UNP Q05397
C	408	GLY	-	expression tag	UNP Q05397
C	409	SER	-	expression tag	UNP Q05397
C	410	GLY	-	expression tag	UNP Q05397
D	408	GLY	-	expression tag	UNP Q05397
D	409	SER	-	expression tag	UNP Q05397
D	410	GLY	-	expression tag	UNP Q05397

- Molecule 2 is 5-[[4-(pyridin-3-ylmethylamino)-5-(trifluoromethyl)pyrimidin-2-yl]amino]-1,3-dihydroindol-2-one (CCD ID: PVT) (formula: C₁₉H₁₅F₃N₆O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			29	19	3	6	1		
2	B	1	Total	C	F	N	O	0	1
			58	38	6	12	2		
2	C	1	Total	C	F	N	O	0	1
			58	38	6	12	2		
2	D	1	Total	C	F	N	O	0	1
			58	38	6	12	2		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

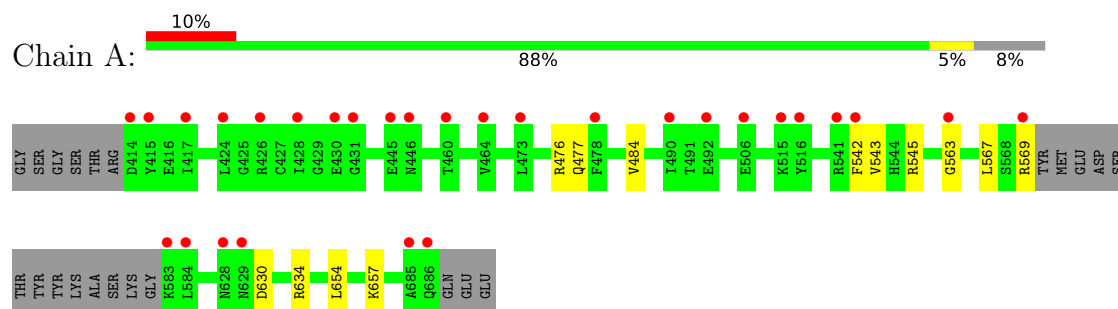
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	192	Total	O	0	0
			192	192		
4	B	182	Total	O	0	0
			182	182		
4	C	74	Total	O	0	0
			74	74		
4	D	174	Total	O	0	0
			174	174		

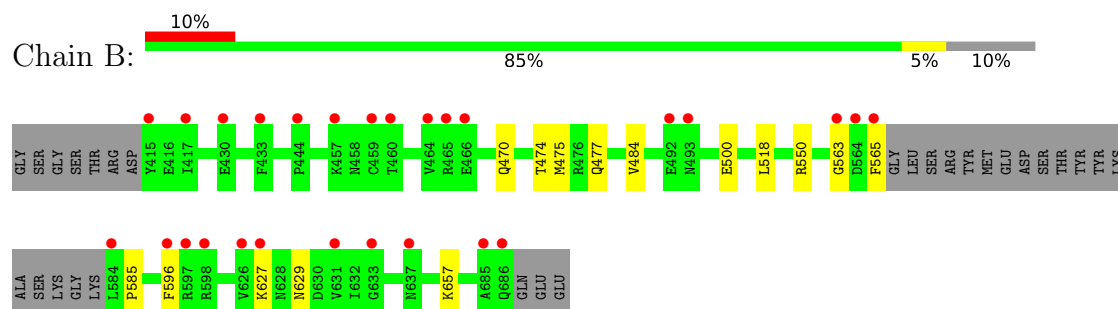
3 Residue-property plots [i](#)

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.

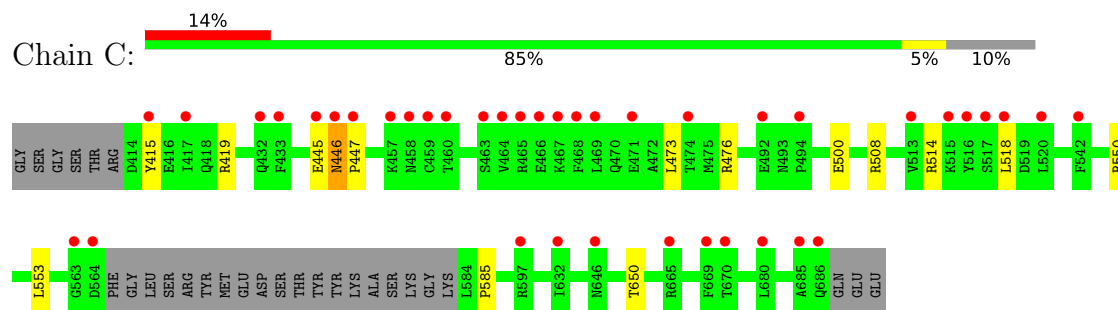
• Molecule 1: Focal adhesion kinase 1



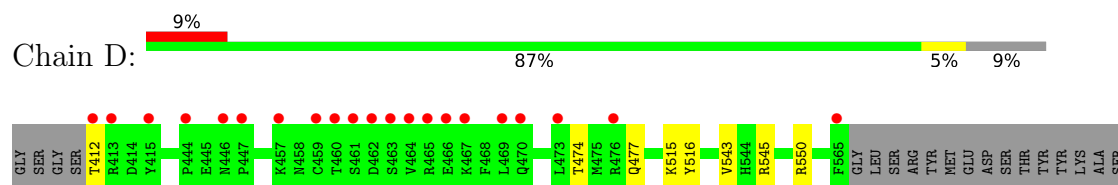
• Molecule 1: Focal adhesion kinase 1



• Molecule 1: Focal adhesion kinase 1



• Molecule 1: Focal adhesion kinase 1





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	114.17Å 76.08Å 172.85Å 90.00° 102.53° 90.00°	Depositor
Resolution (Å)	84.37 – 1.81 84.37 – 1.81	Depositor EDS
% Data completeness (in resolution range)	92.0 (84.37-1.81) 61.5 (84.37-1.81)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 1.81Å)	Xtriage
Refinement program	BUSTER 2.11.7 (19-MAR-2020)	Depositor
R, R_{free}	0.192 , 0.215 0.201 , 0.224	Depositor DCC
R_{free} test set	4095 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	30.0	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 42.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9136	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% 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: PVT, SO4

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	0.76	0/2160	1.04	0/2919
1	B	0.72	0/2117	0.98	0/2863
1	C	0.65	0/2100	1.04	2/2842 (0.1%)
1	D	0.77	0/2115	1.03	0/2860
All	All	0.73	0/8492	1.02	2/11484 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	476	ARG	CA-C-N	6.50	129.64	120.28
1	C	476	ARG	C-N-CA	6.50	129.64	120.28

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	2114	0	2119	7	0
1	B	2068	0	2080	10	0
1	C	2055	0	2057	10	0
1	D	2069	0	2074	8	0
2	A	29	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	58	0	0	1	0
2	C	58	0	0	2	0
2	D	58	0	0	1	0
3	B	5	0	0	0	0
4	A	192	0	0	0	0
4	B	182	0	0	0	0
4	C	74	0	0	0	0
4	D	174	0	0	1	0
All	All	9136	0	8330	36	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 2.

All (36) 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:C:446:ASN:HB3	1:C:447:PRO:HD2	1.67	0.76
1:B:484:VAL:HG12	1:B:563:GLY:HA2	1.75	0.68
1:A:654:LEU:HA	1:A:657:LYS:HE2	1.75	0.67
1:D:474:THR:O	1:D:477:GLN:HG2	1.96	0.66
1:B:484:VAL:CG1	1:B:563:GLY:HA2	2.28	0.63
1:B:474:THR:O	1:B:477:GLN:HG2	1.99	0.62
1:A:630:ASP:O	1:A:634:ARG:HG3	2.01	0.61
1:C:500:GLU:O	1:C:500:GLU:HG3	2.05	0.55
1:B:627[A]:LYS:HZ3	1:B:629:ASN:HB2	1.73	0.52
2:B:701[B]:PVT:N14	2:B:701[B]:PVT:C20	2.73	0.51
1:C:445:GLU:O	1:C:446:ASN:HB2	2.11	0.51
1:A:484:VAL:HG22	1:A:563:GLY:HA2	1.93	0.51
1:A:545:ARG:HD3	1:A:567:LEU:O	2.11	0.50
1:B:627[A]:LYS:NZ	1:B:629:ASN:HB2	2.26	0.50
1:C:514:ARG:HB3	1:C:518:LEU:HG	1.94	0.49
1:D:515:LYS:HB3	4:D:919:HOH:O	2.14	0.47
1:D:550:ARG:HH22	1:D:585:PRO:HB3	1.79	0.47
2:A:701:PVT:N14	2:A:701:PVT:C20	2.78	0.47
1:C:550:ARG:HH22	1:C:585:PRO:HB3	1.80	0.46
1:A:542:PHE:CD1	1:A:569:ARG:HB2	2.51	0.45
1:D:550:ARG:NH2	1:D:585:PRO:HB3	2.32	0.45
1:C:553:LEU:HD13	2:C:701[A]:PVT:C11	2.48	0.44
1:C:419:ARG:HD2	1:D:516:TYR:HB2	1.99	0.44
1:D:671:GLU:HG3	1:D:675:GLN:HE21	1.83	0.43
1:B:475:MET:SD	1:B:565:PHE:CD1	3.12	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:588:TRP:HH2	1:D:621:LYS:HE3	1.83	0.42
1:B:484:VAL:HG23	1:B:500:GLU:HG2	2.01	0.42
1:C:415:TYR:OH	1:C:473:LEU:HD23	2.19	0.42
2:D:701[A]:PVT:N14	2:D:701[A]:PVT:C20	2.82	0.42
1:B:550:ARG:HH22	1:B:585:PRO:HB3	1.85	0.41
1:A:543:VAL:HG12	1:A:545:ARG:HG3	2.03	0.41
1:B:657:LYS:HE2	1:B:657:LYS:HB3	1.91	0.41
1:C:550:ARG:NH2	1:C:585:PRO:HB3	2.35	0.41
1:C:553:LEU:HD13	2:C:701[B]:PVT:C11	2.51	0.41
1:D:543:VAL:HG12	1:D:545:ARG:HG3	2.03	0.41
1:A:476:ARG:HD3	1:B:596:PHE:CE1	2.57	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	260/282 (92%)	256 (98%)	4 (2%)	0	100	100
1	B	254/282 (90%)	252 (99%)	2 (1%)	0	100	100
1	C	252/282 (89%)	248 (98%)	3 (1%)	1 (0%)	30	19
1	D	253/282 (90%)	251 (99%)	2 (1%)	0	100	100
All	All	1019/1128 (90%)	1007 (99%)	11 (1%)	1 (0%)	48	38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	446	ASN

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	235/249 (94%)	234 (100%)	1 (0%)	84	81
1	B	230/249 (92%)	228 (99%)	2 (1%)	70	62
1	C	228/249 (92%)	226 (99%)	2 (1%)	70	62
1	D	229/249 (92%)	228 (100%)	1 (0%)	84	81
All	All	922/996 (93%)	916 (99%)	6 (1%)	76	70

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	477	GLN
1	B	470	GLN
1	B	518	LEU
1	C	508	ARG
1	C	650	THR
1	D	412	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	418	GLN
1	A	446	ASN
1	B	418	GLN
1	B	438	GLN
1	B	629	ASN
1	C	470	GLN
1	D	432	GLN
1	D	493	ASN
1	D	629	ASN
1	D	675	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

8 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	PVT	C	701[B]	-	32,32,32	0.21	0	44,46,46	0.63	1 (2%)
2	PVT	B	701[A]	-	32,32,32	0.21	0	44,46,46	0.66	1 (2%)
2	PVT	D	701[A]	-	32,32,32	0.21	0	44,46,46	0.61	1 (2%)
2	PVT	C	701[A]	-	32,32,32	0.21	0	44,46,46	0.67	1 (2%)
2	PVT	A	701	-	32,32,32	0.25	0	44,46,46	0.63	1 (2%)
2	PVT	D	701[B]	-	32,32,32	0.22	0	44,46,46	0.65	1 (2%)
2	PVT	B	701[B]	-	32,32,32	0.21	0	44,46,46	0.63	1 (2%)
3	SO4	B	702	-	4,4,4	0.27	0	6,6,6	0.29	0

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 Torsions and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PVT	C	701[B]	-	-	0/15/23/23	0/4/4/4

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PVT	B	701[A]	-	-	0/15/23/23	0/4/4/4
2	PVT	D	701[A]	-	-	0/15/23/23	0/4/4/4
2	PVT	C	701[A]	-	-	0/15/23/23	0/4/4/4
2	PVT	A	701	-	-	0/15/23/23	0/4/4/4
2	PVT	D	701[B]	-	-	0/15/23/23	0/4/4/4
2	PVT	B	701[B]	-	-	0/15/23/23	0/4/4/4

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	701[A]	PVT	C15-C12-C13	3.21	124.16	120.06
2	C	701[A]	PVT	C15-C12-C13	3.13	124.06	120.06
2	A	701	PVT	C15-C12-C13	3.11	124.03	120.06
2	C	701[B]	PVT	C15-C12-C13	3.10	124.02	120.06
2	B	701[B]	PVT	C15-C12-C13	3.05	123.96	120.06
2	D	701[B]	PVT	C15-C12-C13	2.99	123.88	120.06
2	D	701[A]	PVT	C15-C12-C13	2.97	123.86	120.06

There are no chirality outliers.

There are no torsion outliers.

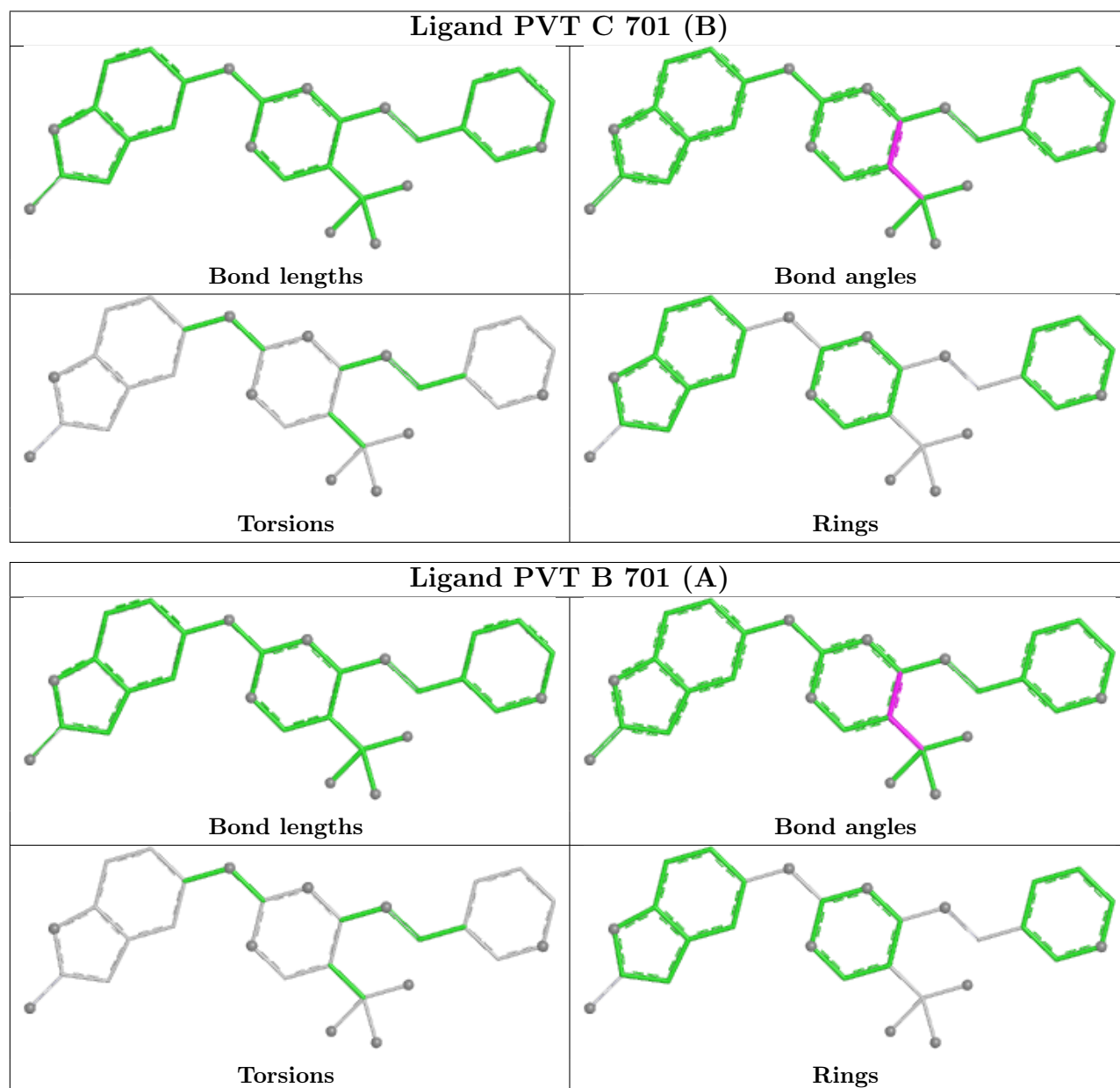
There are no ring outliers.

5 monomers are involved in 5 short contacts:

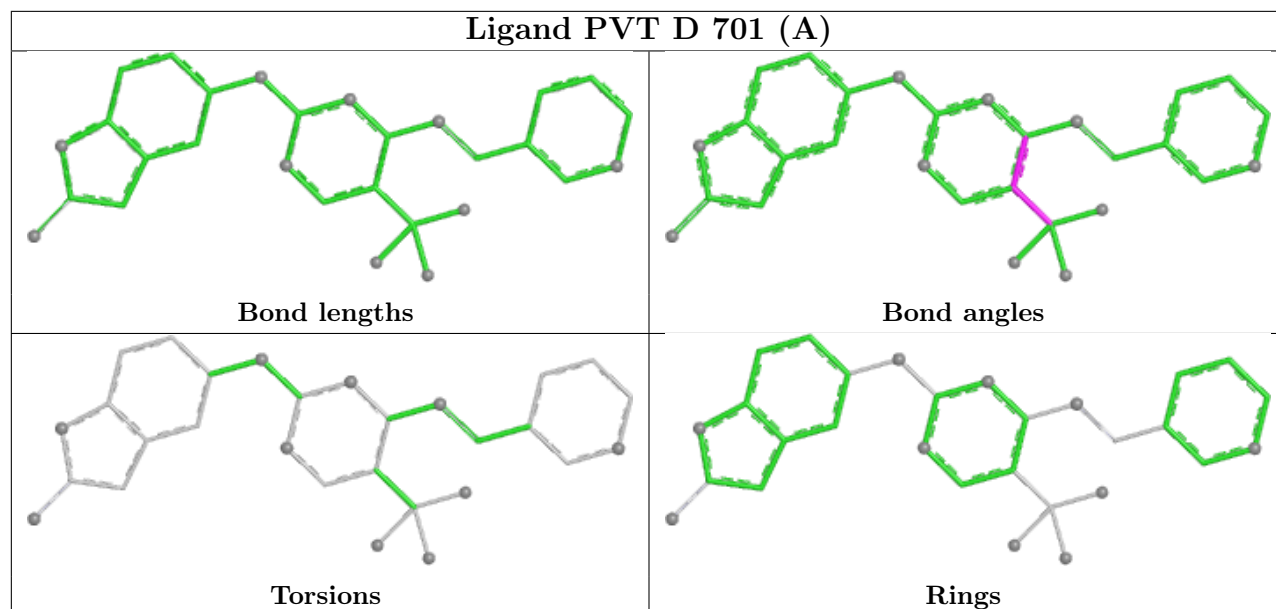
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	701[B]	PVT	1	0
2	D	701[A]	PVT	1	0
2	C	701[A]	PVT	1	0
2	A	701	PVT	1	0
2	B	701[B]	PVT	1	0

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.

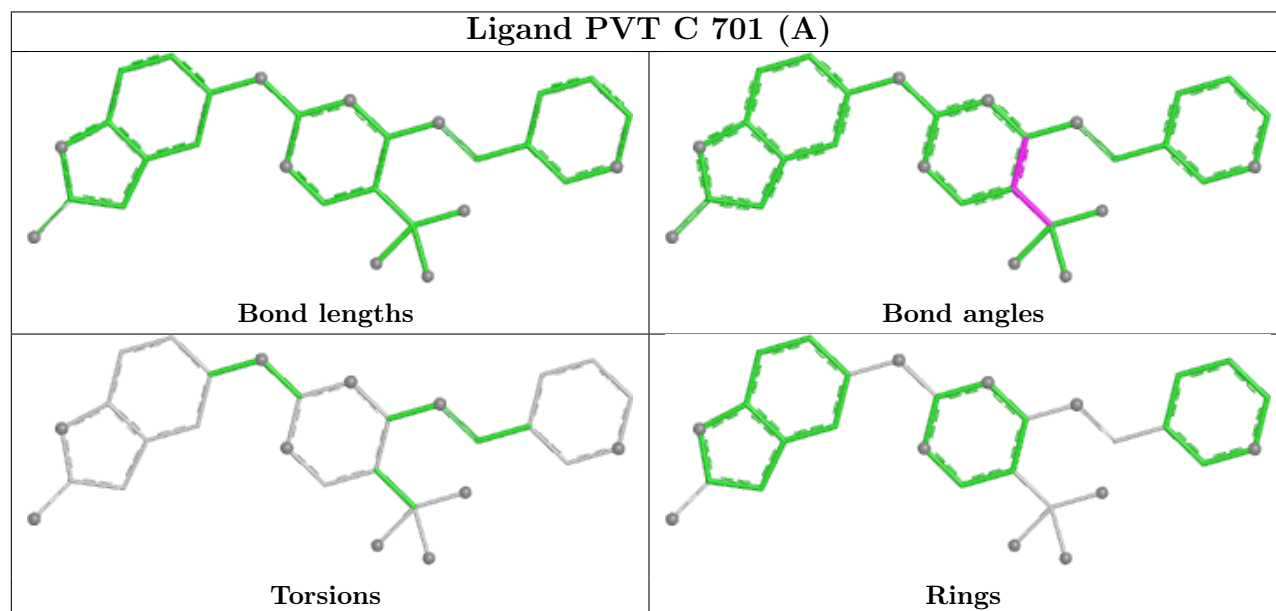
The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

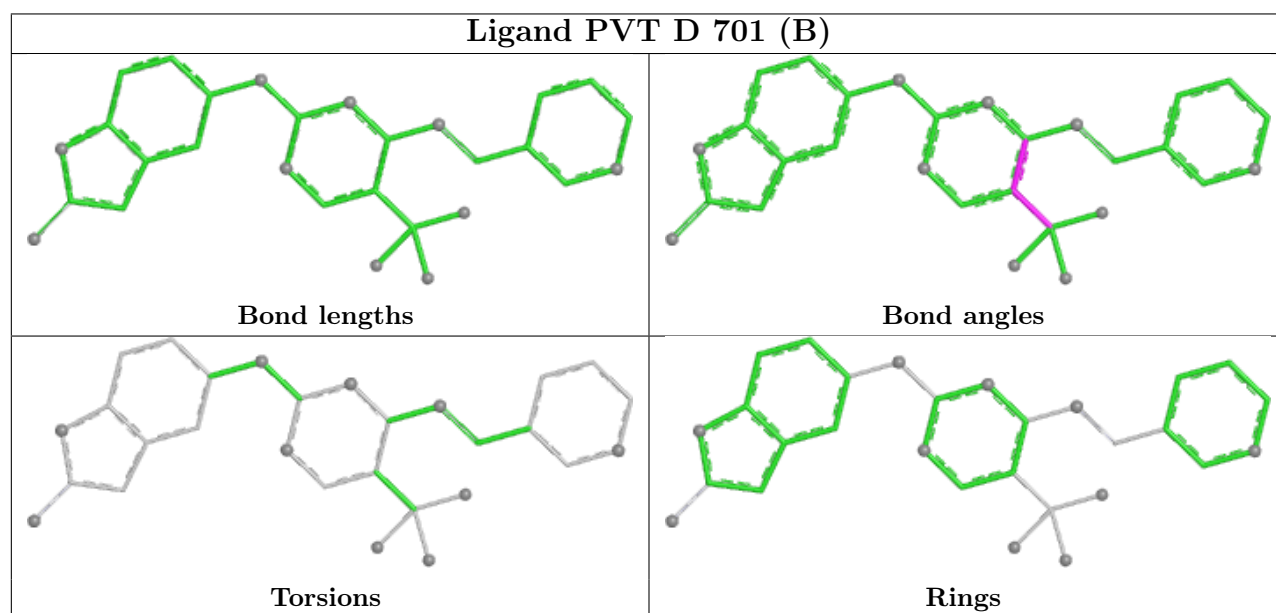
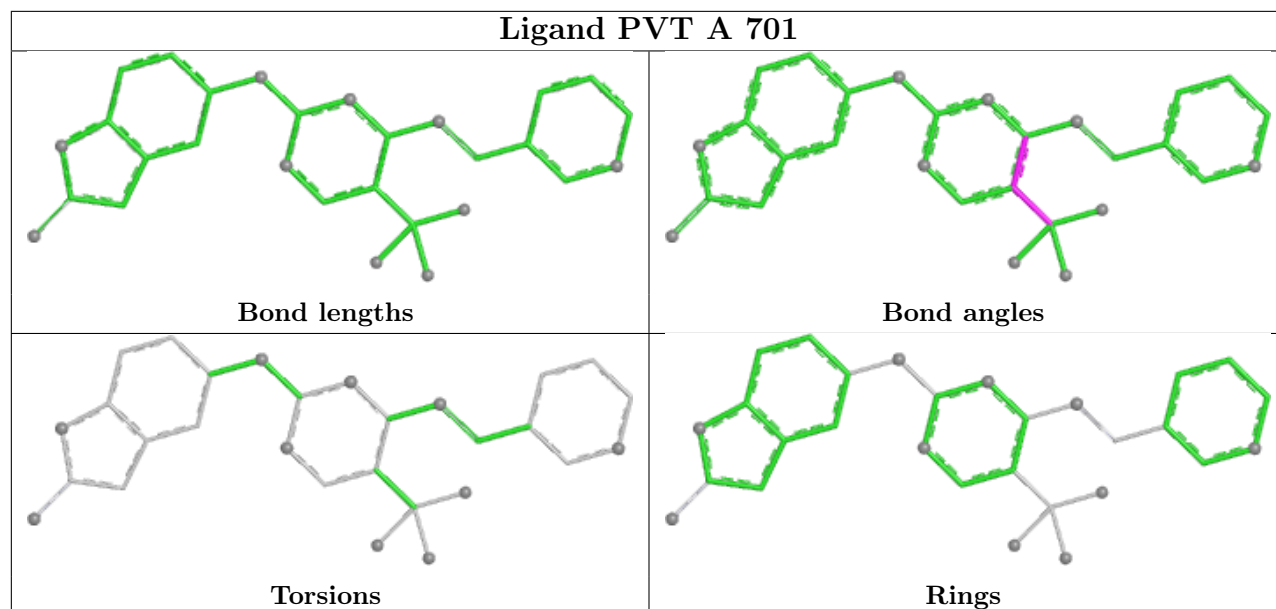


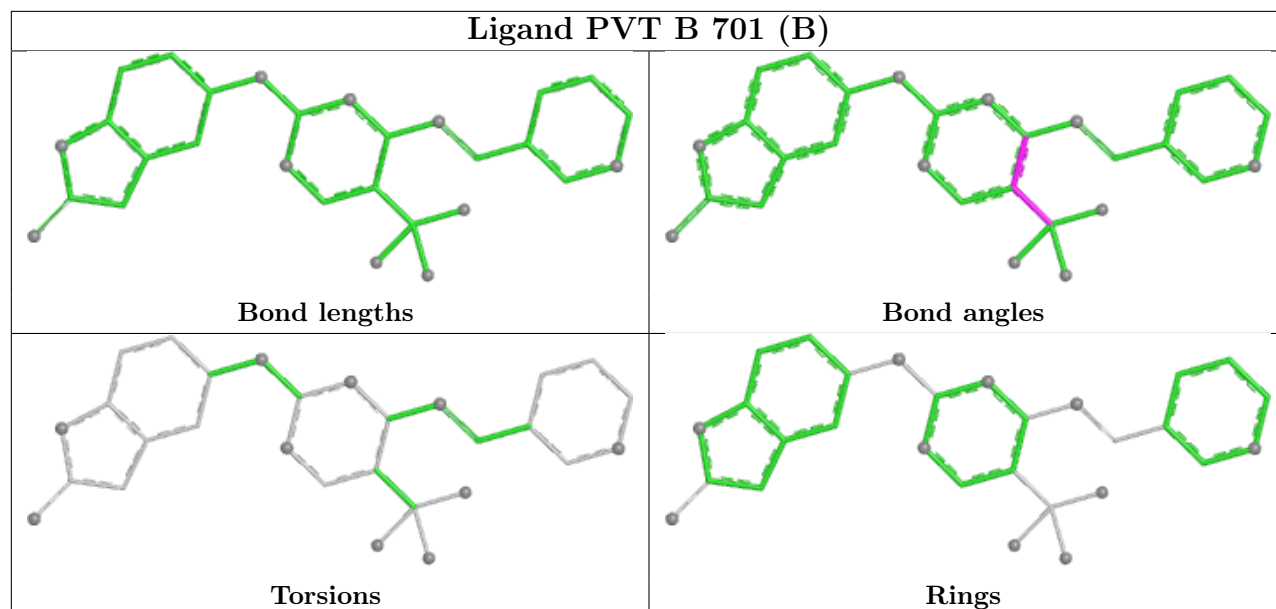
Ligand PVT D 701 (A)



Ligand PVT C 701 (A)







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	260/282 (92%)	0.60	29 (11%)	10	11	13, 33, 56, 72	4 (1%)
1	B	254/282 (90%)	0.44	27 (10%)	11	12	12, 32, 62, 77	4 (1%)
1	C	254/282 (90%)	1.11	40 (15%)	5	5	20, 45, 81, 93	2 (0%)
1	D	257/282 (91%)	0.34	25 (9%)	13	15	21, 32, 60, 71	0
All	All	1025/1128 (90%)	0.62	121 (11%)	9	10	12, 37, 64, 93	10 (0%)

All (121) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	516	TYR	6.3
1	C	460	THR	5.0
1	B	627[A]	LYS	4.4
1	B	685	ALA	4.3
1	B	444	PRO	4.1
1	B	415	TYR	4.1
1	B	460	THR	4.0
1	C	494	PRO	3.9
1	C	415	TYR	3.9
1	C	518	LEU	3.9
1	D	415	TYR	3.8
1	C	685	ALA	3.8
1	B	563	GLY	3.8
1	A	629	ASN	3.8
1	C	542	PHE	3.7
1	D	460	THR	3.7
1	D	412	THR	3.7
1	B	564	ASP	3.6
1	B	596	PHE	3.5
1	C	597	ARG	3.5
1	D	470	GLN	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	564	ASP	3.3
1	C	464	VAL	3.3
1	D	686	GLN	3.3
1	C	665	ARG	3.3
1	C	459	CYS	3.3
1	D	459	CYS	3.3
1	C	632	ILE	3.3
1	A	685	ALA	3.3
1	B	631	VAL	3.2
1	C	686	GLN	3.1
1	B	466	GLU	3.1
1	C	467	LYS	3.0
1	B	417	ILE	3.0
1	C	458	ASN	3.0
1	C	515	LYS	3.0
1	B	686	GLN	2.9
1	A	569	ARG	2.9
1	D	463	SER	2.9
1	A	686	GLN	2.9
1	D	624	GLN	2.9
1	C	445	GLU	2.9
1	A	430	GLU	2.9
1	A	628	ASN	2.8
1	D	464	VAL	2.8
1	B	565	PHE	2.8
1	C	471	GLU	2.8
1	D	565	PHE	2.8
1	C	432	GLN	2.8
1	A	563	GLY	2.8
1	C	646	ASN	2.7
1	A	415	TYR	2.7
1	C	433	PHE	2.7
1	A	584	LEU	2.7
1	D	467	LYS	2.7
1	D	465	ARG	2.6
1	C	469	LEU	2.6
1	B	464	VAL	2.6
1	B	459	CYS	2.6
1	C	517	SER	2.6
1	A	478	PHE	2.6
1	C	516	TYR	2.5
1	D	444	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	430	GLU	2.5
1	D	584	LEU	2.5
1	B	465	ARG	2.5
1	C	457	LYS	2.5
1	D	447	PRO	2.4
1	B	493	ASN	2.4
1	A	542	PHE	2.4
1	B	433	PHE	2.4
1	D	457	LYS	2.4
1	A	426	ARG	2.3
1	C	520	LEU	2.3
1	A	445	GLU	2.3
1	C	446	ASN	2.3
1	B	626	VAL	2.3
1	D	476	ARG	2.3
1	A	473	LEU	2.3
1	C	417	ILE	2.3
1	D	413	ARG	2.3
1	C	466	GLU	2.3
1	B	584	LEU	2.3
1	C	468	PHE	2.3
1	A	414	ASP	2.3
1	A	431	GLY	2.2
1	B	597	ARG	2.2
1	C	465	ARG	2.2
1	A	424	LEU	2.2
1	D	473	LEU	2.2
1	A	428	ILE	2.2
1	D	466	GLU	2.2
1	C	474	THR	2.2
1	C	463	SER	2.2
1	B	598	ARG	2.2
1	A	460	THR	2.2
1	C	680	LEU	2.2
1	D	446	ASN	2.1
1	A	490	ILE	2.1
1	A	464	VAL	2.1
1	B	633	GLY	2.1
1	A	583	LYS	2.1
1	A	446	ASN	2.1
1	D	685	ALA	2.1
1	C	669	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	417	ILE	2.1
1	D	461	SER	2.1
1	A	541	ARG	2.1
1	C	563	GLY	2.1
1	A	506	GLU	2.1
1	B	492	GLU	2.1
1	C	492	GLU	2.1
1	B	637	ASN	2.1
1	A	492	GLU	2.0
1	B	457	LYS	2.0
1	C	670	THR	2.0
1	D	469	LEU	2.0
1	D	462	ASP	2.0
1	A	515	LYS	2.0
1	C	447	PRO	2.0
1	C	513	VAL	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
3	SO4	B	702	5/5	0.85	0.15	58,59,59,59	0
2	PVT	D	701[B]	29/29	0.89	0.10	27,28,35,35	29
2	PVT	D	701[A]	29/29	0.89	0.10	26,28,34,34	29
2	PVT	C	701[B]	29/29	0.90	0.10	29,30,31,31	29
2	PVT	C	701[A]	29/29	0.90	0.10	36,36,37,37	29
2	PVT	A	701	29/29	0.92	0.09	27,32,35,36	0
2	PVT	B	701[B]	29/29	0.93	0.09	21,25,28,29	29

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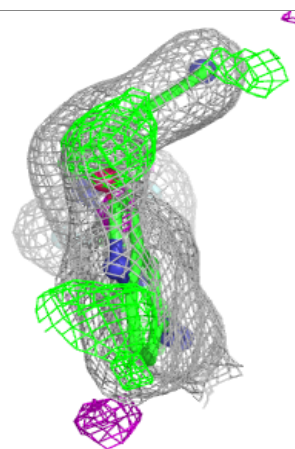
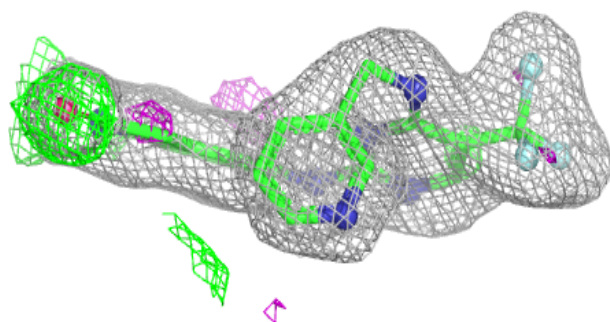
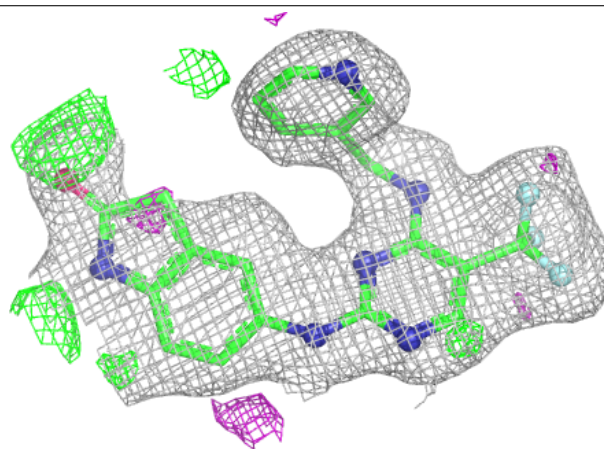
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PVT	B	701[A]	29/29	0.93	0.09	26,29,31,31	29

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.

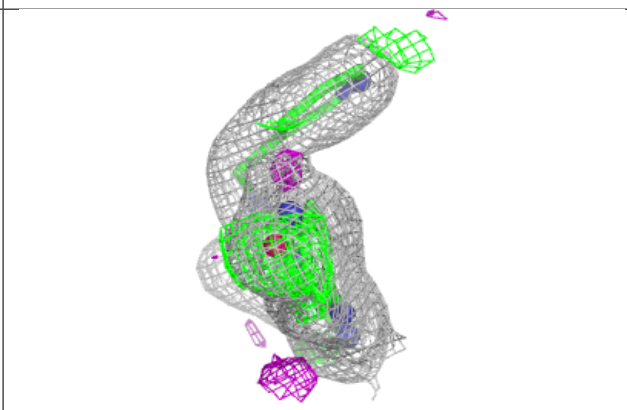
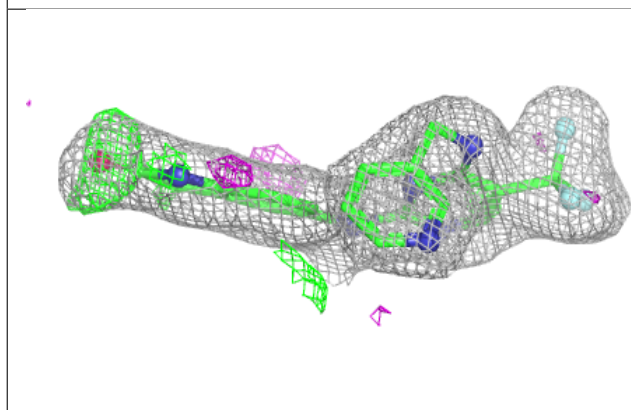
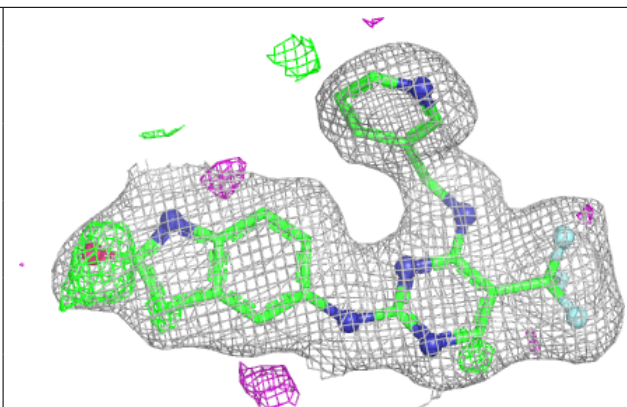
Electron density around PVT D 701 (B):

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

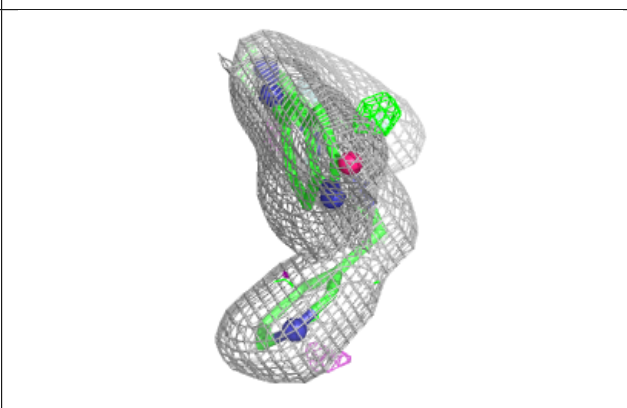
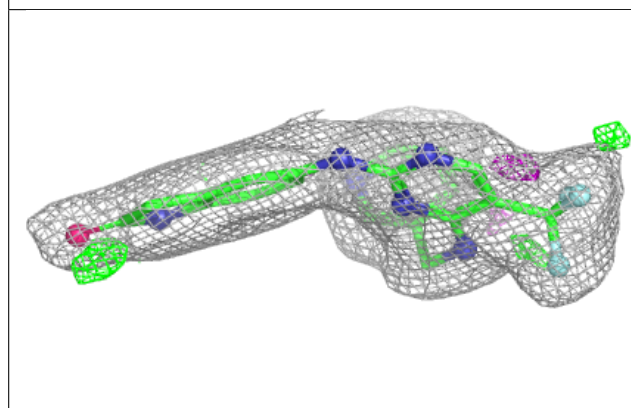
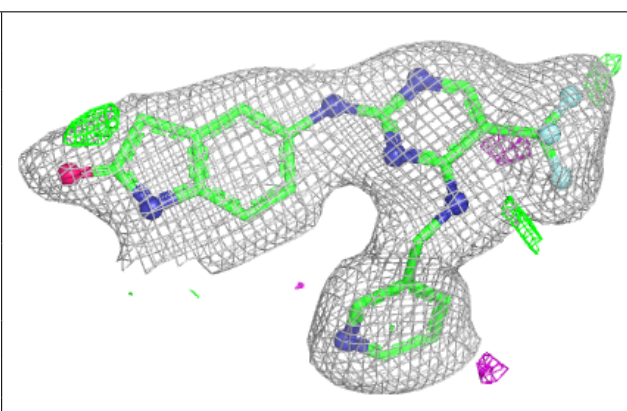


Electron density around PVT D 701 (A):

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

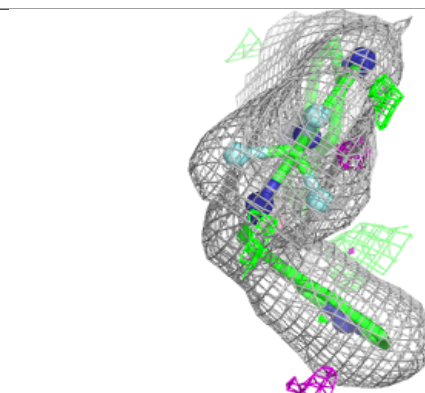
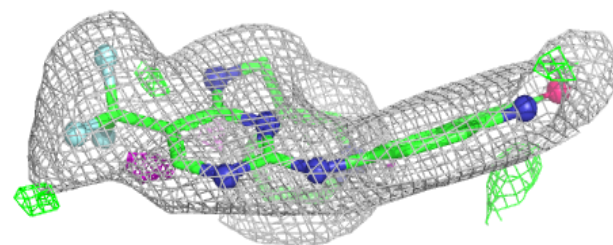
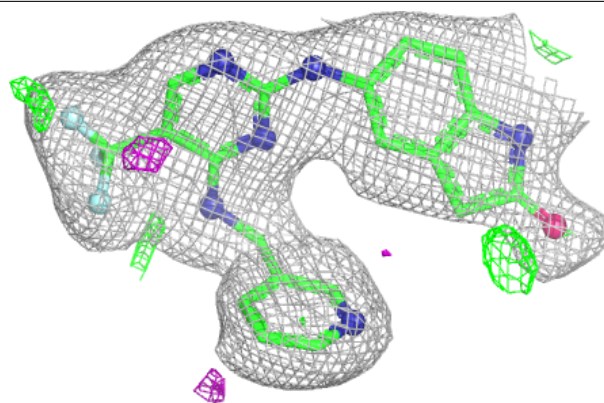
**Electron density around PVT C 701 (B):**

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

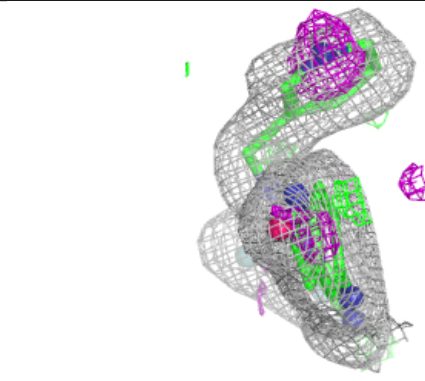
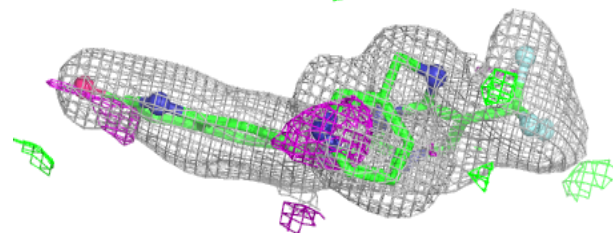
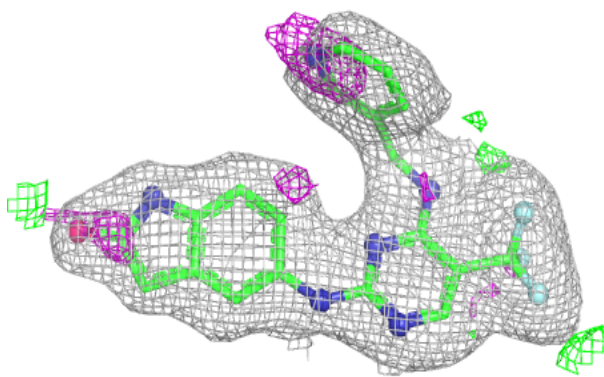


Electron density around PVT C 701 (A):

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

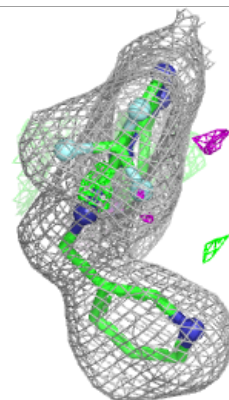
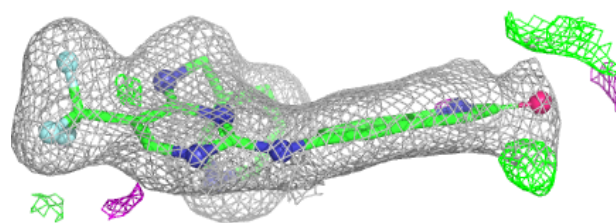
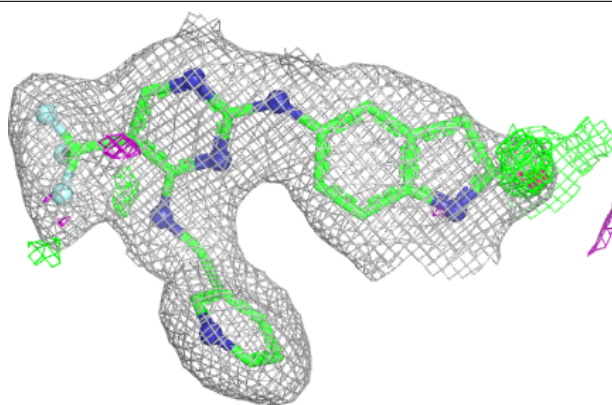
**Electron density around PVT A 701:**

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

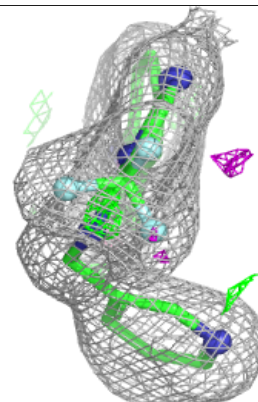
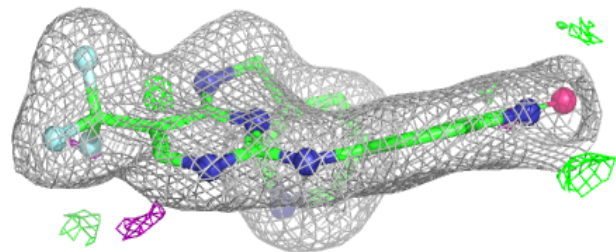
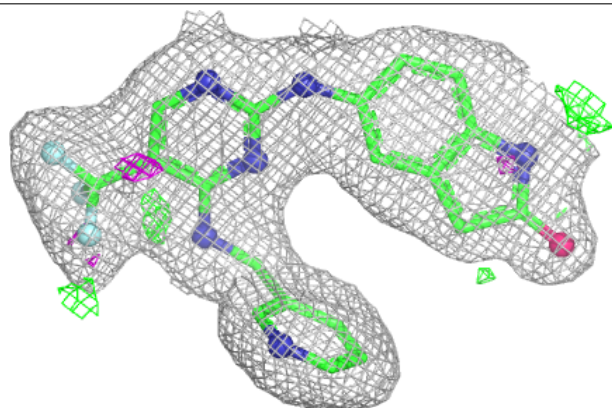


Electron density around PVT B 701 (B):

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

**Electron density around PVT B 701 (A):**

$2mF_o-DF_c$ (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.