



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2026 – 02:42 AM UTC

PDB ID : 9C9V / pdb_00009c9v
Title : HBV capsid with compound 3i
Authors : Olland, A.M.; Suto, R.K.; Fontano, E.; Colussi, T.
Deposited on : 2024-06-16
Resolution : 2.63 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

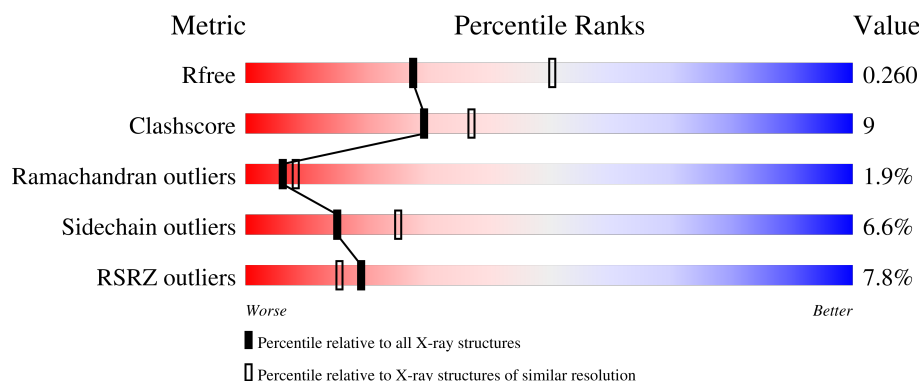
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.63 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	2053 (2.66-2.62)
Clashscore	190562	2097 (2.66-2.62)
Ramachandran outliers	187476	2066 (2.66-2.62)
Sidechain outliers	187428	2066 (2.66-2.62)
RSRZ outliers	180081	2052 (2.66-2.62)

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	167	5% 60% 21% .. 17%
1	B	167	7% 57% 19% .. 19%
1	C	167	6% 55% 22% .. 21%
1	D	167	8% 57% 19% .. 22%
1	E	167	5% 60% 15% .. 22%

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Mol	Chain	Length	Quality of chain
1	F	167	6% 49% 22% • 25%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6459 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 Capsid protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	138	Total	C	N	O	S	0	1	0
			1093	708	183	197	5			
1	B	135	Total	C	N	O	S	0	0	0
			1058	688	176	189	5			
1	C	132	Total	C	N	O	S	0	1	0
			1043	675	174	189	5			
1	D	131	Total	C	N	O	S	0	0	0
			1032	673	171	183	5			
1	E	130	Total	C	N	O	S	0	2	0
			1048	680	172	191	5			
1	F	126	Total	C	N	O	S	0	2	0
			1011	655	168	183	5			

There are 114 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	MET	-	initiating methionine	UNP L7R9I1
A	-1	GLY	-	expression tag	UNP L7R9I1
A	0	SER	-	expression tag	UNP L7R9I1
A	132	ALA	TYR	engineered mutation	UNP L7R9I1
A	150	LYS	-	expression tag	UNP L7R9I1
A	151	LEU	-	expression tag	UNP L7R9I1
A	152	GLU	-	expression tag	UNP L7R9I1
A	153	ASN	-	expression tag	UNP L7R9I1
A	154	LEU	-	expression tag	UNP L7R9I1
A	155	TYR	-	expression tag	UNP L7R9I1
A	156	PHE	-	expression tag	UNP L7R9I1
A	157	GLN	-	expression tag	UNP L7R9I1
A	158	GLY	-	expression tag	UNP L7R9I1
A	159	HIS	-	expression tag	UNP L7R9I1
A	160	HIS	-	expression tag	UNP L7R9I1
A	161	HIS	-	expression tag	UNP L7R9I1
A	162	HIS	-	expression tag	UNP L7R9I1

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Chain	Residue	Modelled	Actual	Comment	Reference
A	163	HIS	-	expression tag	UNP L7R9I1
A	164	HIS	-	expression tag	UNP L7R9I1
B	-2	MET	-	initiating methionine	UNP L7R9I1
B	-1	GLY	-	expression tag	UNP L7R9I1
B	0	SER	-	expression tag	UNP L7R9I1
B	132	ALA	TYR	engineered mutation	UNP L7R9I1
B	150	LYS	-	expression tag	UNP L7R9I1
B	151	LEU	-	expression tag	UNP L7R9I1
B	152	GLU	-	expression tag	UNP L7R9I1
B	153	ASN	-	expression tag	UNP L7R9I1
B	154	LEU	-	expression tag	UNP L7R9I1
B	155	TYR	-	expression tag	UNP L7R9I1
B	156	PHE	-	expression tag	UNP L7R9I1
B	157	GLN	-	expression tag	UNP L7R9I1
B	158	GLY	-	expression tag	UNP L7R9I1
B	159	HIS	-	expression tag	UNP L7R9I1
B	160	HIS	-	expression tag	UNP L7R9I1
B	161	HIS	-	expression tag	UNP L7R9I1
B	162	HIS	-	expression tag	UNP L7R9I1
B	163	HIS	-	expression tag	UNP L7R9I1
B	164	HIS	-	expression tag	UNP L7R9I1
C	-2	MET	-	initiating methionine	UNP L7R9I1
C	-1	GLY	-	expression tag	UNP L7R9I1
C	0	SER	-	expression tag	UNP L7R9I1
C	132	ALA	TYR	engineered mutation	UNP L7R9I1
C	150	LYS	-	expression tag	UNP L7R9I1
C	151	LEU	-	expression tag	UNP L7R9I1
C	152	GLU	-	expression tag	UNP L7R9I1
C	153	ASN	-	expression tag	UNP L7R9I1
C	154	LEU	-	expression tag	UNP L7R9I1
C	155	TYR	-	expression tag	UNP L7R9I1
C	156	PHE	-	expression tag	UNP L7R9I1
C	157	GLN	-	expression tag	UNP L7R9I1
C	158	GLY	-	expression tag	UNP L7R9I1
C	159	HIS	-	expression tag	UNP L7R9I1
C	160	HIS	-	expression tag	UNP L7R9I1
C	161	HIS	-	expression tag	UNP L7R9I1
C	162	HIS	-	expression tag	UNP L7R9I1
C	163	HIS	-	expression tag	UNP L7R9I1
C	164	HIS	-	expression tag	UNP L7R9I1
D	-2	MET	-	initiating methionine	UNP L7R9I1
D	-1	GLY	-	expression tag	UNP L7R9I1

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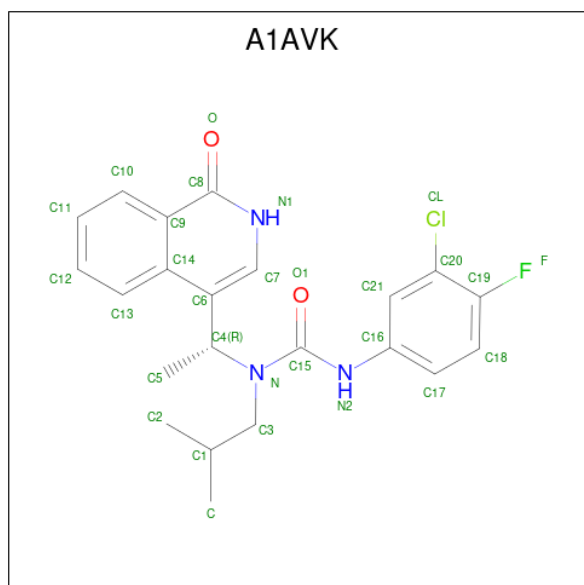
Chain	Residue	Modelled	Actual	Comment	Reference
D	0	SER	-	expression tag	UNP L7R9I1
D	132	ALA	TYR	engineered mutation	UNP L7R9I1
D	150	LYS	-	expression tag	UNP L7R9I1
D	151	LEU	-	expression tag	UNP L7R9I1
D	152	GLU	-	expression tag	UNP L7R9I1
D	153	ASN	-	expression tag	UNP L7R9I1
D	154	LEU	-	expression tag	UNP L7R9I1
D	155	TYR	-	expression tag	UNP L7R9I1
D	156	PHE	-	expression tag	UNP L7R9I1
D	157	GLN	-	expression tag	UNP L7R9I1
D	158	GLY	-	expression tag	UNP L7R9I1
D	159	HIS	-	expression tag	UNP L7R9I1
D	160	HIS	-	expression tag	UNP L7R9I1
D	161	HIS	-	expression tag	UNP L7R9I1
D	162	HIS	-	expression tag	UNP L7R9I1
D	163	HIS	-	expression tag	UNP L7R9I1
D	164	HIS	-	expression tag	UNP L7R9I1
E	-2	MET	-	initiating methionine	UNP L7R9I1
E	-1	GLY	-	expression tag	UNP L7R9I1
E	0	SER	-	expression tag	UNP L7R9I1
E	132	ALA	TYR	engineered mutation	UNP L7R9I1
E	150	LYS	-	expression tag	UNP L7R9I1
E	151	LEU	-	expression tag	UNP L7R9I1
E	152	GLU	-	expression tag	UNP L7R9I1
E	153	ASN	-	expression tag	UNP L7R9I1
E	154	LEU	-	expression tag	UNP L7R9I1
E	155	TYR	-	expression tag	UNP L7R9I1
E	156	PHE	-	expression tag	UNP L7R9I1
E	157	GLN	-	expression tag	UNP L7R9I1
E	158	GLY	-	expression tag	UNP L7R9I1
E	159	HIS	-	expression tag	UNP L7R9I1
E	160	HIS	-	expression tag	UNP L7R9I1
E	161	HIS	-	expression tag	UNP L7R9I1
E	162	HIS	-	expression tag	UNP L7R9I1
E	163	HIS	-	expression tag	UNP L7R9I1
E	164	HIS	-	expression tag	UNP L7R9I1
F	-2	MET	-	initiating methionine	UNP L7R9I1
F	-1	GLY	-	expression tag	UNP L7R9I1
F	0	SER	-	expression tag	UNP L7R9I1
F	132	ALA	TYR	engineered mutation	UNP L7R9I1
F	150	LYS	-	expression tag	UNP L7R9I1
F	151	LEU	-	expression tag	UNP L7R9I1

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Chain	Residue	Modelled	Actual	Comment	Reference
F	152	GLU	-	expression tag	UNP L7R9I1
F	153	ASN	-	expression tag	UNP L7R9I1
F	154	LEU	-	expression tag	UNP L7R9I1
F	155	TYR	-	expression tag	UNP L7R9I1
F	156	PHE	-	expression tag	UNP L7R9I1
F	157	GLN	-	expression tag	UNP L7R9I1
F	158	GLY	-	expression tag	UNP L7R9I1
F	159	HIS	-	expression tag	UNP L7R9I1
F	160	HIS	-	expression tag	UNP L7R9I1
F	161	HIS	-	expression tag	UNP L7R9I1
F	162	HIS	-	expression tag	UNP L7R9I1
F	163	HIS	-	expression tag	UNP L7R9I1
F	164	HIS	-	expression tag	UNP L7R9I1

- Molecule 2 is N'-(3-chloro-4-fluorophenyl)-N-(2-methylpropyl)-N-[(1R)-1-(1-oxo-1,2-dihydroisoquinolin-4-yl)ethyl]urea (CCD ID: A1AVK) (formula: C₂₂H₂₃ClFN₃O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	Cl	F	N	O	0	0
			29	22	1	1	3	2		
2	B	1	Total	C	Cl	F	N	O	0	0
			29	22	1	1	3	2		
2	C	1	Total	C	Cl	F	N	O	0	0
			29	22	1	1	3	2		
2	D	1	Total	C	Cl	F	N	O	0	0
			29	22	1	1	3	2		

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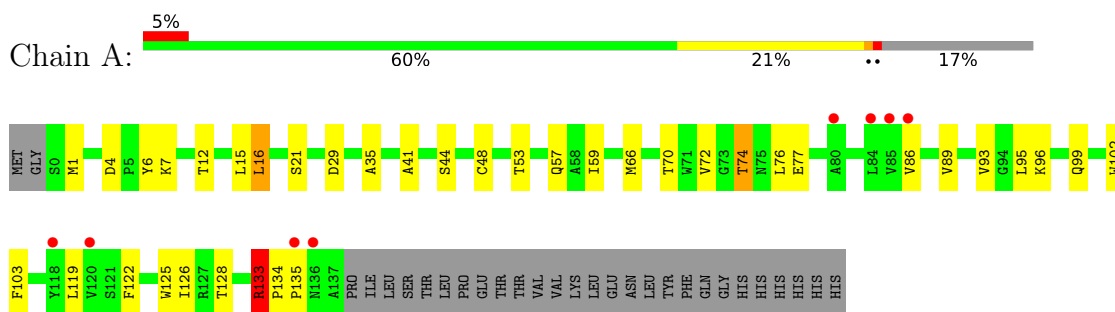
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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	E	1	Total	C	Cl	F	N	O	0	0
			29	22	1	1	3	2		
2	F	1	Total	C	Cl	F	N	O	0	0
			29	22	1	1	3	2		

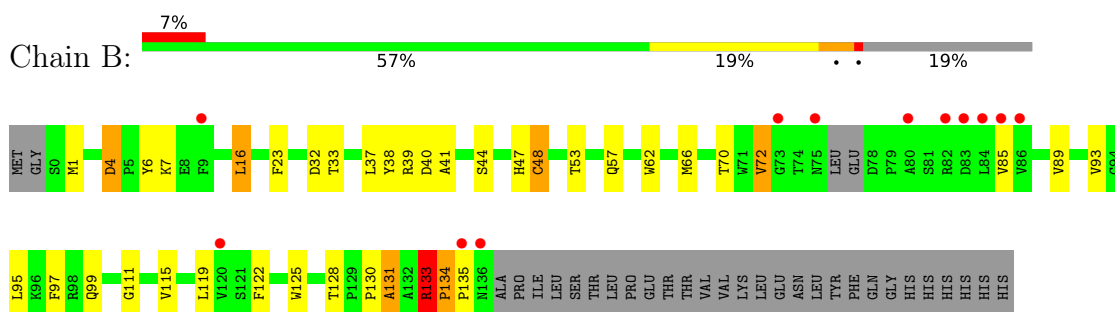
3 Residue-property plots [i](#)

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.

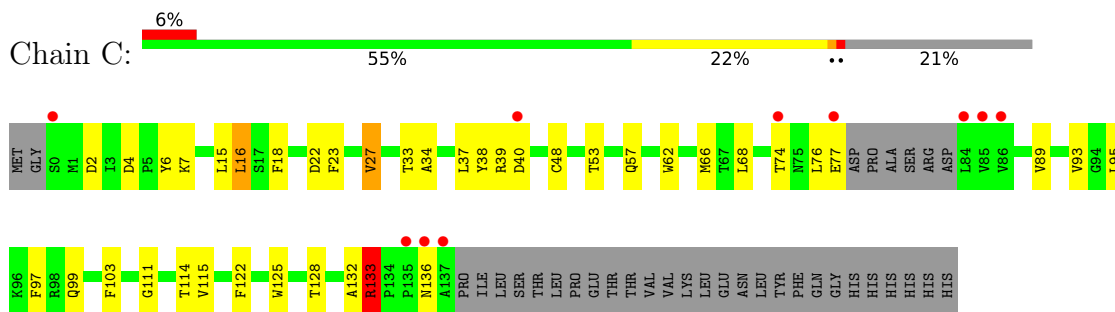
- Molecule 1: Capsid protein



- Molecule 1: Capsid protein

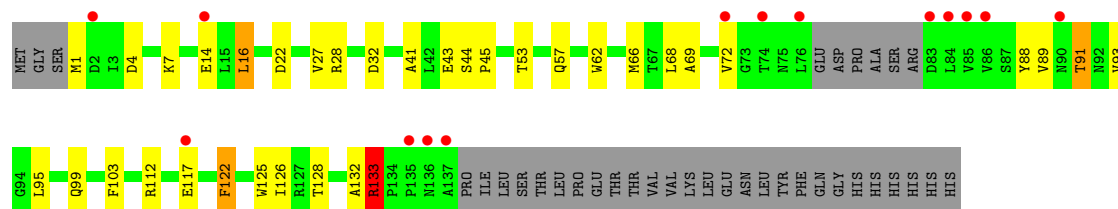


- Molecule 1: Capsid protein

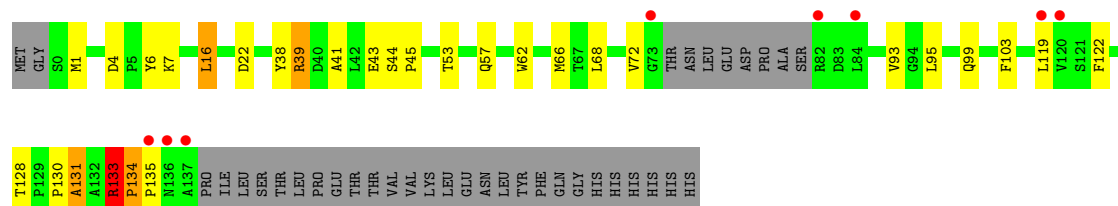


- Molecule 1: Capsid protein

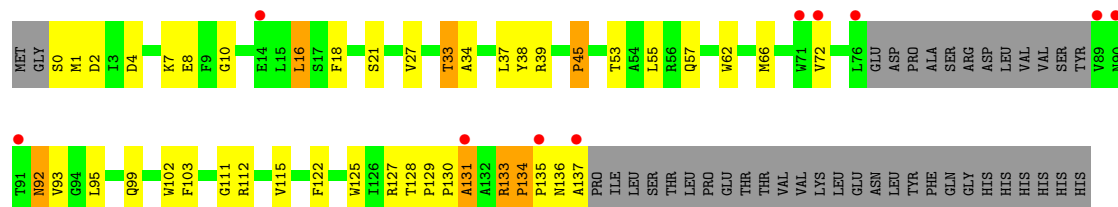




• Molecule 1: Capsid protein



• Molecule 1: Capsid protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	151.94Å 87.92Å 108.50Å 90.00° 102.75° 90.00°	Depositor
Resolution (Å)	48.41 – 2.63 48.41 – 2.63	Depositor EDS
% Data completeness (in resolution range)	98.6 (48.41-2.63) 98.6 (48.41-2.63)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.224 , 0.258 0.231 , 0.260	Depositor DCC
R_{free} test set	2055 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	56.5	Xtriage
Anisotropy	0.274	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 42.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6459	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.33% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: A1AVK

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	0.62	0/1126	1.27	4/1543 (0.3%)
1	B	0.63	0/1090	1.26	6/1494 (0.4%)
1	C	0.61	0/1074	1.25	7/1470 (0.5%)
1	D	0.61	0/1063	1.27	8/1455 (0.5%)
1	E	0.62	0/1079	1.22	4/1476 (0.3%)
1	F	0.62	0/1041	1.27	4/1423 (0.3%)
All	All	0.62	0/6473	1.26	33/8861 (0.4%)

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	B	0	1
1	E	0	2
All	All	0	3

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	129	PRO	N-CA-CB	8.25	107.81	103.19
1	F	122	PHE	CA-CB-CG	-7.86	105.94	113.80
1	D	27	VAL	N-CA-CB	7.00	119.53	110.57
1	E	122	PHE	CA-CB-CG	-6.86	106.94	113.80
1	D	122	PHE	CA-CB-CG	-6.78	107.02	113.80
1	B	133	ARG	CG-CD-NE	6.69	126.72	112.00
1	C	27	VAL	N-CA-CB	6.28	118.61	110.57
1	B	4	ASP	CA-CB-CG	6.27	118.87	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	133	ARG	CB-CA-C	6.20	117.30	108.68
1	B	97	PHE	CA-CB-CG	-6.07	107.73	113.80
1	C	97	PHE	CA-CB-CG	-5.93	107.87	113.80
1	C	122	PHE	CA-CB-CG	-5.84	107.96	113.80
1	C	22	ASP	CA-CB-CG	5.78	118.38	112.60
1	A	59	ILE	O-C-N	5.75	127.55	121.91
1	E	135	PRO	N-CA-C	5.69	121.03	113.40
1	A	29	ASP	CA-CB-CG	5.64	118.24	112.60
1	D	27	VAL	N-CA-C	-5.64	105.23	110.53
1	C	114	THR	CA-CB-OG1	-5.61	101.18	109.60
1	A	133	ARG	CB-CA-C	5.61	117.44	109.08
1	D	22	ASP	CA-CB-CG	5.60	118.20	112.60
1	A	12	THR	CA-CB-OG1	-5.58	101.23	109.60
1	D	32	ASP	CA-CB-CG	5.56	118.16	112.60
1	B	32	ASP	CA-CB-CG	5.43	118.03	112.60
1	E	22	ASP	CA-CB-CG	5.41	118.01	112.60
1	E	134	PRO	N-CA-C	5.33	117.21	110.70
1	B	122	PHE	CA-CB-CG	-5.25	108.55	113.80
1	B	70	THR	CA-CB-OG1	-5.18	101.83	109.60
1	D	117	GLU	N-CA-CB	5.17	117.68	109.82
1	F	55	LEU	N-CA-CB	-5.13	102.58	110.12
1	D	133	ARG	CB-CA-C	5.08	116.03	108.87
1	F	33	THR	CA-CB-OG1	-5.07	102.00	109.60
1	C	74	THR	CA-CB-OG1	-5.07	102.00	109.60
1	D	14	GLU	CB-CG-CD	5.02	121.13	112.60

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	133	ARG	Sidechain
1	E	133	ARG	Sidechain
1	E	39	ARG	Sidechain

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	1093	0	1064	17	0
1	B	1058	0	1023	22	0
1	C	1043	0	994	20	0
1	D	1032	0	1003	17	0
1	E	1048	0	1014	14	0
1	F	1011	0	973	22	0
2	A	29	0	0	0	0
2	B	29	0	0	1	0
2	C	29	0	0	0	0
2	D	29	0	0	1	0
2	E	29	0	0	0	0
2	F	29	0	0	0	0
All	All	6459	0	6071	108	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 9.

All (108) 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:E:95:LEU:O	1:E:99:GLN:HG3	1.91	0.70
1:B:125:TRP:CZ2	1:B:133:ARG:HD2	2.28	0.69
1:A:4:ASP:HB3	1:A:7:LYS:HB2	1.78	0.66
1:C:128:THR:O	1:C:133:ARG:NH1	2.30	0.65
1:E:62:TRP:CZ2	1:E:66:MET:HE3	2.32	0.65
1:B:128:THR:O	1:B:133:ARG:NH1	2.29	0.64
1:F:95:LEU:O	1:F:99:GLN:HG3	1.98	0.64
1:B:66:MET:CE	1:B:89:VAL:HG11	2.28	0.63
1:A:95:LEU:O	1:A:99:GLN:HG3	1.99	0.62
1:C:125:TRP:CE2	1:C:133:ARG:HD2	2.35	0.62
1:D:53:THR:O	1:D:57:GLN:HG2	2.00	0.62
1:D:66:MET:HE1	1:D:89:VAL:HG11	1.82	0.61
1:F:128:THR:O	1:F:133:ARG:NH1	2.34	0.61
1:A:66:MET:HE1	1:A:89:VAL:HG11	1.82	0.60
1:C:53:THR:O	1:C:57:GLN:HG2	2.01	0.60
1:F:111:GLY:O	1:F:115:VAL:HG23	2.02	0.59
1:D:69:ALA:HA	1:D:72:VAL:HG12	1.85	0.59
1:C:2:ASP:N	1:D:43:GLU:OE2	2.28	0.58
1:D:88:TYR:HA	1:D:91:THR:OG1	2.04	0.58
1:F:53:THR:O	1:F:57:GLN:HG2	2.04	0.58
1:D:95:LEU:O	1:D:99:GLN:HG3	2.04	0.57
1:B:53:THR:O	1:B:57:GLN:HG2	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:72:VAL:HG11	1:B:85:VAL:HG11	1.88	0.56
1:A:70:THR:O	1:A:74:THR:HG22	2.05	0.56
1:F:92:ASN:N	1:F:92:ASN:OD1	2.39	0.56
1:B:62:TRP:CZ2	1:B:66:MET:HE3	2.42	0.55
1:B:66:MET:HE1	1:B:89:VAL:HG11	1.88	0.55
1:B:95:LEU:O	1:B:99:GLN:HG3	2.07	0.54
1:E:53:THR:O	1:E:57:GLN:HG2	2.07	0.54
1:A:125:TRP:CE2	1:A:133:ARG:HD2	2.43	0.54
1:F:4:ASP:HB3	1:F:7:LYS:HB2	1.90	0.54
1:E:130:PRO:O	1:E:131:ALA:C	2.51	0.54
1:C:4:ASP:HB3	1:C:7:LYS:HB2	1.89	0.54
1:D:68:LEU:HD23	1:D:68:LEU:O	2.09	0.53
1:A:53:THR:O	1:A:57:GLN:HG2	2.09	0.52
1:C:111:GLY:O	1:C:115:VAL:HG23	2.08	0.52
1:D:41:ALA:O	1:D:44:SER:HB3	2.08	0.52
1:E:16:LEU:HD21	1:E:103:PHE:HB2	1.92	0.52
1:C:16:LEU:HD21	1:C:103:PHE:HB2	1.91	0.52
1:D:16:LEU:HD21	1:D:103:PHE:HB2	1.91	0.52
1:B:119:LEU:C	1:B:119:LEU:HD23	2.34	0.52
1:C:16:LEU:HB3	1:C:99:GLN:OE1	2.10	0.51
1:B:23:PHE:C	1:B:23:PHE:CD1	2.88	0.51
1:C:95:LEU:O	1:C:99:GLN:HG3	2.11	0.51
1:D:4:ASP:HB3	1:D:7:LYS:HB2	1.93	0.51
1:A:128:THR:O	1:A:133:ARG:NH1	2.44	0.50
1:B:4:ASP:HB3	1:B:7:LYS:HB2	1.94	0.50
1:C:34:ALA:O	1:C:38:TYR:HB2	2.12	0.50
1:B:41:ALA:O	1:B:44:SER:HB3	2.11	0.50
1:B:111:GLY:O	1:B:115:VAL:HG23	2.11	0.49
1:C:6:TYR:CZ	1:C:16:LEU:HD22	2.48	0.48
1:F:125:TRP:CE2	1:F:133:ARG:HD2	2.48	0.48
1:B:134:PRO:HB3	1:B:135:PRO:HD2	1.96	0.48
1:B:16:LEU:HB3	1:B:99:GLN:OE1	2.13	0.47
1:F:10:GLY:HA3	1:F:112:ARG:NH2	2.29	0.47
1:C:76:LEU:O	1:C:77:GLU:CB	2.61	0.47
1:A:41:ALA:O	1:A:44:SER:HB3	2.14	0.47
1:C:68:LEU:HD22	1:D:68:LEU:HD22	1.97	0.47
1:C:62:TRP:CZ2	1:C:66:MET:HE3	2.51	0.46
1:B:6:TYR:CZ	1:B:16:LEU:HD22	2.51	0.46
1:A:16:LEU:HB3	1:A:99:GLN:OE1	2.16	0.46
1:F:130:PRO:O	1:F:131:ALA:C	2.58	0.46
1:D:112:ARG:CB	1:D:112:ARG:HH11	2.28	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:201:A1AVK:O1	2:B:201:A1AVK:C17	2.63	0.46
1:A:119:LEU:HD23	1:A:119:LEU:C	2.41	0.46
1:C:16:LEU:HD23	1:C:99:GLN:HB3	1.98	0.46
1:E:128:THR:O	1:E:133:ARG:NH1	2.49	0.46
1:C:15:LEU:O	1:C:18:PHE:HB3	2.16	0.45
1:E:41:ALA:O	1:E:44:SER:HB3	2.16	0.45
1:E:119:LEU:HD23	1:E:119:LEU:C	2.41	0.45
1:E:43:GLU:OE2	1:F:2:ASP:N	2.39	0.45
1:B:130:PRO:O	1:B:131:ALA:C	2.59	0.45
1:F:18:PHE:CE1	1:F:127:ARG:NH1	2.85	0.45
1:C:33:THR:HG22	1:C:37:LEU:HD12	1.99	0.45
1:D:122:PHE:O	1:D:126:ILE:HG22	2.17	0.44
1:C:66:MET:CE	1:C:89:VAL:HG11	2.47	0.44
1:D:128:THR:O	1:D:133:ARG:NH1	2.49	0.44
1:B:33:THR:HG22	1:B:37:LEU:HD12	2.00	0.44
1:F:62:TRP:CZ2	1:F:66:MET:HE3	2.52	0.44
1:E:7:LYS:HE3	1:F:45:PRO:HB3	1.99	0.44
1:B:47:HIS:O	1:B:48:CYS:HB2	2.16	0.44
1:B:134:PRO:CB	1:B:135:PRO:HD2	2.48	0.44
1:C:23:PHE:CD1	1:C:23:PHE:C	2.96	0.43
1:F:125:TRP:CZ2	1:F:133:ARG:HD2	2.53	0.43
1:D:125:TRP:CE2	1:D:133:ARG:HD2	2.54	0.43
1:A:125:TRP:CZ2	1:A:133:ARG:HD2	2.53	0.43
1:A:16:LEU:HD21	1:A:103:PHE:HB2	2.01	0.42
1:A:6:TYR:CZ	1:A:16:LEU:HD22	2.55	0.42
1:F:134:PRO:CB	1:F:135:PRO:CD	2.98	0.42
1:F:16:LEU:HD21	1:F:103:PHE:HB2	2.00	0.42
1:F:102:TRP:O	1:F:103:PHE:C	2.61	0.42
1:A:74:THR:O	1:A:74:THR:OG1	2.34	0.41
2:D:201:A1AVK:O1	2:D:201:A1AVK:C17	2.68	0.41
1:E:4:ASP:HB3	1:E:7:LYS:HB2	2.01	0.41
1:D:62:TRP:CZ2	1:D:66:MET:HE3	2.55	0.41
1:E:6:TYR:CZ	1:E:16:LEU:HD22	2.54	0.41
1:D:88:TYR:CA	1:D:91:THR:OG1	2.68	0.41
1:E:45:PRO:O	1:F:8:GLU:HA	2.20	0.41
1:B:38:TYR:O	1:B:39:ARG:C	2.63	0.41
1:E:38:TYR:O	1:E:39:ARG:C	2.64	0.41
1:F:34:ALA:O	1:F:38:TYR:HB2	2.21	0.41
1:C:38:TYR:O	1:C:39:ARG:C	2.63	0.41
1:A:122:PHE:O	1:A:126:ILE:HG22	2.21	0.41
1:A:102:TRP:O	1:A:103:PHE:C	2.62	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:ALA:HB2	1:B:1:MET:HE2	2.03	0.40
1:F:33:THR:HG22	1:F:37:LEU:HD12	2.03	0.40
1:F:128:THR:HB	1:F:133:ARG:HB2	2.03	0.40
1:F:136:ASN:HA	1:F:137:ALA:HA	1.86	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	137/167 (82%)	124 (90%)	9 (7%)	4 (3%)	3	4
1	B	131/167 (78%)	123 (94%)	5 (4%)	3 (2%)	5	6
1	C	129/167 (77%)	112 (87%)	13 (10%)	4 (3%)	3	4
1	D	127/167 (76%)	117 (92%)	9 (7%)	1 (1%)	16	24
1	E	128/167 (77%)	117 (91%)	9 (7%)	2 (2%)	7	10
1	F	124/167 (74%)	115 (93%)	7 (6%)	2 (2%)	7	10
All	All	776/1002 (77%)	708 (91%)	52 (7%)	16 (2%)	6	7

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	77	GLU
1	A	134	PRO
1	E	134	PRO
1	F	131	ALA
1	A	48	CYS
1	B	48	CYS
1	C	48	CYS
1	E	131	ALA

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Mol	Chain	Res	Type
1	F	134	PRO
1	C	40[A]	ASP
1	C	40[B]	ASP
1	C	132	ALA
1	D	132	ALA
1	A	135	PRO
1	B	131	ALA
1	B	134	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	118/147 (80%)	107 (91%)	11 (9%)	8	12
1	B	113/147 (77%)	109 (96%)	4 (4%)	32	51
1	C	110/147 (75%)	105 (96%)	5 (4%)	24	40
1	D	110/147 (75%)	103 (94%)	7 (6%)	16	26
1	E	114/147 (78%)	108 (95%)	6 (5%)	20	34
1	F	108/147 (74%)	97 (90%)	11 (10%)	7	10
All	All	673/882 (76%)	629 (94%)	44 (6%)	15	26

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	15	LEU
1	A	16	LEU
1	A	21	SER
1	A	72	VAL
1	A	74	THR
1	A	76	LEU
1	A	86	VAL
1	A	93	VAL
1	A	96	LYS

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Mol	Chain	Res	Type
1	A	133	ARG
1	B	16	LEU
1	B	40	ASP
1	B	72	VAL
1	B	93	VAL
1	C	16	LEU
1	C	27	VAL
1	C	93	VAL
1	C	133	ARG
1	C	136	ASN
1	D	1	MET
1	D	16	LEU
1	D	28	ARG
1	D	45	PRO
1	D	91	THR
1	D	93	VAL
1	D	133	ARG
1	E	1	MET
1	E	16	LEU
1	E	68	LEU
1	E	72	VAL
1	E	93	VAL
1	E	133	ARG
1	F	0	SER
1	F	1	MET
1	F	16	LEU
1	F	21	SER
1	F	27	VAL
1	F	39	ARG
1	F	45	PRO
1	F	72	VAL
1	F	92	ASN
1	F	93	VAL
1	F	133	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	ASN
1	B	92	ASN
1	E	92	ASN
1	F	136	ASN

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 ⓘ

6 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	A1AVK	F	201	-	31,31,31	1.46	3 (9%)	36,44,44	2.28	15 (41%)
2	A1AVK	A	201	-	31,31,31	1.57	4 (12%)	36,44,44	2.38	12 (33%)
2	A1AVK	D	201	-	31,31,31	1.59	4 (12%)	36,44,44	2.22	13 (36%)
2	A1AVK	C	201	-	31,31,31	1.57	5 (16%)	36,44,44	2.32	10 (27%)
2	A1AVK	E	201	-	31,31,31	1.58	5 (16%)	36,44,44	2.24	11 (30%)
2	A1AVK	B	201	-	31,31,31	1.78	5 (16%)	36,44,44	2.34	11 (30%)

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	A1AVK	F	201	-	-	5/18/20/20	0/3/3/3
2	A1AVK	A	201	-	-	5/18/20/20	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	A1AVK	D	201	-	-	5/18/20/20	0/3/3/3
2	A1AVK	C	201	-	-	3/18/20/20	0/3/3/3
2	A1AVK	E	201	-	-	6/18/20/20	0/3/3/3
2	A1AVK	B	201	-	-	5/18/20/20	0/3/3/3

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	201	A1AVK	C14-C9	5.94	1.50	1.40
2	F	201	A1AVK	C14-C9	5.53	1.49	1.40
2	D	201	A1AVK	C14-C9	5.49	1.49	1.40
2	C	201	A1AVK	C14-C9	5.47	1.49	1.40
2	E	201	A1AVK	C14-C9	4.98	1.48	1.40
2	A	201	A1AVK	C14-C9	4.69	1.48	1.40
2	A	201	A1AVK	C7-C6	4.45	1.40	1.35
2	B	201	A1AVK	C7-C6	4.06	1.40	1.35
2	C	201	A1AVK	C7-C6	3.87	1.39	1.35
2	D	201	A1AVK	C7-C6	3.41	1.39	1.35
2	E	201	A1AVK	C7-C6	3.38	1.39	1.35
2	E	201	A1AVK	C20-CL	3.35	1.81	1.73
2	D	201	A1AVK	C20-CL	3.22	1.81	1.73
2	B	201	A1AVK	C20-CL	3.22	1.81	1.73
2	E	201	A1AVK	C13-C14	-3.17	1.35	1.39
2	F	201	A1AVK	C20-CL	3.12	1.81	1.73
2	C	201	A1AVK	C13-C14	-2.65	1.35	1.39
2	C	201	A1AVK	C20-CL	2.56	1.79	1.73
2	A	201	A1AVK	C20-CL	2.48	1.79	1.73
2	F	201	A1AVK	C7-C6	2.31	1.38	1.35
2	C	201	A1AVK	C11-C10	-2.28	1.35	1.38
2	E	201	A1AVK	O1-C15	-2.27	1.19	1.23
2	B	201	A1AVK	C13-C14	-2.25	1.36	1.39
2	B	201	A1AVK	C15-N2	-2.22	1.33	1.37
2	D	201	A1AVK	C13-C14	-2.11	1.36	1.39
2	A	201	A1AVK	C12-C13	2.05	1.42	1.38

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	201	A1AVK	N2-C15-N	8.42	125.06	115.54
2	A	201	A1AVK	C9-C8-N1	8.08	121.12	114.33
2	C	201	A1AVK	C5-C4-N	-7.61	103.79	112.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	201	A1AVK	N2-C15-N	6.99	123.44	115.54
2	D	201	A1AVK	C5-C4-N	-6.23	105.35	112.38
2	B	201	A1AVK	C5-C4-N	-6.12	105.47	112.38
2	D	201	A1AVK	N2-C15-N	5.70	121.98	115.54
2	F	201	A1AVK	C10-C9-C14	5.28	125.19	119.26
2	D	201	A1AVK	C9-C8-N1	5.20	118.70	114.33
2	E	201	A1AVK	C9-C8-N1	5.13	118.64	114.33
2	A	201	A1AVK	C5-C4-N	-4.54	107.26	112.38
2	C	201	A1AVK	C10-C9-C14	4.53	124.34	119.26
2	F	201	A1AVK	O-C8-N1	4.45	125.86	120.44
2	F	201	A1AVK	C11-C12-C13	4.41	125.68	120.24
2	B	201	A1AVK	C9-C8-N1	4.29	117.93	114.33
2	C	201	A1AVK	C11-C12-C13	4.20	125.42	120.24
2	A	201	A1AVK	N2-C15-N	4.18	120.27	115.54
2	E	201	A1AVK	C10-C9-C14	4.03	123.78	119.26
2	B	201	A1AVK	C11-C12-C13	4.01	125.19	120.24
2	C	201	A1AVK	O-C8-N1	3.71	124.95	120.44
2	F	201	A1AVK	O-C8-C9	-3.64	116.77	123.27
2	F	201	A1AVK	C9-C8-N1	3.62	117.37	114.33
2	C	201	A1AVK	C3-N-C4	-3.46	113.32	117.00
2	B	201	A1AVK	C13-C14-C9	-3.44	115.40	119.26
2	A	201	A1AVK	C14-C9-C8	-3.34	116.05	119.56
2	F	201	A1AVK	C5-C4-N	-3.24	108.73	112.38
2	A	201	A1AVK	C13-C14-C9	-3.23	115.64	119.26
2	F	201	A1AVK	N2-C15-N	3.13	119.07	115.54
2	F	201	A1AVK	C3-N-C4	-3.08	113.73	117.00
2	D	201	A1AVK	C11-C12-C13	3.07	124.02	120.24
2	C	201	A1AVK	C13-C14-C9	-2.98	115.92	119.26
2	C	201	A1AVK	N2-C15-N	2.97	118.90	115.54
2	E	201	A1AVK	C17-C16-C21	-2.92	116.12	119.66
2	B	201	A1AVK	C17-C16-C21	-2.90	116.15	119.66
2	D	201	A1AVK	C21-C20-CL	2.88	123.16	118.45
2	A	201	A1AVK	C16-N2-C15	-2.86	120.26	126.04
2	A	201	A1AVK	C19-C20-CL	-2.83	116.22	119.79
2	A	201	A1AVK	C11-C12-C13	2.80	123.69	120.24
2	E	201	A1AVK	C11-C12-C13	2.77	123.66	120.24
2	E	201	A1AVK	O-C8-C9	-2.77	118.33	123.27
2	D	201	A1AVK	C19-C20-CL	-2.74	116.33	119.79
2	F	201	A1AVK	C13-C14-C9	-2.57	116.38	119.26
2	C	201	A1AVK	C17-C16-C21	-2.56	116.56	119.66
2	E	201	A1AVK	O1-C15-N2	-2.55	117.87	123.46
2	D	201	A1AVK	C16-N2-C15	-2.50	120.98	126.04

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	201	A1AVK	C12-C11-C10	-2.50	117.16	120.24
2	F	201	A1AVK	C18-C19-C20	-2.44	119.11	121.36
2	B	201	A1AVK	C10-C9-C14	2.42	121.98	119.26
2	F	201	A1AVK	C21-C20-CL	2.37	122.33	118.45
2	D	201	A1AVK	C14-C9-C8	-2.37	117.08	119.56
2	D	201	A1AVK	C2-C1-C3	-2.36	101.43	110.99
2	B	201	A1AVK	C21-C20-CL	2.35	122.30	118.45
2	D	201	A1AVK	C18-C17-C16	2.34	122.99	120.30
2	B	201	A1AVK	C14-C9-C8	-2.33	117.12	119.56
2	F	201	A1AVK	C14-C9-C8	-2.31	117.14	119.56
2	E	201	A1AVK	C14-C9-C8	-2.30	117.14	119.56
2	D	201	A1AVK	C10-C9-C14	2.28	121.82	119.26
2	E	201	A1AVK	C13-C14-C9	-2.27	116.71	119.26
2	B	201	A1AVK	C16-N2-C15	-2.25	121.48	126.04
2	A	201	A1AVK	C21-C20-CL	2.25	122.14	118.45
2	C	201	A1AVK	O-C8-C9	-2.25	119.25	123.27
2	C	201	A1AVK	C18-C17-C16	2.24	122.88	120.30
2	D	201	A1AVK	O-C8-C9	-2.24	119.27	123.27
2	F	201	A1AVK	C17-C16-C21	-2.23	116.96	119.66
2	B	201	A1AVK	C12-C11-C10	-2.17	117.56	120.24
2	D	201	A1AVK	C-C1-C3	2.15	119.72	110.99
2	A	201	A1AVK	F-C19-C20	-2.14	116.37	118.96
2	E	201	A1AVK	O-C8-N1	2.13	123.03	120.44
2	F	201	A1AVK	C19-C20-CL	-2.09	117.15	119.79
2	F	201	A1AVK	C11-C10-C9	-2.07	116.05	119.80
2	E	201	A1AVK	C16-C21-C20	2.04	122.01	119.54
2	A	201	A1AVK	O-C8-C9	-2.03	119.64	123.27

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	201	A1AVK	O1-C15-N-C4
2	E	201	A1AVK	O1-C15-N-C4
2	F	201	A1AVK	C-C1-C3-N
2	A	201	A1AVK	N-C15-N2-C16
2	B	201	A1AVK	N-C15-N2-C16
2	D	201	A1AVK	N-C15-N2-C16
2	F	201	A1AVK	N-C15-N2-C16
2	A	201	A1AVK	C2-C1-C3-N
2	C	201	A1AVK	N-C15-N2-C16
2	A	201	A1AVK	O1-C15-N-C4

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Mol	Chain	Res	Type	Atoms
2	B	201	A1AVK	O1-C15-N-C4
2	E	201	A1AVK	N-C15-N2-C16
2	A	201	A1AVK	C1-C3-N-C4
2	C	201	A1AVK	C1-C3-N-C15
2	F	201	A1AVK	C2-C1-C3-N
2	A	201	A1AVK	O1-C15-N2-C16
2	B	201	A1AVK	O1-C15-N-C3
2	E	201	A1AVK	O1-C15-N-C3
2	D	201	A1AVK	O1-C15-N-C3
2	E	201	A1AVK	C2-C1-C3-N
2	B	201	A1AVK	C1-C3-N-C4
2	C	201	A1AVK	C1-C3-N-C4
2	D	201	A1AVK	C1-C3-N-C4
2	E	201	A1AVK	C1-C3-N-C4
2	F	201	A1AVK	C1-C3-N-C4
2	B	201	A1AVK	C1-C3-N-C15
2	E	201	A1AVK	C1-C3-N-C15
2	F	201	A1AVK	C1-C3-N-C15
2	D	201	A1AVK	O1-C15-N2-C16

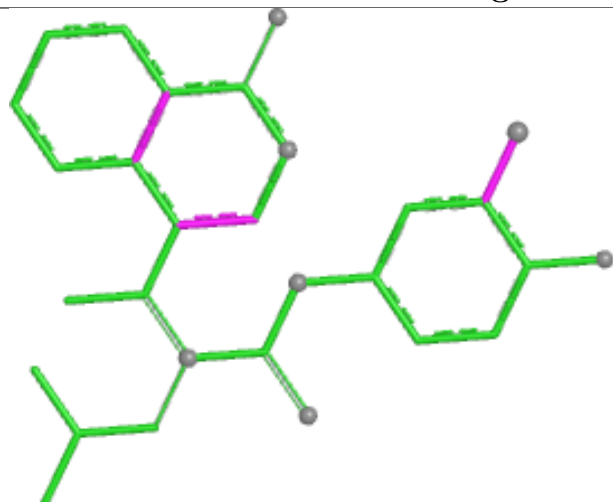
There are no ring outliers.

2 monomers are involved in 2 short contacts:

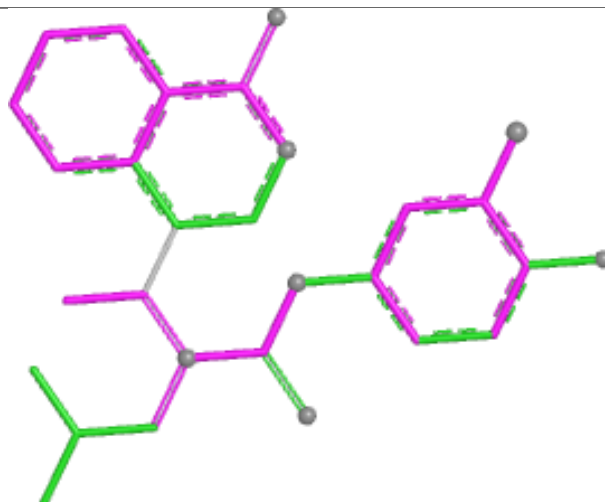
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	201	A1AVK	1	0
2	B	201	A1AVK	1	0

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.

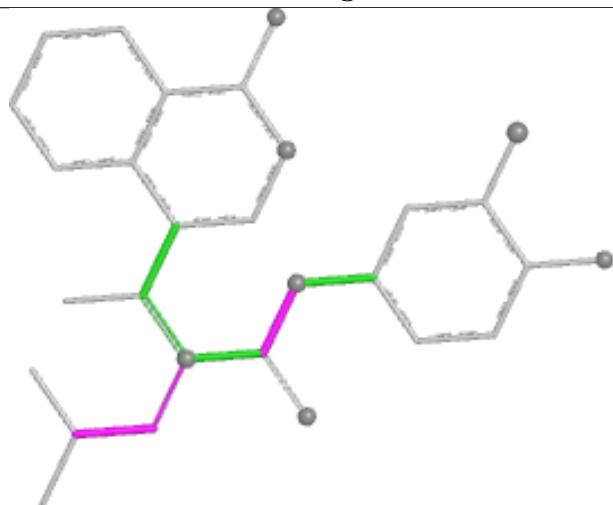
Ligand A1AVK F 201



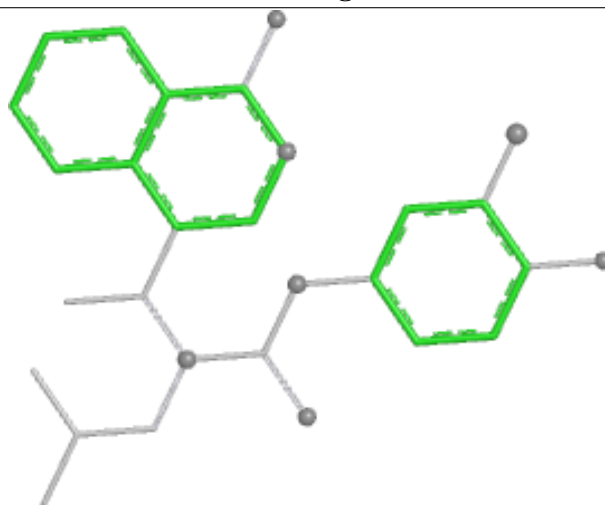
Bond lengths



Bond angles

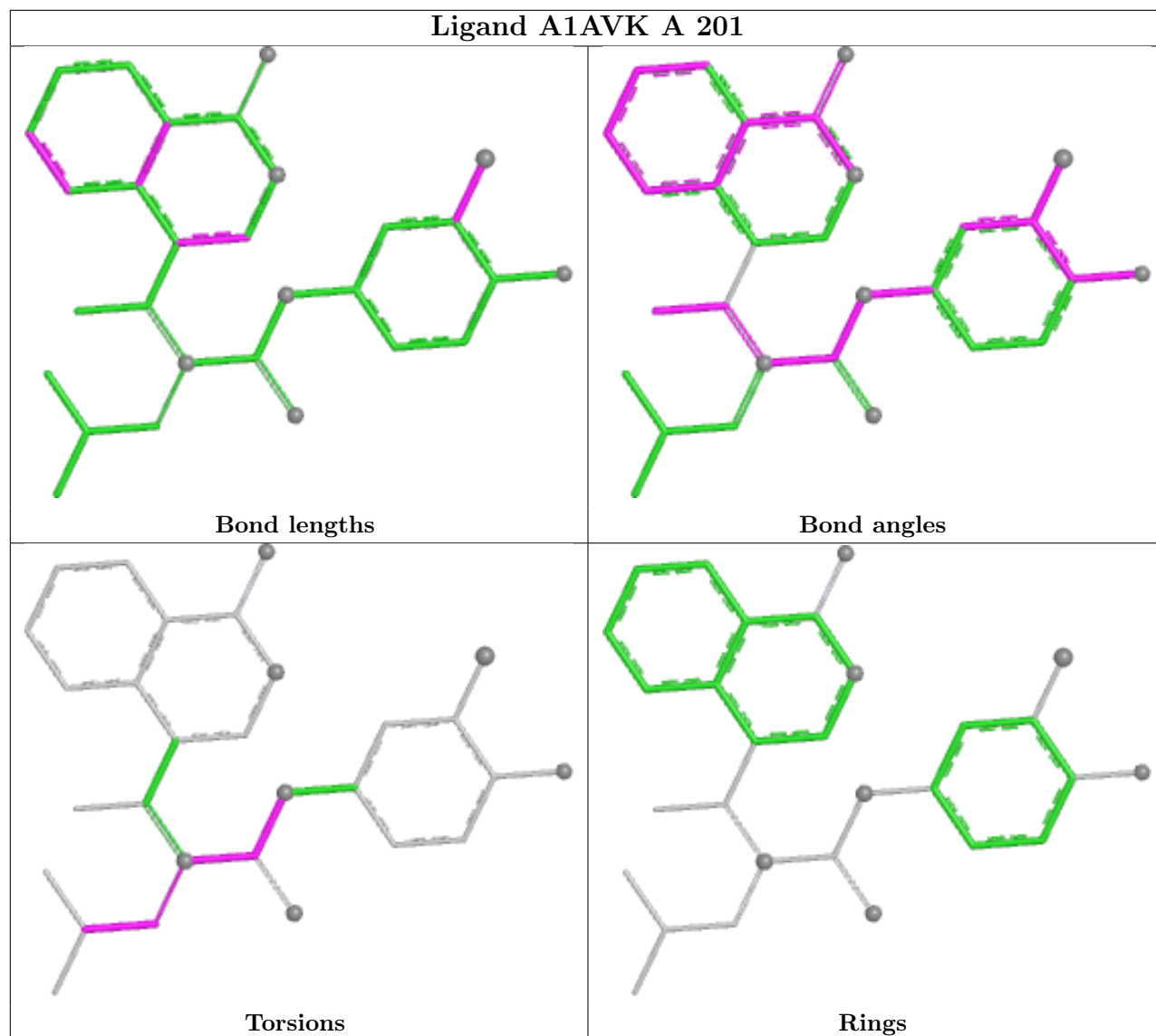


Torsions

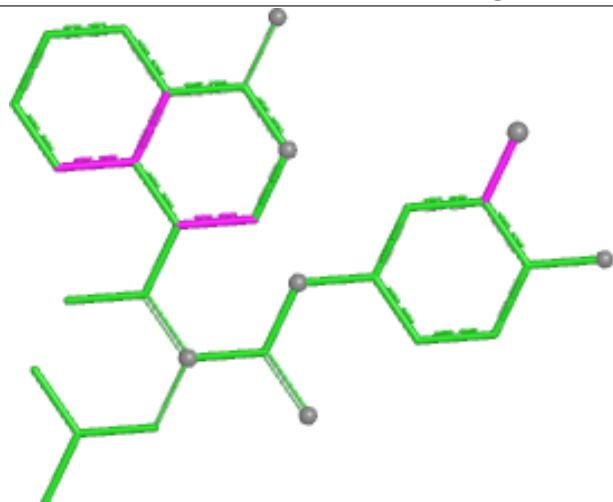


Rings

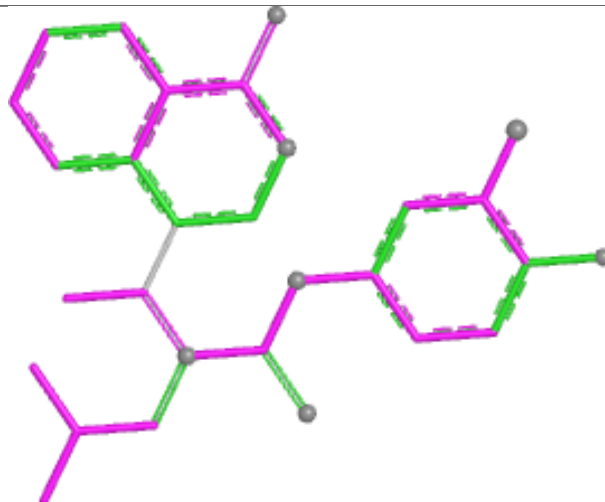
Ligand A1AVK A 201



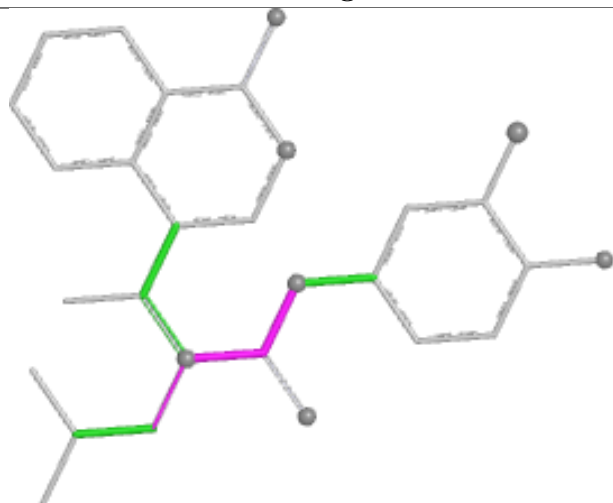
Ligand A1AVK D 201



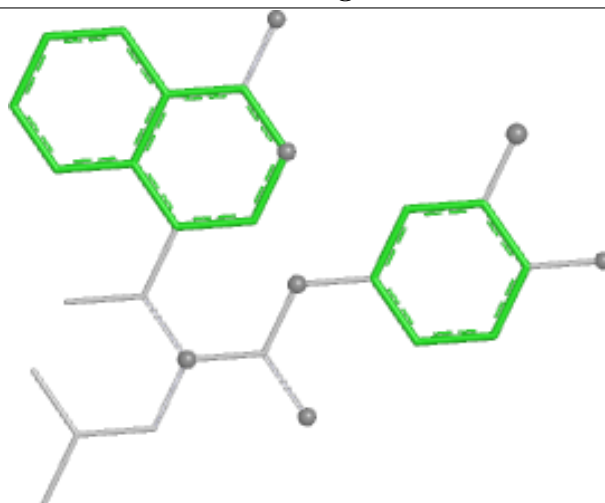
Bond lengths



Bond angles

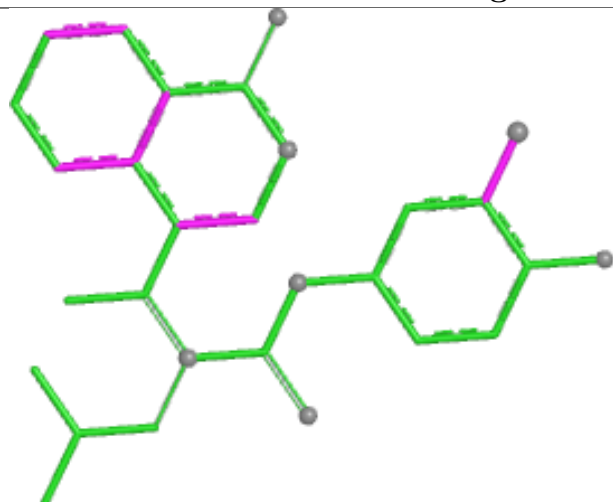


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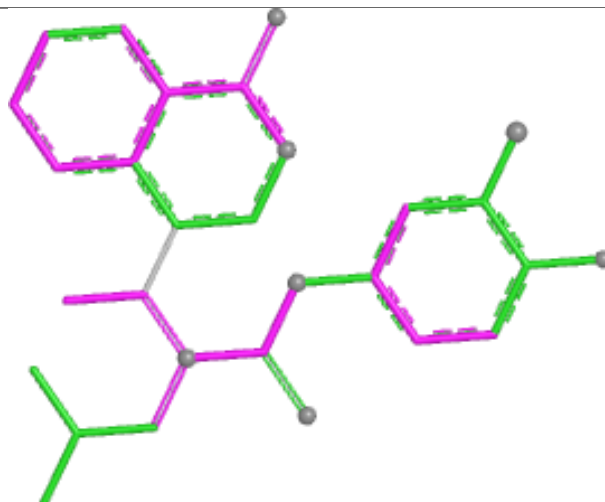


Rings

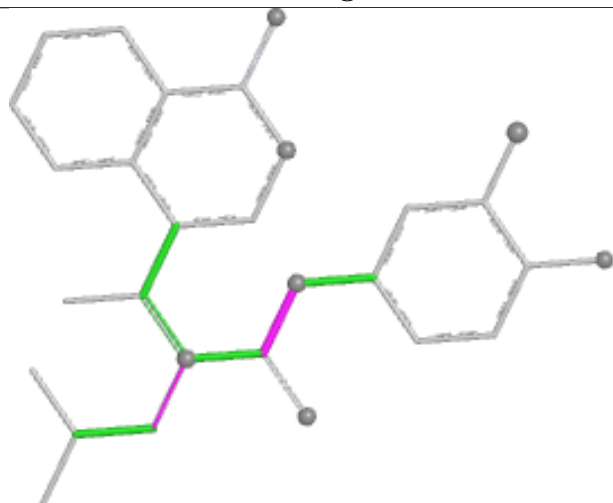
Ligand A1AVK C 201



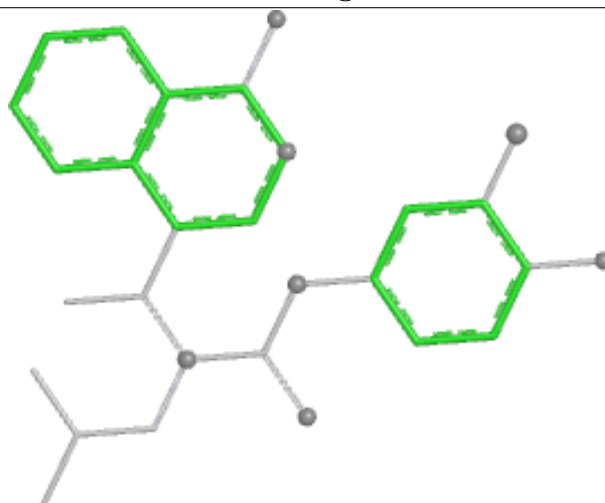
Bond lengths



Bond angles

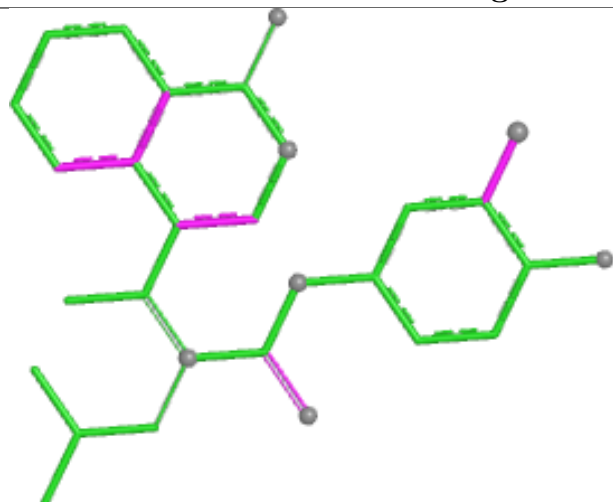


Torsions

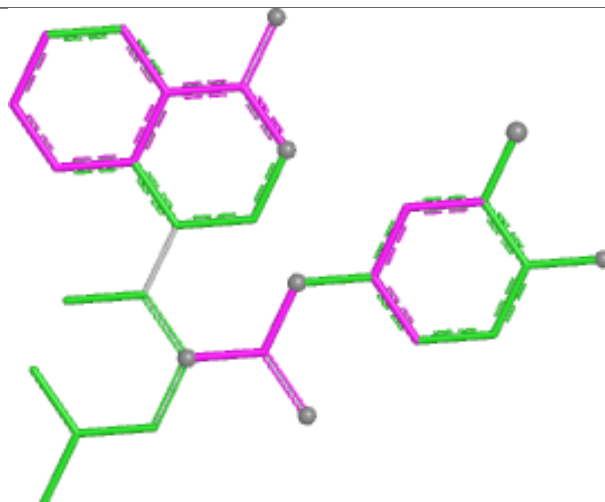


Rings

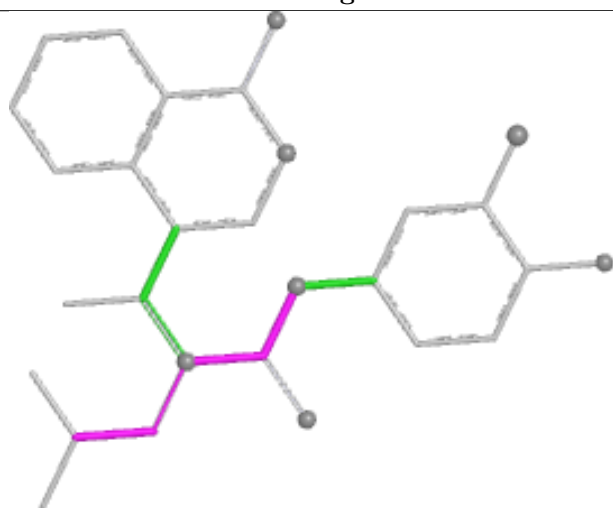
Ligand A1AVK E 201



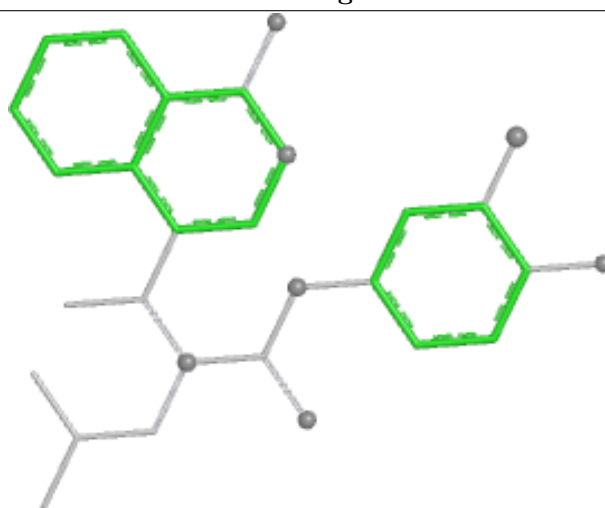
Bond lengths



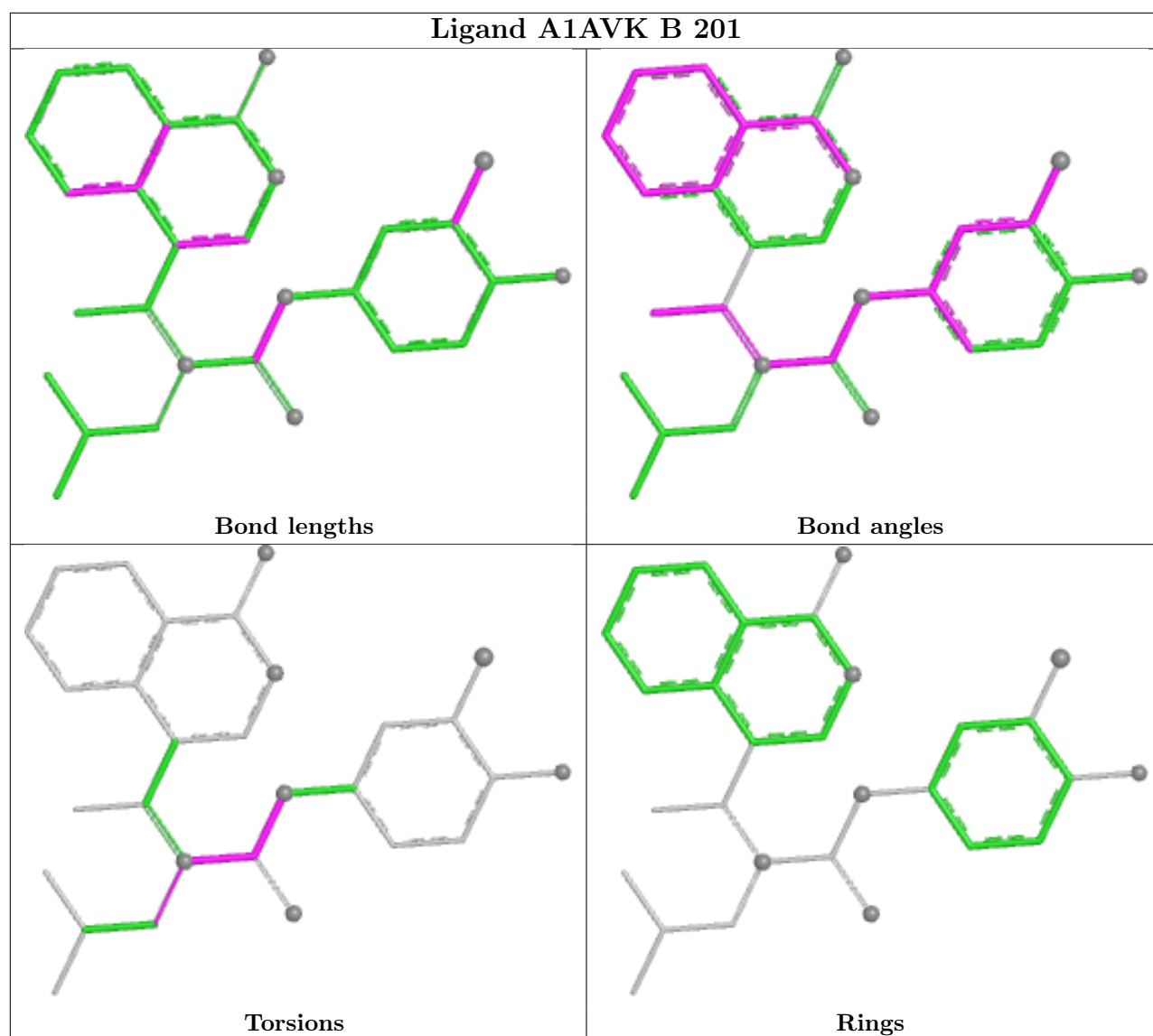
Bond angles



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, 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	138/167 (82%)	0.28	8 (5%) 29 24	35, 64, 136, 158	1 (0%)
1	B	135/167 (80%)	0.32	12 (8%) 15 12	46, 64, 158, 199	0
1	C	132/167 (79%)	0.30	10 (7%) 20 16	30, 63, 135, 154	1 (0%)
1	D	131/167 (78%)	0.41	14 (10%) 11 9	46, 66, 143, 185	0
1	E	130/167 (77%)	0.29	8 (6%) 26 22	29, 64, 117, 138	2 (1%)
1	F	126/167 (75%)	0.32	10 (7%) 18 15	33, 62, 122, 138	2 (1%)
All	All	792/1002 (79%)	0.32	62 (7%) 19 15	29, 64, 134, 199	6 (0%)

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	136	ASN	6.9
1	D	76	LEU	6.0
1	C	137	ALA	6.0
1	D	86	VAL	5.5
1	F	89	VAL	4.5
1	B	82	ARG	4.3
1	D	83	ASP	4.3
1	D	85	VAL	4.3
1	E	73	GLY	4.1
1	B	85	VAL	4.0
1	C	77	GLU	3.9
1	D	137	ALA	3.9
1	A	86	VAL	3.6
1	E	84	LEU	3.6
1	B	84	LEU	3.5
1	A	80	ALA	3.4
1	A	85	VAL	3.3
1	B	86	VAL	3.3
1	F	137	ALA	3.3

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Mol	Chain	Res	Type	RSRZ
1	E	82	ARG	3.2
1	D	72	VAL	3.2
1	C	84	LEU	3.1
1	E	136	ASN	3.0
1	E	135	PRO	3.0
1	F	135	PRO	3.0
1	C	0	SER	2.9
1	F	91	THR	2.9
1	C	74	THR	2.9
1	B	136	ASN	2.8
1	F	14[A]	GLU	2.8
1	D	90	ASN	2.7
1	B	83	ASP	2.7
1	A	84	LEU	2.7
1	A	120	VAL	2.7
1	F	72	VAL	2.7
1	C	135	PRO	2.7
1	D	135	PRO	2.7
1	F	131	ALA	2.6
1	B	73	GLY	2.6
1	F	90	ASN	2.6
1	C	86	VAL	2.5
1	D	84	LEU	2.4
1	B	75	ASN	2.4
1	A	135	PRO	2.4
1	E	137	ALA	2.4
1	E	119	LEU	2.3
1	F	76	LEU	2.3
1	B	135	PRO	2.3
1	D	14	GLU	2.2
1	C	40[A]	ASP	2.2
1	D	2	ASP	2.2
1	A	118	TYR	2.2
1	C	136	ASN	2.2
1	C	85	VAL	2.2
1	B	80	ALA	2.1
1	A	136	ASN	2.1
1	B	9	PHE	2.1
1	D	117	GLU	2.1
1	D	74	THR	2.1
1	F	71	TRP	2.0
1	B	120	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
1	E	120	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](#)

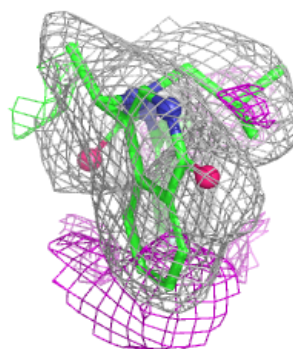
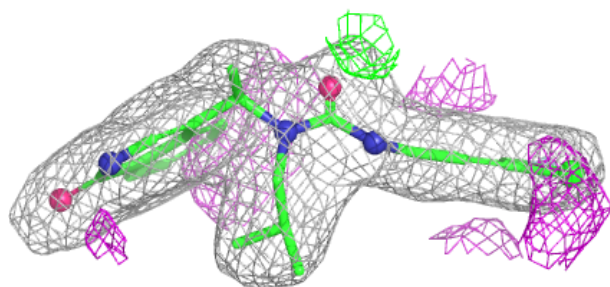
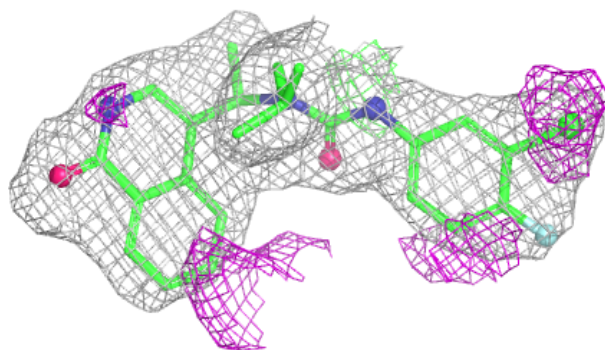
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	A1AVK	F	201	29/29	0.92	0.10	46,51,63,73	0
2	A1AVK	E	201	29/29	0.93	0.10	48,57,69,81	0
2	A1AVK	D	201	29/29	0.94	0.10	44,55,70,72	0
2	A1AVK	A	201	29/29	0.94	0.09	42,55,69,77	0
2	A1AVK	C	201	29/29	0.94	0.10	46,57,68,75	0
2	A1AVK	B	201	29/29	0.95	0.09	46,58,70,72	0

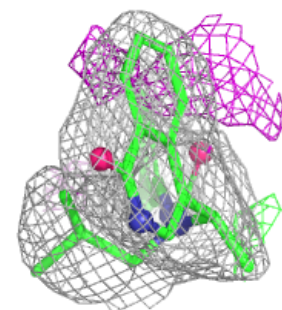
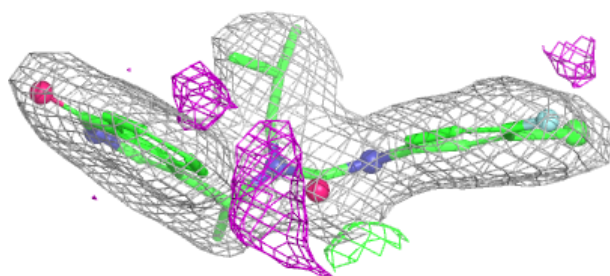
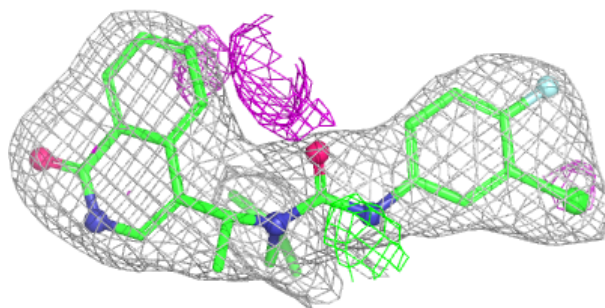
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around A1AVK F 201:

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

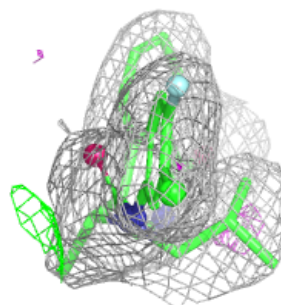
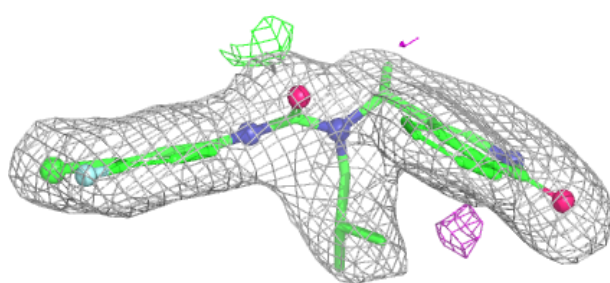
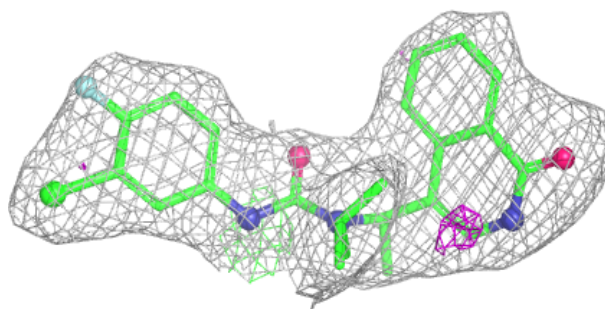
**Electron density around A1AVK E 201:**

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

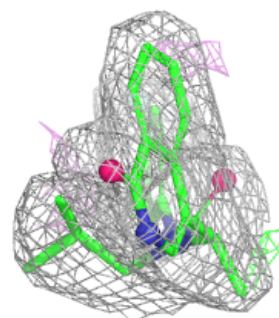
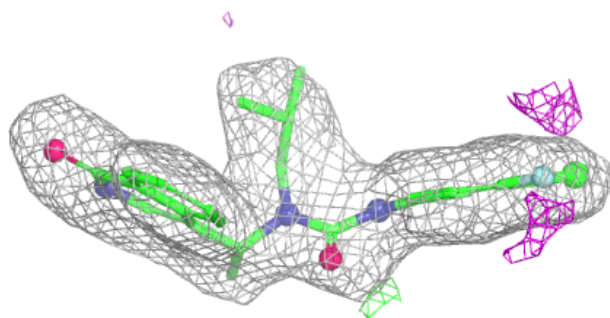
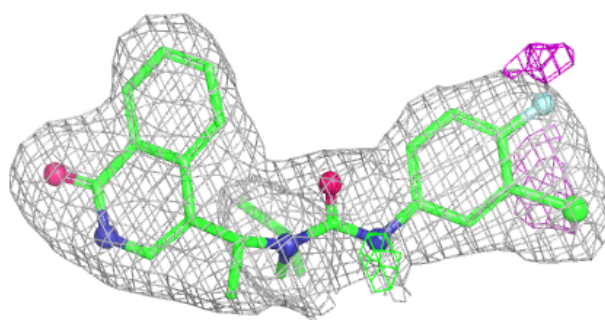


Electron density around A1AVK D 201:

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

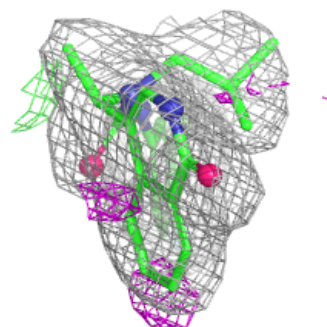
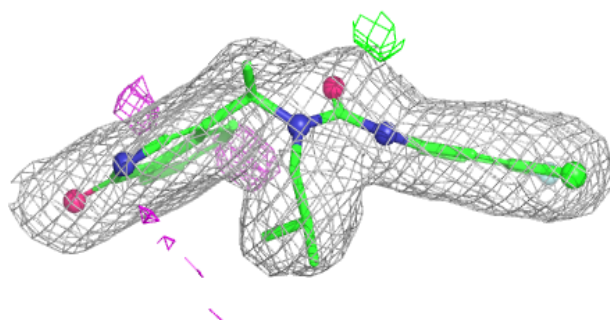
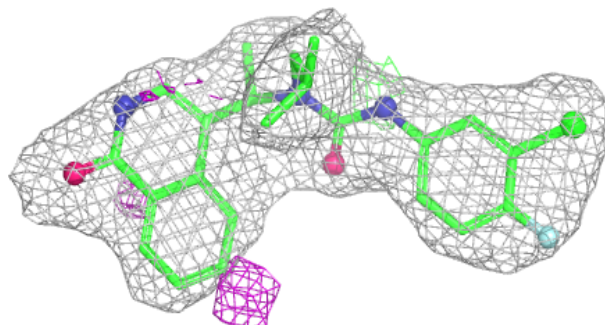
**Electron density around A1AVK A 201:**

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

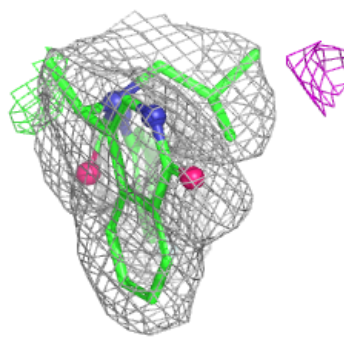
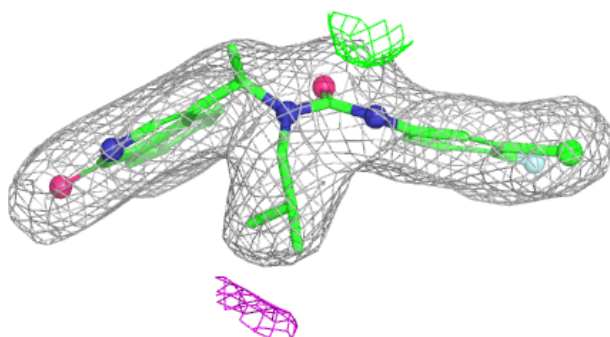
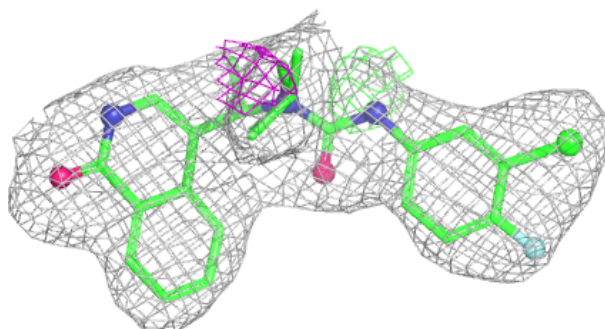


Electron density around A1AVK C 201:

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

**Electron density around A1AVK B 201:**

$2mF_o-DF_c$ (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.