



# Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 11:56 pm BST

PDB ID : 4IDV  
Title : Crystal Structure of NIK with compound 4-{3-[2-amino-5-(2-methoxyethoxy)pyrimidin-4-yl]-1H-indol-5-yl}-2-methylbut-3-yn-2-ol (13V)  
Authors : Liu, J.; Sudom, A.; Wang, Z.  
Deposited on : 2012-12-13  
Resolution : 2.90 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](mailto: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.

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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 : 2.11  
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) : 2.11

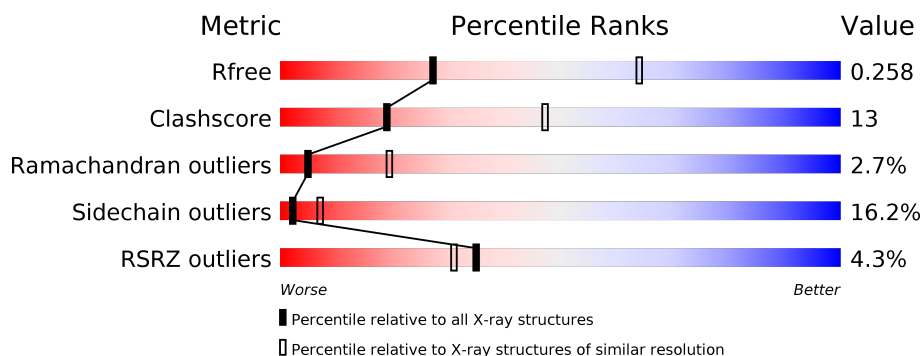
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	356	<div> <div>3%</div> <div> <div></div> <div>65%</div> <div>23%</div> <div>• • 6%</div> </div> </div>
1	B	356	<div> <div>4%</div> <div> <div></div> <div>60%</div> <div>25%</div> <div>7% • 6%</div> </div> </div>
1	C	356	<div> <div>6%</div> <div> <div></div> <div>64%</div> <div>23%</div> <div>7% • 6%</div> </div> </div>
1	D	356	<div> <div>4%</div> <div> <div></div> <div>62%</div> <div>24%</div> <div>6% • 6%</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10534 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.

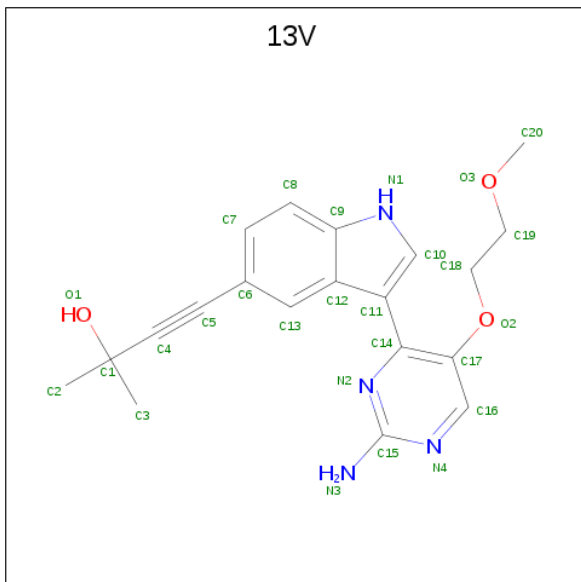
- Molecule 1 is a protein called Mitogen-activated protein kinase kinase kinase 14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	335	Total	C	N	O	S	0	1	0
			2595	1636	463	478	18			
1	B	335	Total	C	N	O	S	0	1	0
			2595	1636	463	478	18			
1	C	335	Total	C	N	O	S	0	1	0
			2595	1636	463	478	18			
1	D	335	Total	C	N	O	S	0	1	0
			2595	1636	463	478	18			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	325	GLY	-	EXPRESSION TAG	UNP Q99558
A	326	ALA	-	EXPRESSION TAG	UNP Q99558
A	327	MET	-	EXPRESSION TAG	UNP Q99558
A	328	GLY	-	EXPRESSION TAG	UNP Q99558
A	329	SER	-	EXPRESSION TAG	UNP Q99558
B	325	GLY	-	EXPRESSION TAG	UNP Q99558
B	326	ALA	-	EXPRESSION TAG	UNP Q99558
B	327	MET	-	EXPRESSION TAG	UNP Q99558
B	328	GLY	-	EXPRESSION TAG	UNP Q99558
B	329	SER	-	EXPRESSION TAG	UNP Q99558
C	325	GLY	-	EXPRESSION TAG	UNP Q99558
C	326	ALA	-	EXPRESSION TAG	UNP Q99558
C	327	MET	-	EXPRESSION TAG	UNP Q99558
C	328	GLY	-	EXPRESSION TAG	UNP Q99558
C	329	SER	-	EXPRESSION TAG	UNP Q99558
D	325	GLY	-	EXPRESSION TAG	UNP Q99558
D	326	ALA	-	EXPRESSION TAG	UNP Q99558
D	327	MET	-	EXPRESSION TAG	UNP Q99558
D	328	GLY	-	EXPRESSION TAG	UNP Q99558
D	329	SER	-	EXPRESSION TAG	UNP Q99558

- Molecule 2 is 4-{3-[2-amino-5-(2-methoxyethoxy)pyrimidin-4-yl]-1H-indol-5-yl}-2-methylbut-3-yn-2-ol (three-letter code: 13V) (formula: C<sub>20</sub>H<sub>22</sub>N<sub>4</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			27	20	4	3		
2	B	1	Total	C	N	O	0	0
			27	20	4	3		
2	C	1	Total	C	N	O	0	0
			27	20	4	3		
2	D	1	Total	C	N	O	0	0
			27	20	4	3		

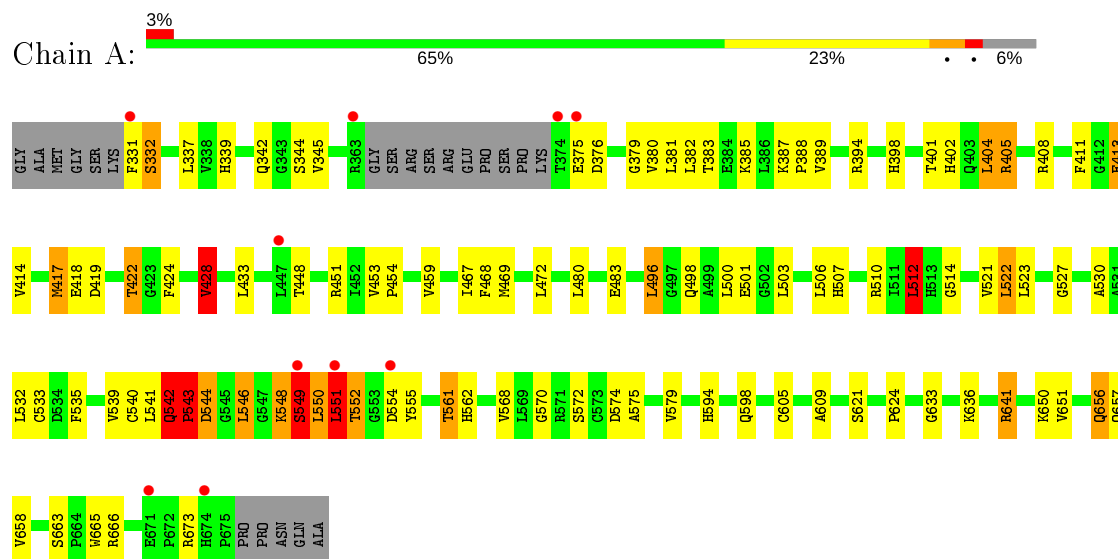
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	9	Total	O	0	0
			9	9		
3	B	12	Total	O	0	0
			12	12		
3	C	8	Total	O	0	0
			8	8		
3	D	17	Total	O	0	0
			17	17		

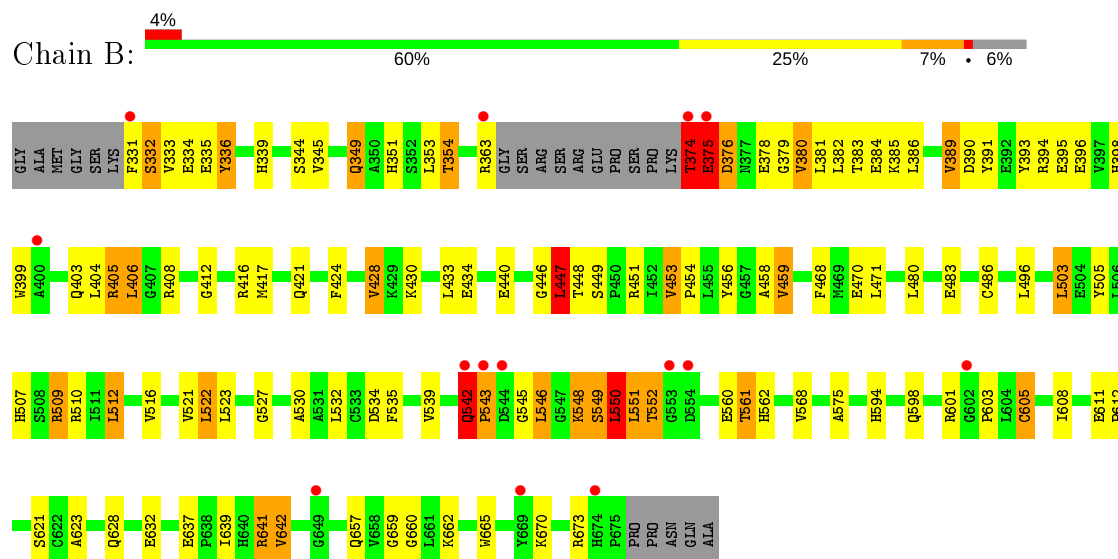
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

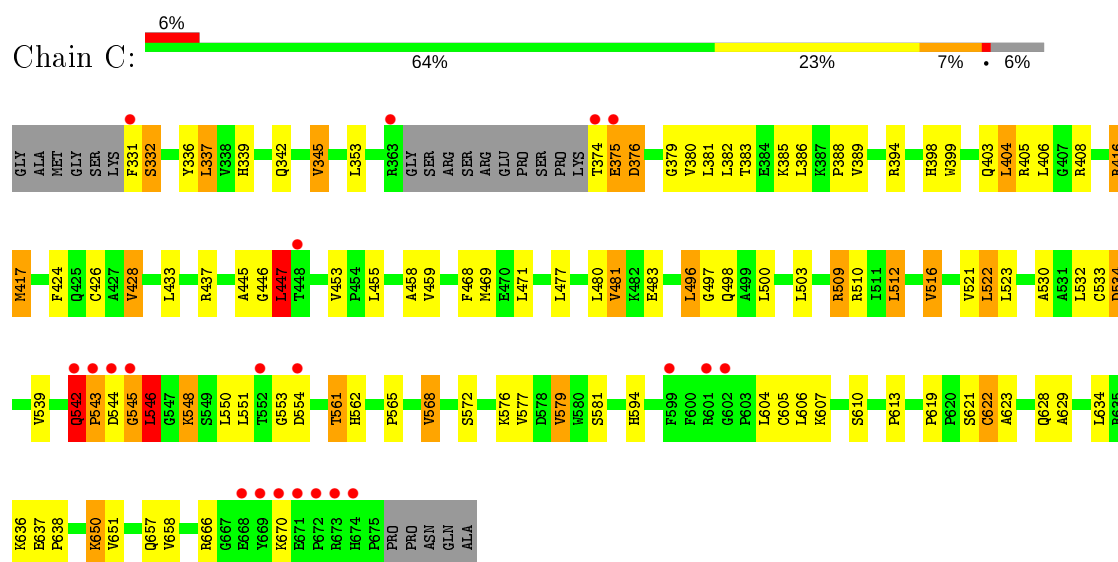
- Molecule 1: Mitogen-activated protein kinase kinase kinase 14



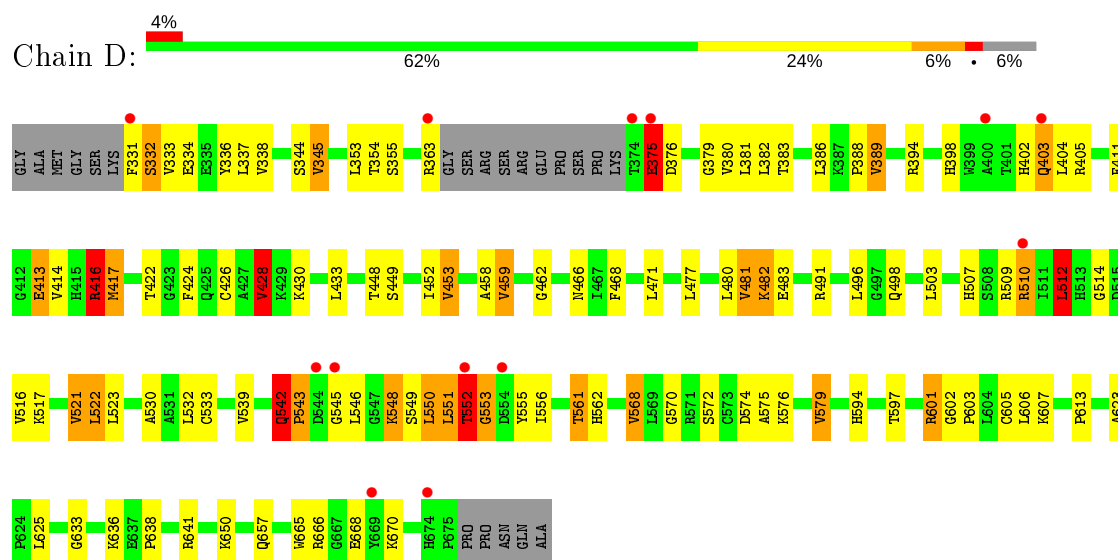
- Molecule 1: Mitogen-activated protein kinase kinase kinase 14



- Molecule 1: Mitogen-activated protein kinase kinase kinase 14



- Molecule 1: Mitogen-activated protein kinase kinase 14



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	45.38Å 146.01Å 254.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.14 – 2.90 32.12 – 2.90	Depositor EDS
% Data completeness (in resolution range)	80.7 (32.14-2.90) 80.8 (32.12-2.90)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	0.14	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.88 (at 2.90Å)	Xtriage
Refinement program	PHENIX, REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.203 , 0.269 0.203 , 0.258	Depositor DCC
$R_{free}$ test set	1572 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	43.8	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 39.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	10534	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.25% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: 13V

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	0.58	0/2656	0.83	2/3597 (0.1%)
1	B	0.55	0/2656	0.83	3/3597 (0.1%)
1	C	0.57	0/2656	0.83	4/3597 (0.1%)
1	D	0.58	0/2656	0.84	7/3597 (0.2%)
All	All	0.57	0/10624	0.83	16/14388 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	3
1	C	0	1
1	D	0	1
All	All	0	8

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	512	LEU	CA-CB-CG	8.04	133.79	115.30
1	C	542	GLN	C-N-CD	-6.47	106.37	120.60
1	C	546	LEU	CA-CB-CG	6.39	130.00	115.30
1	D	416	ARG	NE-CZ-NH1	5.96	123.28	120.30
1	D	428	VAL	CB-CA-C	-5.91	100.18	111.40
1	D	512	LEU	CA-CB-CG	5.81	128.66	115.30
1	A	428	VAL	CB-CA-C	-5.59	100.78	111.40
1	B	641	ARG	NE-CZ-NH1	5.53	123.06	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	510	ARG	NE-CZ-NH2	5.53	123.06	120.30
1	D	542	GLN	C-N-CD	-5.51	108.47	120.60
1	B	546	LEU	CA-CB-CG	5.47	127.89	115.30
1	D	482	LYS	CD-CE-NZ	5.28	123.85	111.70
1	C	437	ARG	NE-CZ-NH1	5.11	122.85	120.30
1	C	428	VAL	CB-CA-C	-5.04	101.83	111.40
1	B	550	LEU	CA-CB-CG	5.03	126.87	115.30
1	D	553	GLY	N-CA-C	-5.01	100.58	113.10

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	543	PRO	Peptide
1	A	544	ASP	Peptide
1	A	551	LEU	Peptide
1	B	374	THR	Peptide
1	B	375	GLU	Peptide
1	B	447	LEU	Peptide
1	C	374	THR	Peptide
1	D	552	THR	Peptide

## 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	2595	0	2563	60	1
1	B	2595	0	2563	71	0
1	C	2595	0	2563	69	0
1	D	2595	0	2563	79	1
2	A	27	0	22	2	0
2	B	27	0	22	2	0
2	C	27	0	22	4	0
2	D	27	0	22	3	0
3	A	9	0	0	0	0
3	B	12	0	0	0	0
3	C	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	17	0	0	0	0
All	All	10534	0	10340	270	1

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 13.

All (270) 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:D:561:THR:HG21	1:D:594:HIS:ND1	1.86	0.91
1:C:545:GLY:O	1:C:546:LEU:HG	1.70	0.90
1:A:542:GLN:HB2	1:A:543:PRO:CD	2.02	0.89
1:C:417:MET:HE2	1:C:426:CYS:SG	2.12	0.89
1:A:394:ARG:H	1:A:398:HIS:HD2	1.21	0.85
1:B:394:ARG:H	1:B:398:HIS:HD2	1.20	0.85
1:A:598:GLN:OE1	1:A:673:ARG:NH1	2.09	0.85
1:A:387:LYS:HD3	1:A:388:PRO:HD2	1.58	0.84
1:A:385:LYS:O	1:A:422:THR:HG21	1.76	0.84
1:A:542:GLN:HB2	1:A:543:PRO:HD3	1.58	0.83
1:A:453:VAL:HG23	1:A:532:LEU:O	1.81	0.80
1:D:507:HIS:CE1	1:D:575:ALA:HB1	2.16	0.80
1:A:483:GLU:OE2	1:A:666:ARG:NH1	2.14	0.80
1:C:561:THR:HG21	1:C:594:HIS:HB3	1.64	0.80
1:A:422:THR:HG22	1:A:424:PHE:H	1.46	0.79
1:D:576:LYS:NZ	1:D:641:ARG:O	2.16	0.78
1:C:417:MET:CE	1:C:426:CYS:SG	2.71	0.78
1:B:598:GLN:OE1	1:B:673:ARG:NH1	2.18	0.77
1:C:623:ALA:HB1	1:C:657:GLN:HE22	1.52	0.75
1:C:394:ARG:H	1:C:398:HIS:HD2	1.35	0.74
1:A:561:THR:HG21	1:A:594:HIS:ND1	2.03	0.74
1:B:332:SER:O	1:B:334:GLU:N	2.21	0.74
1:D:394:ARG:H	1:D:398:HIS:HD2	1.34	0.73
1:B:623:ALA:HB1	1:B:657:GLN:NE2	2.04	0.72
1:D:459:VAL:HG13	1:D:468:PHE:HE2	1.55	0.70
1:D:552:THR:OG1	1:D:553:GLY:HA3	1.92	0.70
1:C:483:GLU:OE2	1:C:666:ARG:NH1	2.24	0.70
1:D:417:MET:CE	1:D:426:CYS:SG	2.81	0.69
1:A:345:VAL:HG22	1:A:388:PRO:HB3	1.75	0.69
1:A:633:GLY:O	1:A:641:ARG:HD3	1.92	0.68
1:A:548:LYS:HE3	1:A:549:SER:N	2.08	0.68
1:D:417:MET:HE2	1:D:426:CYS:SG	2.34	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:613:PRO:HG3	1:D:636:LYS:HG3	1.77	0.67
1:D:411:PHE:HB2	1:D:413:GLU:HG3	1.77	0.66
1:D:483:GLU:OE2	1:D:666:ARG:NH1	2.29	0.66
1:B:560:GLU:OE2	1:C:337:LEU:HD23	1.95	0.65
1:C:542:GLN:HB2	1:C:543:PRO:C	2.18	0.63
1:A:568:VAL:CG2	1:A:605:CYS:HB2	2.30	0.62
1:C:496:LEU:HD13	1:C:651:VAL:HG11	1.81	0.62
1:D:477:LEU:HB2	1:D:521:VAL:HG22	1.82	0.62
1:B:416:ARG:CZ	1:B:471:LEU:HD21	2.31	0.61
1:C:339:HIS:HA	1:C:342:GLN:HE21	1.65	0.61
1:A:428:VAL:HG13	1:A:468:PHE:CE1	2.35	0.61
1:C:417:MET:HG2	1:C:428:VAL:CG2	2.30	0.61
1:B:512:LEU:HD23	1:B:512:LEU:C	2.21	0.61
1:C:459:VAL:HG13	1:C:468:PHE:HE2	1.65	0.61
1:C:345:VAL:HA	1:C:379:GLY:O	2.02	0.60
1:D:507:HIS:ND1	1:D:575:ALA:HB1	2.16	0.59
1:A:546:LEU:HG	1:A:574:ASP:HB3	1.85	0.58
1:D:345:VAL:HG22	1:D:388:PRO:HB3	1.85	0.58
1:C:453:VAL:HG23	1:C:532:LEU:O	2.04	0.58
1:C:353:LEU:HD11	1:C:458:ALA:HB3	1.86	0.57
1:D:394:ARG:H	1:D:398:HIS:CD2	2.20	0.57
1:A:394:ARG:H	1:A:398:HIS:CD2	2.12	0.57
1:B:545:GLY:O	1:B:546:LEU:HG	2.04	0.57
1:B:459:VAL:HG13	1:B:468:PHE:HE2	1.69	0.57
1:B:542:GLN:HB2	1:B:543:PRO:CD	2.35	0.56
1:B:405:ARG:HH11	1:B:406:LEU:H	1.53	0.56
1:A:550:LEU:O	1:A:551:LEU:HB2	2.06	0.56
1:B:561:THR:HG21	1:B:594:HIS:HB3	1.87	0.56
1:A:522:LEU:HD22	1:A:533:CYS:HB3	1.88	0.56
1:A:404:LEU:HD12	1:A:405:ARG:O	2.05	0.56
1:A:498:GLN:NE2	1:A:530:ALA:H	2.03	0.56
1:D:633:GLY:O	1:D:641:ARG:HD3	2.06	0.56
1:A:402:HIS:O	1:A:404:LEU:HA	2.06	0.56
1:B:345:VAL:HG12	1:B:378:GLU:HA	1.87	0.56
1:A:546:LEU:HB3	1:A:572:SER:OG	2.06	0.56
1:C:498:GLN:NE2	1:C:530:ALA:H	2.04	0.55
1:C:512:LEU:HD22	1:C:539:VAL:HG13	1.88	0.55
1:D:331:PHE:O	1:D:332:SER:C	2.44	0.55
1:C:345:VAL:HG22	1:C:388:PRO:HB3	1.88	0.55
1:D:512:LEU:HD13	1:D:539:VAL:HG22	1.89	0.55
1:A:331:PHE:CD2	1:A:332:SER:HA	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:611:GLU:HB3	1:B:612:PRO:HD2	1.88	0.54
1:A:568:VAL:HG21	1:A:605:CYS:HB2	1.89	0.54
1:D:428:VAL:HG13	1:D:468:PHE:CE1	2.42	0.54
1:D:498:GLN:NE2	1:D:530:ALA:H	2.06	0.54
1:C:545:GLY:O	1:C:546:LEU:CG	2.48	0.54
2:D:701:13V:H11	2:D:701:13V:C18	2.38	0.54
2:D:701:13V:H5	2:D:701:13V:H11	1.89	0.54
1:C:416:ARG:NE	1:C:471:LEU:HD21	2.22	0.54
1:B:345:VAL:HG21	1:B:393:TYR:HB2	1.89	0.53
1:C:619:PRO:HD2	1:C:622:CYS:SG	2.48	0.53
1:A:512:LEU:HD13	1:A:539:VAL:HG22	1.90	0.53
1:D:389:VAL:HG22	1:D:398:HIS:CE1	2.43	0.53
1:B:345:VAL:HA	1:B:379:GLY:O	2.08	0.53
1:B:385:LYS:HB3	1:B:424:PHE:CG	2.43	0.53
1:D:422:THR:HG22	1:D:424:PHE:HB2	1.91	0.53
1:A:408:ARG:HB3	2:A:701:13V:H10	1.90	0.53
1:C:445:ALA:HA	1:C:455:LEU:HD23	1.91	0.53
1:B:395:GLU:OE2	1:B:430:LYS:NZ	2.41	0.52
1:B:550:LEU:O	1:B:551:LEU:HB2	2.08	0.52
1:D:491:ARG:CZ	1:D:665:TRP:HB3	2.40	0.52
1:B:375:GLU:O	1:B:376:ASP:O	2.28	0.52
1:C:417:MET:HG2	1:C:428:VAL:HG22	1.91	0.52
1:B:454:PRO:HG2	1:B:470:GLU:OE1	2.09	0.52
1:B:507:HIS:CE1	1:B:575:ALA:HB1	2.45	0.52
1:B:637:GLU:OE2	1:B:639:ILE:HD13	2.10	0.52
1:A:411:PHE:HB2	1:A:413:GLU:HG3	1.91	0.52
1:C:394:ARG:H	1:C:398:HIS:CD2	2.23	0.52
1:A:624:PRO:HG2	1:A:657:GLN:HE22	1.75	0.51
1:C:399:TRP:CZ3	1:C:428:VAL:HG21	2.45	0.51
1:A:419:ASP:OD1	1:A:422:THR:HB	2.11	0.51
1:C:516:VAL:HG22	1:C:581:SER:HB2	1.92	0.51
1:B:336:TYR:HE2	1:D:555:TYR:O	1.94	0.51
1:C:446:GLY:O	1:C:447:LEU:C	2.48	0.51
1:B:456:TYR:OH	1:B:470:GLU:HG3	2.11	0.51
1:B:453:VAL:HG22	1:B:532:LEU:O	2.11	0.51
1:C:375:GLU:O	1:C:376:ASP:O	2.27	0.51
1:A:422:THR:CG2	1:A:424:PHE:HB2	2.41	0.50
1:D:477:LEU:CB	1:D:521:VAL:HG22	2.40	0.50
1:B:336:TYR:OH	1:D:552:THR:O	2.29	0.50
1:B:552:THR:O	1:C:336:TYR:OH	2.25	0.50
1:D:542:GLN:HG3	1:D:545:GLY:N	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:623:ALA:HB1	1:C:657:GLN:NE2	2.23	0.50
1:A:339:HIS:HA	1:A:342:GLN:HE21	1.77	0.50
1:D:561:THR:HG22	1:D:562:HIS:ND1	2.26	0.50
1:B:548:LYS:O	1:B:548:LYS:HG3	2.11	0.50
1:D:345:VAL:HA	1:D:379:GLY:O	2.11	0.50
1:D:568:VAL:CG2	1:D:605:CYS:HB2	2.41	0.50
1:D:375:GLU:HB2	1:D:462:GLY:HA2	1.94	0.50
1:A:331:PHE:HD2	1:A:332:SER:HA	1.76	0.49
1:B:408:ARG:HG3	2:B:701:13V:H5	1.94	0.49
1:D:603:PRO:HG2	1:D:606:LEU:HD12	1.93	0.49
1:C:544:ASP:O	1:C:545:GLY:O	2.30	0.49
1:D:522:LEU:HD22	1:D:533:CYS:HB3	1.94	0.49
1:B:331:PHE:O	1:B:332:SER:C	2.50	0.49
1:C:522:LEU:HD22	1:C:533:CYS:HB3	1.95	0.49
1:D:551:LEU:C	1:D:552:THR:HG22	2.32	0.49
1:A:345:VAL:CG2	1:A:388:PRO:HB3	2.41	0.49
1:A:552:THR:HG21	1:A:555:TYR:HB3	1.94	0.49
1:B:354:THR:HG21	1:D:601:ARG:HH22	1.76	0.49
1:B:394:ARG:H	1:B:398:HIS:CD2	2.12	0.49
1:C:550:LEU:HD23	1:C:551:LEU:HD23	1.94	0.49
1:C:561:THR:HG22	1:C:562:HIS:ND1	2.28	0.49
1:B:605:CYS:HA	1:B:608:ILE:HD12	1.95	0.48
1:D:546:LEU:HG	1:D:574:ASP:HB3	1.94	0.48
1:C:550:LEU:HD23	1:C:551:LEU:CD2	2.44	0.48
1:D:430:LYS:NZ	1:D:466:ASN:HD21	2.10	0.48
1:D:402:HIS:HB2	1:D:416:ARG:HG3	1.95	0.48
1:D:428:VAL:HG13	1:D:468:PHE:CD1	2.48	0.48
1:C:516:VAL:HG22	1:C:581:SER:CB	2.43	0.48
1:B:548:LYS:HE3	1:B:549:SER:N	2.28	0.48
1:C:613:PRO:HG3	1:C:636:LYS:HG3	1.96	0.48
1:C:385:LYS:HB3	1:C:424:PHE:CG	2.49	0.47
1:C:512:LEU:C	1:C:512:LEU:HD23	2.35	0.47
1:D:416:ARG:CZ	1:D:471:LEU:HD21	2.44	0.47
1:D:550:LEU:O	1:D:551:LEU:HB2	2.12	0.47
1:A:542:GLN:N	1:A:542:GLN:OE1	2.47	0.47
1:A:561:THR:HG22	1:A:562:HIS:ND1	2.29	0.47
1:A:548:LYS:HE2	1:A:550:LEU:H	1.80	0.47
1:D:623:ALA:HB1	1:D:657:GLN:NE2	2.30	0.47
1:A:512:LEU:HD23	1:A:514:GLY:N	2.29	0.47
1:B:659:GLY:O	1:B:662:LYS:HE3	2.14	0.47
1:C:498:GLN:HE22	1:C:530:ALA:H	1.61	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:403:GLN:HE21	1:C:404:LEU:HD22	1.80	0.47
1:B:601:ARG:HD2	1:C:509:ARG:NH2	2.29	0.47
1:D:402:HIS:O	1:D:404:LEU:HA	2.15	0.47
1:D:517:LYS:HB3	1:D:562:HIS:CD2	2.49	0.47
1:B:561:THR:HG21	1:B:594:HIS:CG	2.50	0.47
1:C:408:ARG:HG3	2:C:701:13V:H7	1.96	0.47
1:A:551:LEU:O	1:A:552:THR:HB	2.15	0.47
1:D:477:LEU:O	1:D:481:VAL:HG13	2.15	0.46
1:A:506:LEU:HD11	1:A:535:PHE:HE2	1.81	0.46
1:C:576:LYS:HA	1:C:579:VAL:HG12	1.96	0.46
1:B:354:THR:HG21	1:D:601:ARG:NH2	2.31	0.46
1:C:561:THR:HG21	1:C:594:HIS:CB	2.41	0.46
1:A:454:PRO:O	1:A:469:MET:HB2	2.16	0.46
1:A:548:LYS:HE2	1:A:550:LEU:N	2.30	0.46
2:A:701:13V:H11	2:A:701:13V:O2	2.15	0.46
1:B:550:LEU:HD23	1:B:550:LEU:O	2.16	0.46
1:D:417:MET:HE3	1:D:426:CYS:SG	2.56	0.46
1:A:496:LEU:HD22	1:A:500:LEU:HD11	1.98	0.46
1:A:548:LYS:C	1:A:548:LYS:HE3	2.37	0.46
1:B:331:PHE:CD2	1:D:551:LEU:HD11	2.50	0.46
1:C:629:ALA:HB2	1:C:650:LYS:HB3	1.97	0.46
1:B:428:VAL:HG13	1:B:468:PHE:CD1	2.51	0.45
1:B:561:THR:HG22	1:B:562:HIS:ND1	2.31	0.45
1:C:408:ARG:CG	2:C:701:13V:H7	2.46	0.45
1:C:548:LYS:HG3	1:C:550:LEU:O	2.16	0.45
1:D:332:SER:O	1:D:334:GLU:N	2.49	0.45
1:A:345:VAL:HA	1:A:379:GLY:O	2.16	0.45
1:C:405:ARG:HG3	1:C:406:LEU:H	1.80	0.45
1:B:389:VAL:HG13	1:B:398:HIS:ND1	2.32	0.45
1:C:607:LYS:HA	1:C:610:SER:HB3	1.99	0.45
1:D:403:GLN:HA	1:D:403:GLN:HE21	1.82	0.45
1:C:496:LEU:HD13	1:C:651:VAL:CG1	2.44	0.45
1:A:331:PHE:O	1:A:332:SER:C	2.54	0.44
1:B:374:THR:O	1:B:375:GLU:C	2.56	0.44
1:B:449:SER:C	1:B:451:ARG:H	2.20	0.44
1:C:375:GLU:O	1:C:376:ASP:C	2.56	0.44
2:C:701:13V:H11	2:C:701:13V:O2	2.17	0.44
1:B:623:ALA:HB1	1:B:657:GLN:HE22	1.79	0.44
1:C:405:ARG:HG2	1:C:416:ARG:NH1	2.32	0.44
1:B:527:GLY:HA3	1:B:665:TRP:CE3	2.53	0.44
1:D:449:SER:HB3	1:D:452:ILE:HG12	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:486:CYS:SG	1:B:670:LYS:HD2	2.58	0.44
1:A:414:VAL:HA	1:A:428:VAL:O	2.18	0.44
1:B:522:LEU:O	1:B:530:ALA:HA	2.17	0.44
1:B:534:ASP:H	2:B:701:13V:C4	2.31	0.43
1:D:542:GLN:HB2	1:D:543:PRO:C	2.38	0.43
1:A:417:MET:HG2	1:A:428:VAL:HG22	2.00	0.43
1:D:561:THR:HG21	1:D:594:HIS:CG	2.50	0.43
1:D:498:GLN:HE22	1:D:530:ALA:H	1.65	0.43
1:D:638:PRO:HA	1:D:641:ARG:HB3	2.00	0.43
1:D:594:HIS:HB2	1:D:597:THR:HG21	2.00	0.43
1:C:565:PRO:O	1:C:568:VAL:HG12	2.19	0.43
1:B:428:VAL:HG13	1:B:468:PHE:CE1	2.54	0.43
1:D:512:LEU:HD23	1:D:514:GLY:N	2.34	0.43
1:A:609:ALA:O	1:A:636:LYS:NZ	2.51	0.43
1:A:539:VAL:HG23	1:A:540:CYS:O	2.19	0.43
1:B:551:LEU:HD11	1:C:331:PHE:HD1	1.84	0.43
1:D:428:VAL:CG1	1:D:468:PHE:CE1	3.02	0.43
1:D:552:THR:OG1	1:D:553:GLY:CA	2.63	0.43
1:D:561:THR:CG2	1:D:562:HIS:ND1	2.82	0.42
1:B:440:GLU:HG3	1:B:535:PHE:C	2.40	0.42
1:B:353:LEU:HD11	1:B:458:ALA:HB3	1.99	0.42
1:D:603:PRO:HG2	1:D:606:LEU:CD1	2.48	0.42
1:B:632:GLU:O	1:B:642:VAL:HG22	2.19	0.42
1:B:412:GLY:HA3	1:C:388:PRO:O	2.20	0.42
1:D:353:LEU:HD11	1:D:458:ALA:HB3	2.00	0.42
1:D:414:VAL:HA	1:D:428:VAL:O	2.19	0.42
1:B:512:LEU:HD22	1:B:539:VAL:HG13	2.00	0.42
1:C:576:LYS:O	1:C:579:VAL:HG13	2.19	0.42
1:A:527:GLY:HA3	1:A:665:TRP:CD2	2.55	0.42
1:B:353:LEU:HA	1:B:380:VAL:HG13	2.01	0.42
1:D:552:THR:HG21	1:D:555:TYR:HB3	2.02	0.42
1:A:507:HIS:CD2	1:A:575:ALA:HB1	2.55	0.42
1:B:345:VAL:HG12	1:B:378:GLU:CA	2.50	0.42
1:D:576:LYS:HA	1:D:579:VAL:HG13	2.01	0.42
1:B:446:GLY:O	1:B:447:LEU:C	2.57	0.42
1:C:613:PRO:HB3	1:C:634:LEU:O	2.20	0.42
1:A:451:ARG:HG3	1:A:501:GLU:HB2	2.01	0.42
1:C:426:CYS:HB2	1:C:469:MET:O	2.19	0.42
1:D:338:VAL:HG22	1:D:355:SER:HB3	2.02	0.42
1:D:623:ALA:HB1	1:D:657:GLN:HE21	1.85	0.42
1:D:666:ARG:HD3	1:D:666:ARG:HA	1.81	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:417:MET:HG2	1:D:428:VAL:HG22	2.02	0.41
1:D:522:LEU:HD13	1:D:533:CYS:SG	2.61	0.41
1:B:503:LEU:HD13	1:B:532:LEU:HD11	2.02	0.41
1:A:422:THR:HG22	1:A:424:PHE:N	2.26	0.41
1:A:656:GLN:HE21	1:A:656:GLN:CA	2.32	0.41
1:C:497:GLY:HA2	1:C:651:VAL:HG21	2.02	0.41
1:C:496:LEU:HD22	1:C:500:LEU:CD1	2.51	0.41
1:B:349:GLN:HB3	1:B:351:HIS:CD2	2.55	0.41
1:B:623:ALA:HB1	1:B:657:GLN:HE21	1.83	0.41
1:A:548:LYS:HD2	1:A:570:GLY:C	2.41	0.41
1:B:390:ASP:O	1:B:391:TYR:HB2	2.20	0.41
1:C:416:ARG:HE	1:C:471:LEU:HD21	1.85	0.41
1:D:552:THR:HG21	1:D:555:TYR:CB	2.50	0.41
1:D:568:VAL:HG22	1:D:605:CYS:HB2	2.02	0.41
1:C:331:PHE:O	1:C:332:SER:C	2.59	0.41
1:A:428:VAL:HA	1:A:467:ILE:O	2.21	0.41
1:C:477:LEU:O	1:C:481:VAL:HG13	2.21	0.41
1:A:496:LEU:HD22	1:A:500:LEU:CD1	2.50	0.41
1:B:449:SER:HB2	1:B:505:TYR:CD1	2.56	0.41
1:D:512:LEU:CD2	1:D:512:LEU:C	2.89	0.41
1:D:602:GLY:HA2	1:D:603:PRO:C	2.41	0.41
1:B:335:GLU:O	1:B:339:HIS:HB2	2.20	0.40
1:B:396:GLU:N	1:B:399:TRP:O	2.54	0.40
1:C:544:ASP:C	1:C:545:GLY:O	2.59	0.40
1:C:534:ASP:H	2:C:701:13V:C5	2.34	0.40
1:D:453:VAL:HG22	1:D:532:LEU:O	2.21	0.40
1:D:548:LYS:HD2	1:D:570:GLY:C	2.40	0.40
1:D:522:LEU:HD21	2:D:701:13V:N2	2.36	0.40
1:A:453:VAL:HG21	1:A:533:CYS:HB3	2.03	0.40
1:B:509:ARG:HH22	1:D:601:ARG:HB3	1.85	0.40
1:B:551:LEU:HD11	1:C:331:PHE:CD1	2.56	0.40
1:C:637:GLU:HA	1:C:638:PRO:HD2	1.94	0.40
1:B:384:GLU:HB2	1:D:482:LYS:HE2	2.02	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:552:THR:O	1:D:336:TYR:OH[3_555]	2.12	0.08



## 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	331/356 (93%)	299 (90%)	23 (7%)	9 (3%)	5	19
1	B	331/356 (93%)	295 (89%)	26 (8%)	10 (3%)	4	17
1	C	331/356 (93%)	296 (89%)	24 (7%)	11 (3%)	4	15
1	D	331/356 (93%)	301 (91%)	24 (7%)	6 (2%)	8	29
All	All	1324/1424 (93%)	1191 (90%)	97 (7%)	36 (3%)	5	19

All (36) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	375	GLU
1	A	542	GLN
1	A	543	PRO
1	A	552	THR
1	B	333	VAL
1	B	376	ASP
1	B	542	GLN
1	B	543	PRO
1	B	551	LEU
1	C	332	SER
1	C	376	ASP
1	C	542	GLN
1	C	543	PRO
1	C	554	ASP
1	D	332	SER
1	D	333	VAL
1	D	543	PRO
1	A	376	ASP
1	B	332	SER
1	B	375	GLU
1	B	660	GLY
1	C	375	GLU

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Mol	Chain	Res	Type
1	C	447	LEU
1	C	545	GLY
1	C	546	LEU
1	D	375	GLU
1	B	549	SER
1	A	332	SER
1	A	551	LEU
1	C	534	ASP
1	D	542	GLN
1	D	551	LEU
1	A	549	SER
1	A	554	ASP
1	C	553	GLY
1	B	603	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	277/293 (94%)	234 (84%)	43 (16%)	2	8
1	B	277/293 (94%)	229 (83%)	48 (17%)	2	6
1	C	277/293 (94%)	237 (86%)	40 (14%)	3	9
1	D	277/293 (94%)	228 (82%)	49 (18%)	2	5
All	All	1108/1172 (94%)	928 (84%)	180 (16%)	2	7

All (180) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	337	LEU
1	A	344	SER
1	A	380	VAL
1	A	381	LEU
1	A	382	LEU
1	A	383	THR
1	A	389	VAL

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Mol	Chain	Res	Type
1	A	401	THR
1	A	404	LEU
1	A	405	ARG
1	A	413	GLU
1	A	417	MET
1	A	418	GLU
1	A	422	THR
1	A	428	VAL
1	A	433	LEU
1	A	448	THR
1	A	459	VAL
1	A	472	LEU
1	A	480	LEU
1	A	496	LEU
1	A	503	LEU
1	A	510	ARG
1	A	512	LEU
1	A	521	VAL
1	A	522	LEU
1	A	523	LEU
1	A	541	LEU
1	A	542	GLN
1	A	544	ASP
1	A	546	LEU
1	A	548	LYS
1	A	549	SER
1	A	550	LEU
1	A	561	THR
1	A	579	VAL
1	A	621	SER
1	A	641	ARG
1	A	650	LYS
1	A	651	VAL
1	A	656	GLN
1	A	658	VAL
1	A	663	SER
1	B	336	TYR
1	B	344	SER
1	B	349	GLN
1	B	354	THR
1	B	363	ARG
1	B	374	THR

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Mol	Chain	Res	Type
1	B	375	GLU
1	B	380	VAL
1	B	381	LEU
1	B	382	LEU
1	B	383	THR
1	B	386	LEU
1	B	389	VAL
1	B	403	GLN
1	B	404	LEU
1	B	405	ARG
1	B	406	LEU
1	B	417	MET
1	B	421	GLN
1	B	428	VAL
1	B	433	LEU
1	B	434	GLU
1	B	447	LEU
1	B	448	THR
1	B	453	VAL
1	B	459	VAL
1	B	480	LEU
1	B	483	GLU
1	B	496	LEU
1	B	503	LEU
1	B	509	ARG
1	B	510	ARG
1	B	512	LEU
1	B	516	VAL
1	B	521	VAL
1	B	522	LEU
1	B	523	LEU
1	B	542	GLN
1	B	548	LYS
1	B	550	LEU
1	B	552	THR
1	B	561	THR
1	B	568	VAL
1	B	605	CYS
1	B	621	SER
1	B	628	GLN
1	B	641	ARG
1	B	642	VAL

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Mol	Chain	Res	Type
1	C	337	LEU
1	C	345	VAL
1	C	380	VAL
1	C	381	LEU
1	C	382	LEU
1	C	383	THR
1	C	386	LEU
1	C	389	VAL
1	C	404	LEU
1	C	416	ARG
1	C	417	MET
1	C	433	LEU
1	C	447	LEU
1	C	480	LEU
1	C	481	VAL
1	C	496	LEU
1	C	503	LEU
1	C	509	ARG
1	C	510	ARG
1	C	512	LEU
1	C	516	VAL
1	C	521	VAL
1	C	522	LEU
1	C	523	LEU
1	C	546	LEU
1	C	548	LYS
1	C	561	THR
1	C	568	VAL
1	C	572	SER
1	C	577	VAL
1	C	579	VAL
1	C	604	LEU
1	C	605	CYS
1	C	606	LEU
1	C	621	SER
1	C	622	CYS
1	C	628	GLN
1	C	650	LYS
1	C	658	VAL
1	C	670	LYS
1	D	337	LEU
1	D	344	SER

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Mol	Chain	Res	Type
1	D	345	VAL
1	D	354	THR
1	D	363	ARG
1	D	375	GLU
1	D	376	ASP
1	D	380	VAL
1	D	381	LEU
1	D	382	LEU
1	D	383	THR
1	D	386	LEU
1	D	389	VAL
1	D	403	GLN
1	D	405	ARG
1	D	413	GLU
1	D	416	ARG
1	D	417	MET
1	D	428	VAL
1	D	433	LEU
1	D	448	THR
1	D	453	VAL
1	D	459	VAL
1	D	480	LEU
1	D	481	VAL
1	D	496	LEU
1	D	503	LEU
1	D	509	ARG
1	D	510	ARG
1	D	512	LEU
1	D	516	VAL
1	D	521	VAL
1	D	522	LEU
1	D	523	LEU
1	D	548	LYS
1	D	549	SER
1	D	550	LEU
1	D	552	THR
1	D	556	ILE
1	D	561	THR
1	D	568	VAL
1	D	572	SER
1	D	579	VAL
1	D	601	ARG

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Mol	Chain	Res	Type
1	D	607	LYS
1	D	625	LEU
1	D	650	LYS
1	D	668	GLU
1	D	670	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	342	GLN
1	A	398	HIS
1	A	403	GLN
1	A	466	ASN
1	A	498	GLN
1	A	628	GLN
1	A	631	GLN
1	A	656	GLN
1	B	342	GLN
1	B	398	HIS
1	B	403	GLN
1	B	415	HIS
1	B	466	ASN
1	B	498	GLN
1	B	529	HIS
1	B	631	GLN
1	B	657	GLN
1	C	342	GLN
1	C	398	HIS
1	C	403	GLN
1	C	498	GLN
1	C	542	GLN
1	C	631	GLN
1	D	342	GLN
1	D	398	HIS
1	D	403	GLN
1	D	415	HIS
1	D	421	GLN
1	D	466	ASN
1	D	498	GLN
1	D	529	HIS
1	D	542	GLN
1	D	631	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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

4 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	13V	D	701	-	27,29,29	1.84	7 (25%)	33,41,41	2.59	5 (15%)
2	13V	B	701	-	27,29,29	1.46	5 (18%)	33,41,41	1.65	6 (18%)
2	13V	C	701	-	27,29,29	1.55	6 (22%)	33,41,41	1.36	4 (12%)
2	13V	A	701	-	27,29,29	1.58	4 (14%)	33,41,41	1.52	6 (18%)

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	13V	D	701	-	-	6/10/15/15	0/3/3/3
2	13V	B	701	-	-	5/10/15/15	0/3/3/3
2	13V	C	701	-	-	4/10/15/15	0/3/3/3
2	13V	A	701	-	-	2/10/15/15	0/3/3/3



All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	701	13V	C5-C4	3.86	1.24	1.19
2	A	701	13V	O1-C1	-3.75	1.38	1.44
2	B	701	13V	O1-C1	-3.54	1.38	1.44
2	D	701	13V	O1-C1	-3.43	1.38	1.44
2	D	701	13V	C17-C14	3.38	1.45	1.40
2	A	701	13V	C15-N4	-3.32	1.31	1.35
2	C	701	13V	O1-C1	-3.19	1.39	1.44
2	A	701	13V	C17-C14	3.12	1.45	1.40
2	C	701	13V	C15-N4	-3.03	1.31	1.35
2	D	701	13V	C15-N4	-2.98	1.31	1.35
2	B	701	13V	C15-N4	-2.81	1.31	1.35
2	D	701	13V	O2-C17	2.78	1.43	1.37
2	D	701	13V	C1-C4	2.69	1.52	1.48
2	C	701	13V	C5-C4	2.56	1.22	1.19
2	D	701	13V	C16-C17	2.39	1.42	1.38
2	C	701	13V	C16-C17	2.36	1.42	1.38
2	C	701	13V	O2-C17	2.34	1.42	1.37
2	B	701	13V	C16-C17	2.28	1.42	1.38
2	C	701	13V	C1-C4	2.19	1.51	1.48
2	B	701	13V	C5-C4	2.10	1.22	1.19
2	B	701	13V	C17-C14	2.03	1.43	1.40
2	A	701	13V	C6-C5	-2.02	1.39	1.44

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	701	13V	C18-O2-C17	8.99	139.69	117.69
2	D	701	13V	C17-C14-N2	-7.95	114.86	121.61
2	D	701	13V	C14-N2-C15	5.97	123.22	117.22
2	B	701	13V	C18-O2-C17	5.89	132.10	117.69
2	A	701	13V	C18-O2-C17	5.45	131.03	117.69
2	C	701	13V	C18-O2-C17	5.03	129.99	117.69
2	B	701	13V	C17-C14-N2	-3.28	118.82	121.61
2	B	701	13V	C11-C14-N2	3.12	119.43	114.93
2	B	701	13V	C2-C1-C4	-2.99	105.18	109.92
2	A	701	13V	C11-C14-N2	2.78	118.94	114.93
2	D	701	13V	C13-C12-C9	-2.69	114.57	118.26
2	C	701	13V	C11-C14-N2	2.53	118.58	114.93
2	D	701	13V	C3-C1-C4	2.36	113.66	109.92
2	A	701	13V	C17-C14-N2	-2.27	119.68	121.61
2	A	701	13V	C7-C8-C9	-2.21	118.06	120.84
2	C	701	13V	C14-N2-C15	2.19	119.43	117.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	701	13V	C14-N2-C15	2.19	119.42	117.22
2	C	701	13V	C17-C14-N2	-2.16	119.77	121.61
2	B	701	13V	C20-O3-C19	2.05	125.73	112.96
2	A	701	13V	C13-C12-C9	-2.02	115.50	118.26
2	A	701	13V	C8-C9-C12	2.01	124.44	120.76

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	701	13V	C1-C4-C5-C6
2	A	701	13V	C1-C4-C5-C6
2	D	701	13V	C14-C17-O2-C18
2	B	701	13V	O2-C18-C19-O3
2	B	701	13V	C1-C4-C5-C6
2	D	701	13V	C16-C17-O2-C18
2	C	701	13V	C1-C4-C5-C6
2	C	701	13V	O2-C18-C19-O3
2	B	701	13V	C14-C17-O2-C18
2	C	701	13V	C18-C19-O3-C20
2	D	701	13V	C19-C18-O2-C17
2	B	701	13V	C16-C17-O2-C18
2	B	701	13V	C18-C19-O3-C20
2	D	701	13V	C18-C19-O3-C20
2	A	701	13V	C18-C19-O3-C20
2	D	701	13V	O2-C18-C19-O3
2	C	701	13V	C16-C17-O2-C18

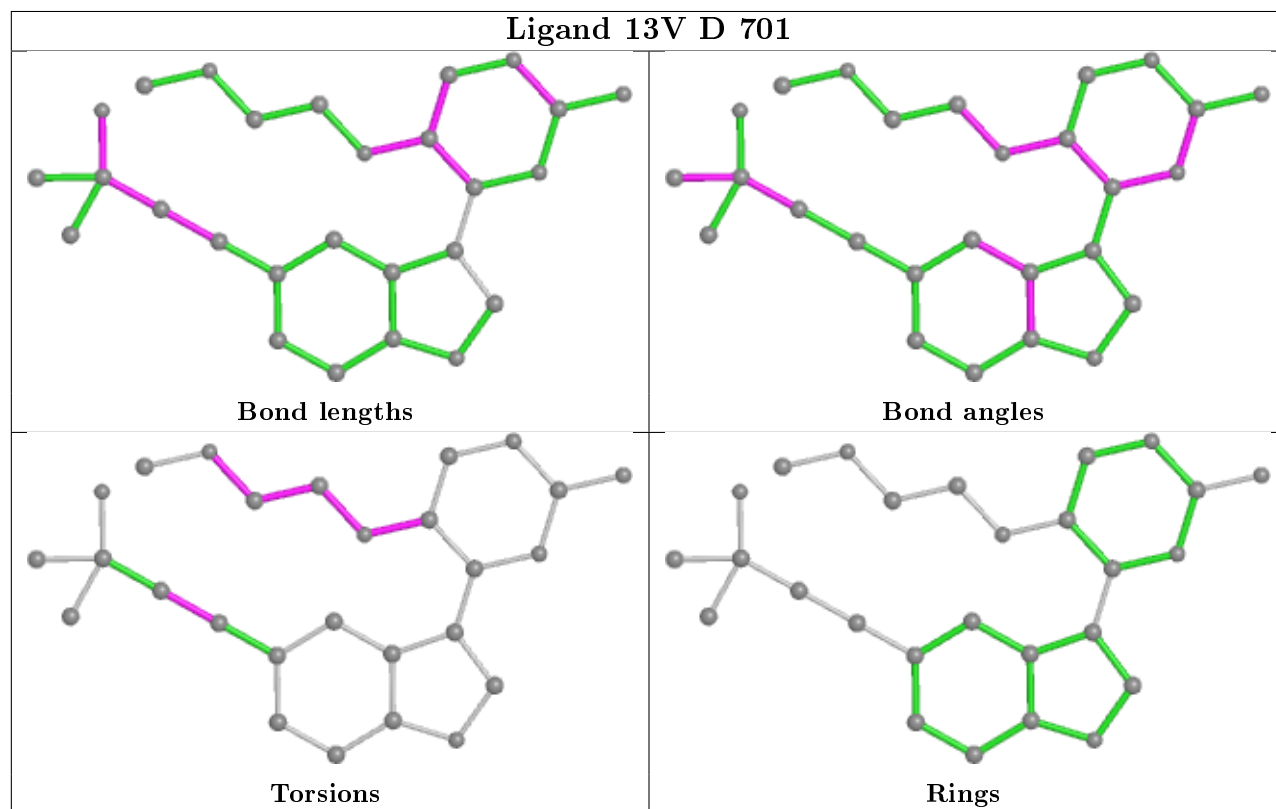
There are no ring outliers.

4 monomers are involved in 11 short contacts:

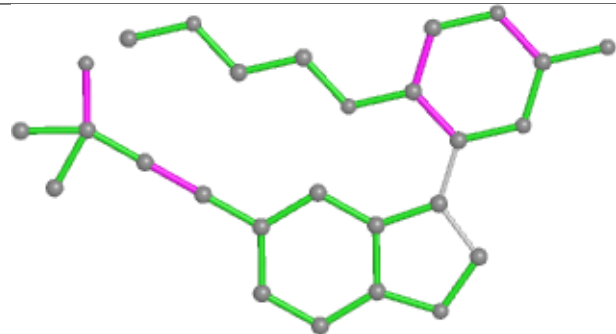
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	701	13V	3	0
2	B	701	13V	2	0
2	C	701	13V	4	0
2	A	701	13V	2	0

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.

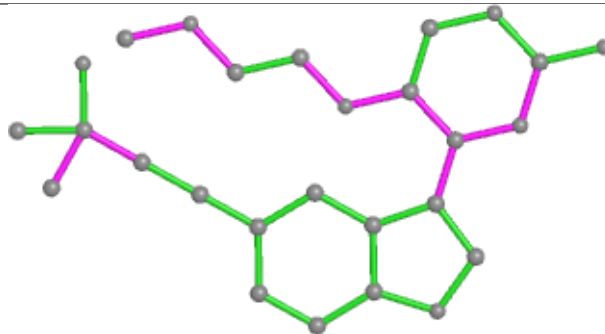
Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD 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.



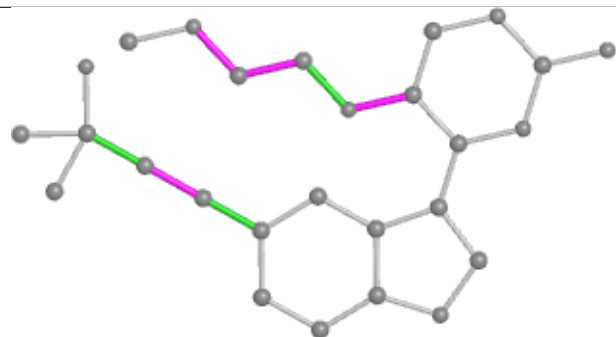
## Ligand 13V B 701



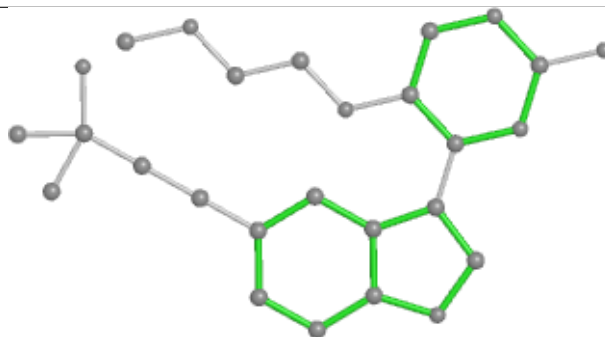
Bond lengths



Bond angles

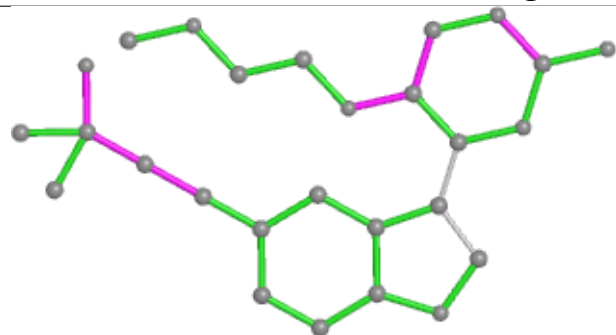


Torsions

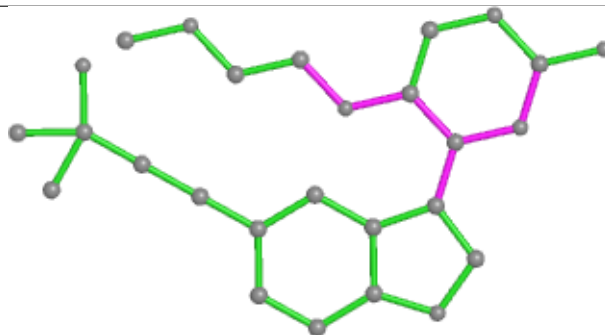


Rings

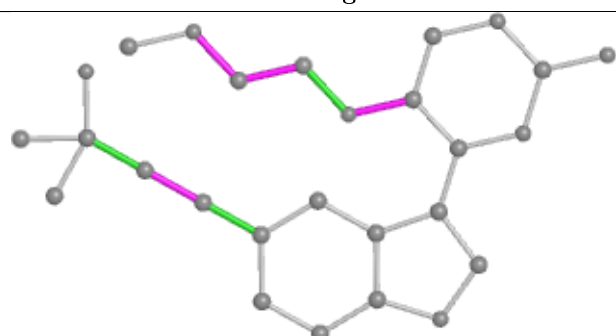
## Ligand 13V C 701



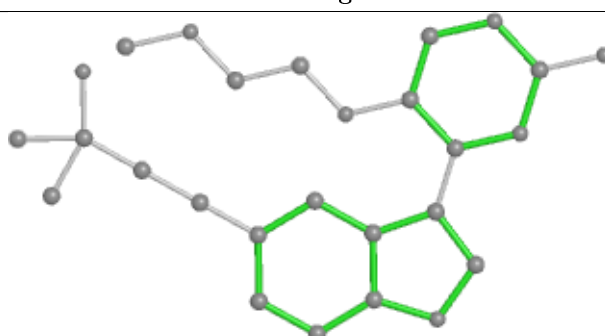
Bond lengths



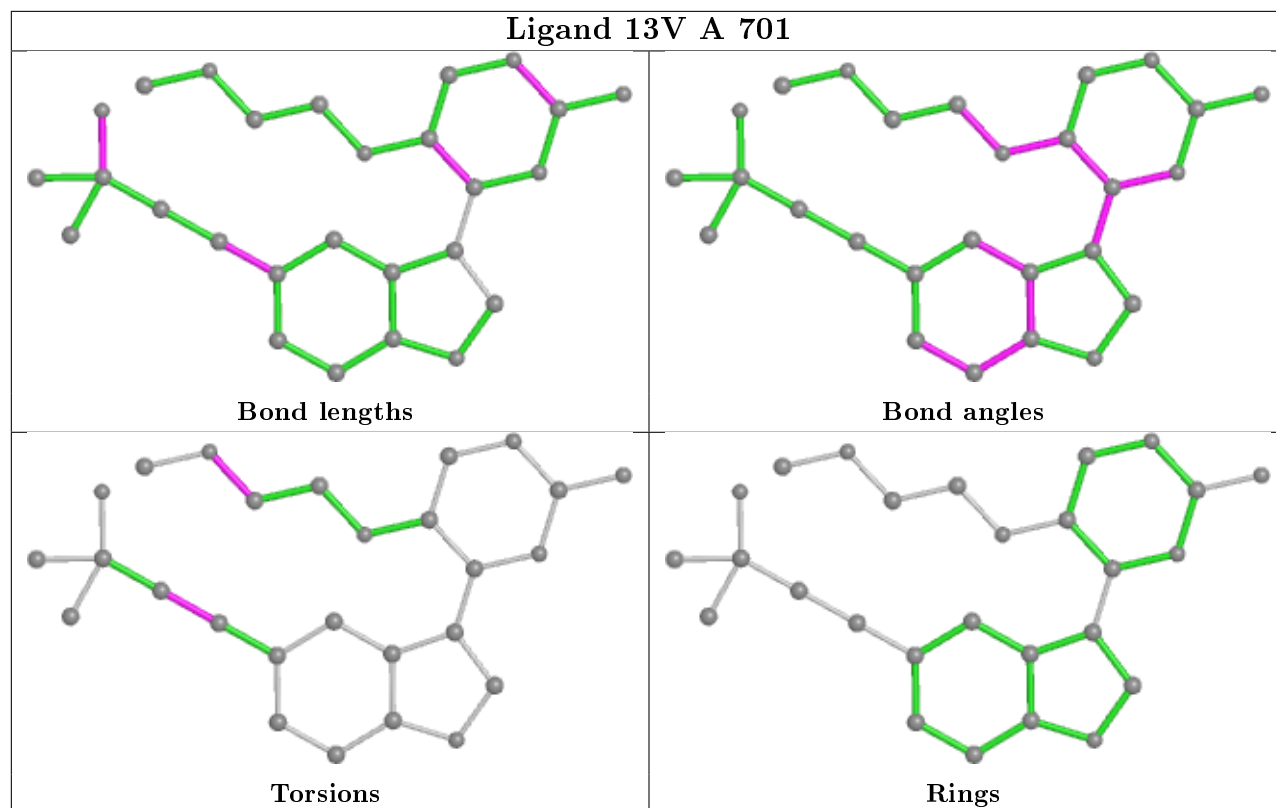
Bond angles



Torsions



Rings



## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	335/356 (94%)	-0.19	10 (2%) 50 45	17, 32, 72, 137	1 (0%)
1	B	335/356 (94%)	-0.04	14 (4%) 36 32	19, 36, 79, 116	1 (0%)
1	C	335/356 (94%)	0.00	21 (6%) 20 16	19, 38, 81, 137	1 (0%)
1	D	335/356 (94%)	-0.16	13 (3%) 39 35	18, 33, 73, 117	1 (0%)
All	All	1340/1424 (94%)	-0.10	58 (4%) 35 31	17, 35, 77, 137	4 (0%)

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	374	THR	7.9
1	A	374	THR	7.5
1	C	331	PHE	6.2
1	A	331	PHE	5.5
1	B	554	ASP	5.1
1	B	543	PRO	4.8
1	D	375	GLU	4.5
1	B	374	THR	4.2
1	D	331	PHE	4.2
1	A	375	GLU	4.1
1	A	363	ARG	3.7
1	C	544	ASP	3.5
1	C	599	PHE	3.5
1	C	674	HIS	3.4
1	B	363	ARG	3.3
1	C	374	THR	3.3
1	D	669	TYR	3.2
1	A	674	HIS	3.2
1	B	375	GLU	3.2
1	A	551	LEU	3.1
1	B	553	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	669	TYR	3.0
1	D	403	GLN	2.9
1	B	544	ASP	2.9
1	D	554	ASP	2.9
1	C	543	PRO	2.8
1	C	554	ASP	2.8
1	C	363	ARG	2.7
1	D	674	HIS	2.7
1	C	542	GLN	2.7
1	D	545	GLY	2.7
1	A	549	SER	2.6
1	C	375	GLU	2.4
1	A	554	ASP	2.3
1	C	671	GLU	2.3
1	B	602	GLY	2.3
1	C	672	PRO	2.3
1	C	602	GLY	2.3
1	B	669	TYR	2.3
1	B	674	HIS	2.3
1	D	400	ALA	2.2
1	C	448	THR	2.2
1	B	542	GLN	2.2
1	B	649	GLY	2.2
1	C	673	ARG	2.2
1	B	331	PHE	2.1
1	C	552	THR	2.1
1	D	363	ARG	2.1
1	A	671	GLU	2.1
1	C	668	GLU	2.1
1	B	400	ALA	2.1
1	C	601	ARG	2.1
1	D	544	ASP	2.0
1	A	447	LEU	2.0
1	D	510	ARG	2.0
1	C	545	GLY	2.0
1	D	552	THR	2.0
1	C	670	LYS	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates [i](#)

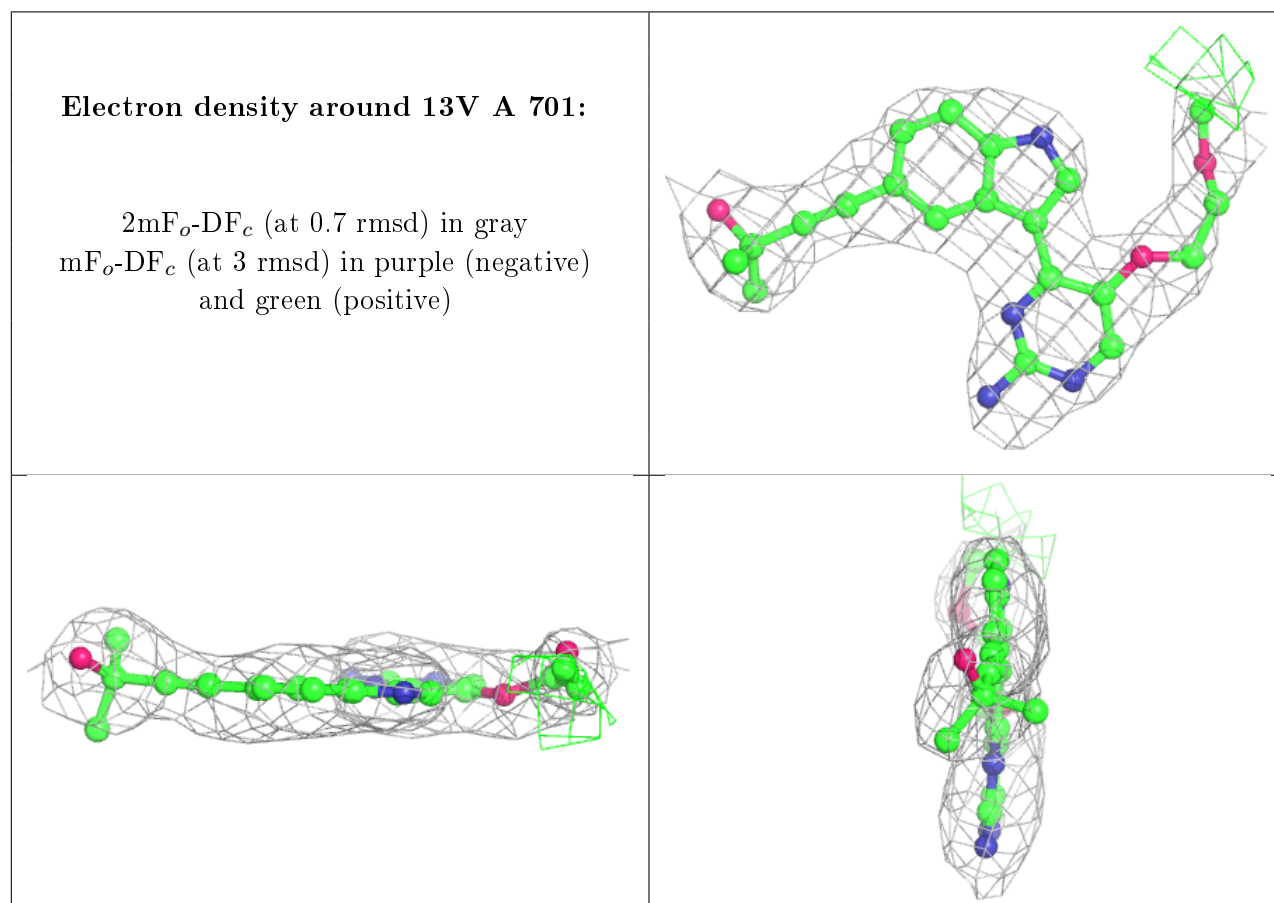
There are no carbohydrates in this entry.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	13V	A	701	27/27	0.94	0.19	29,33,50,53	0
2	13V	C	701	27/27	0.95	0.17	23,32,55,59	0
2	13V	D	701	27/27	0.96	0.20	21,28,43,48	0
2	13V	B	701	27/27	0.96	0.18	24,34,53,61	0

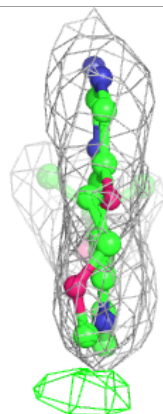
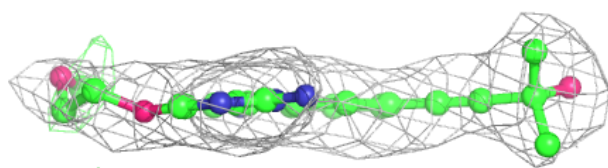
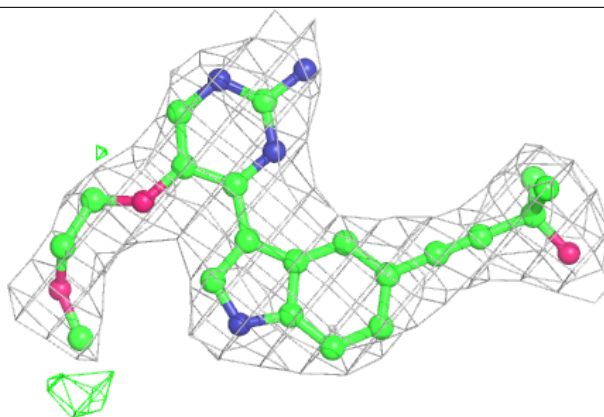
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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



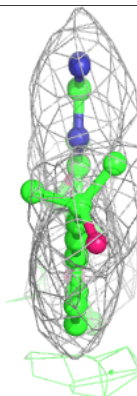
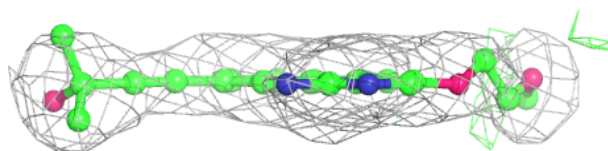
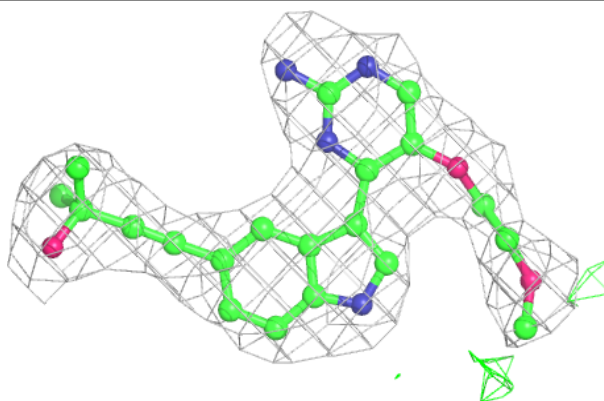


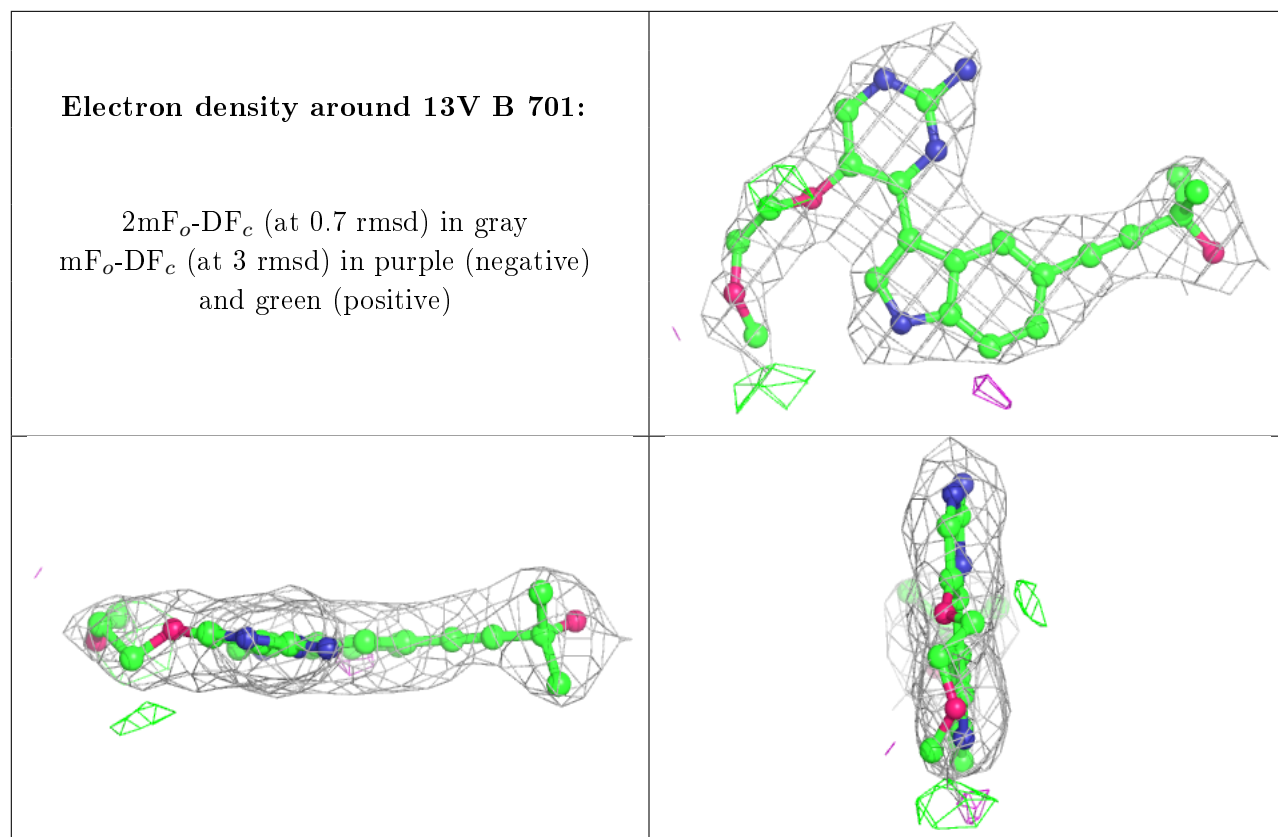
**Electron density around 13V C 701:**

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

**Electron density around 13V D 701:**

$2mF_o-DF_c$  (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.