



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 3, 2025 – 09:15 am GMT

PDB ID : 9GUJ  
Title : Crystal structure of transcription factor NtcA from *Synechococcus elongatus* in complex with its transcriptional co- activator PipX and its target DNA (Crystal II)  
Authors : Llacer, J.L.; Forcada-Nadal, A.; Rubio, V.  
Deposited on : 2024-09-19  
Resolution : 4.30 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.

The types of validation reports are described at

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

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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	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.41

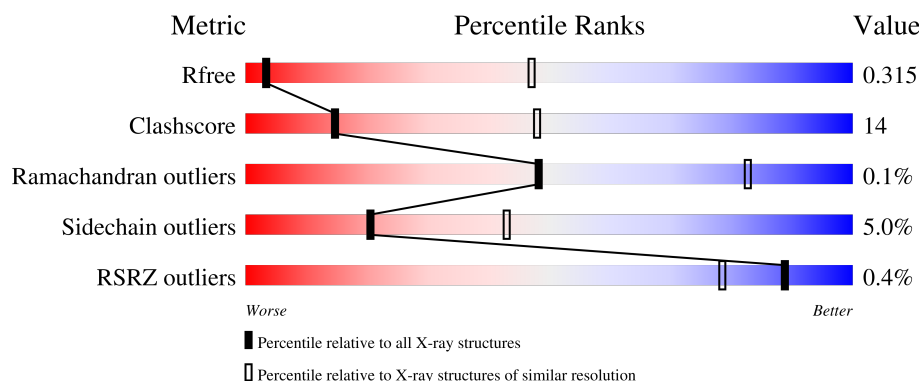
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*





The reported resolution of this entry is 4.30 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))
$R_{free}$	164625	1028 (4.72-3.86)
Clashscore	180529	1030 (4.70-3.90)
Ramachandran outliers	177936	1014 (4.76-3.84)
Sidechain outliers	177891	1022 (4.76-3.82)
RSRZ outliers	164620	1026 (4.72-3.86)

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  $\geq 3$ , 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	222	
1	B	222	
1	D	222	
1	F	222	

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Mol	Chain	Length	Quality of chain
2	C	30	
2	E	30	
2	G	30	
2	I	30	
3	H	89	
3	J	89	
3	K	89	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	AKG	F	301	-	X	-	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11268 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 Global nitrogen regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	213	Total	C	N	O	S	0	0	0
			1665	1070	288	300	7			
1	F	216	Total	C	N	O	S	0	0	0
			1689	1086	290	306	7			
1	A	213	Total	C	N	O	S	0	0	0
			1665	1070	288	300	7			
1	B	215	Total	C	N	O	S	0	0	0
			1685	1084	289	305	7			

- Molecule 2 is a DNA chain called DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	30	Total	C	N	O	P	0	0	0
			615	297	108	180	30			
2	I	30	Total	C	N	O	P	0	0	0
			615	297	108	180	30			
2	C	28	Total	C	N	O	P	0	0	0
			574	277	101	168	28			
2	E	30	Total	C	N	O	P	0	0	0
			615	297	108	180	30			

- Molecule 3 is a protein called PipX.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	J	84	Total	C	N	O	S	0	0	0
			701	451	125	123	2			
3	H	85	Total	C	N	O	S	0	0	0
			703	454	123	124	2			
3	K	84	Total	C	N	O	S	0	0	0
			701	451	125	123	2			

- Molecule 4 is 2-OXOGLUTARIC ACID (three-letter code: AKG) (formula: C<sub>5</sub>H<sub>6</sub>O<sub>5</sub>) (labeled as "Ligand of Interest" by depositor).

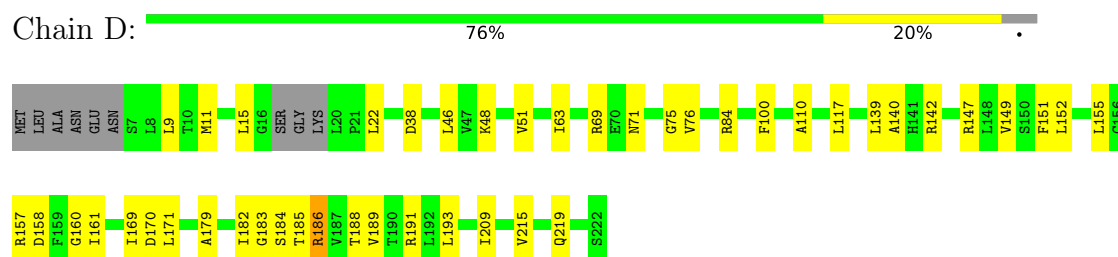


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	C	O	0	0
			10	5	5		
4	F	1	Total	C	O	0	0
			10	5	5		
4	A	1	Total	C	O	0	0
			10	5	5		
4	B	1	Total	C	O	0	0
			10	5	5		

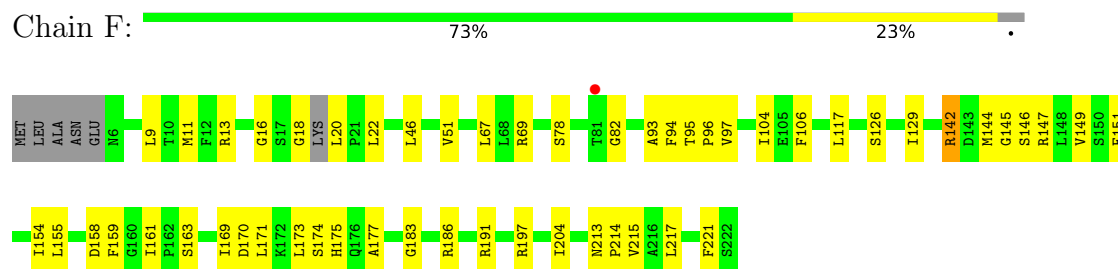
### 3 Residue-property plots [i](#)

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.

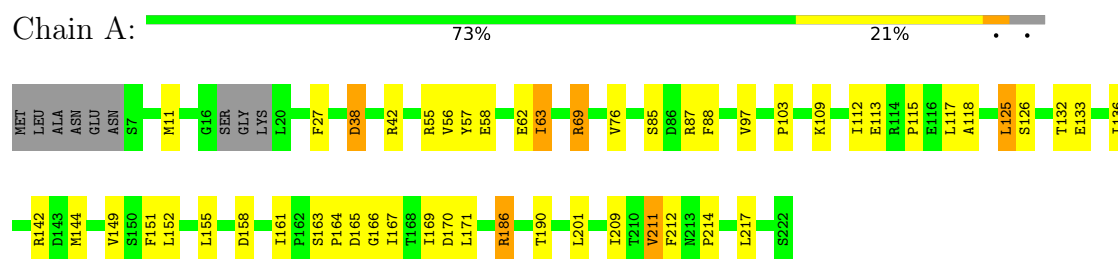
- Molecule 1: Global nitrogen regulator



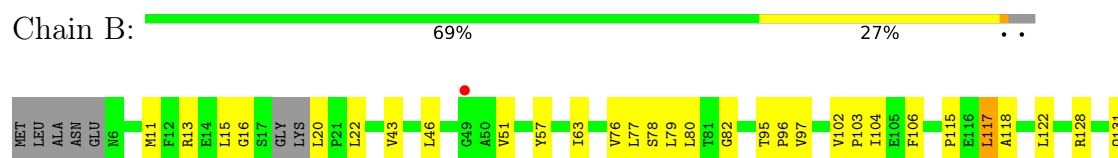
- Molecule 1: Global nitrogen regulator

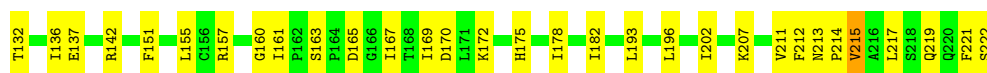


- Molecule 1: Global nitrogen regulator

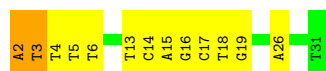


- Molecule 1: Global nitrogen regulator

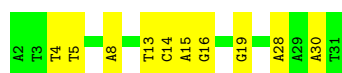




- Molecule 2: DNA



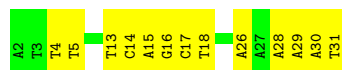
- Molecule 2: DNA



- Molecule 2: DNA



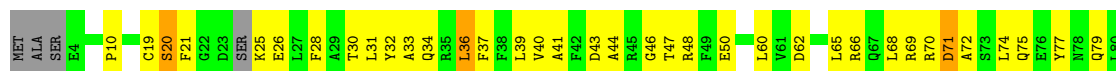
- Molecule 2: DNA



- Molecule 3: PipX



- Molecule 3: PipX



- Molecule 3: PipX





## 4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	165.71Å 165.71Å 177.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	82.86 – 4.30 82.86 – 4.30	Depositor EDS
% Data completeness (in resolution range)	98.5 (82.86-4.30) 98.4 (82.86-4.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.57 (at 4.30Å)	Xtriage
Refinement program	REFMAC 5.8.0352	Depositor
R, $R_{free}$	0.284 , 0.316 0.276 , 0.315	Depositor DCC
$R_{free}$ test set	817 reflections (5.21%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	178.5	Xtriage
Anisotropy	0.088	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 364.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.389 for -h,k,-l	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	11268	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	96.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.24% 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: AKG

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	0.39	0/1692	0.70	0/2286
1	B	0.44	0/1712	0.72	0/2311
1	D	0.40	0/1692	0.69	0/2286
1	F	0.43	0/1716	0.68	0/2316
2	C	0.68	0/643	0.97	0/990
2	E	0.72	0/689	1.05	2/1061 (0.2%)
2	G	0.70	0/689	1.03	2/1061 (0.2%)
2	I	0.72	0/689	1.03	1/1061 (0.1%)
3	H	0.46	0/717	0.72	0/965
3	J	0.49	0/715	0.79	1/961 (0.1%)
3	K	0.44	0/715	0.71	0/961
All	All	0.51	0/11669	0.80	6/16259 (0.0%)

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
3	H	0	2
3	J	0	2
All	All	0	4

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	3	DT	P-O3'-C3'	-8.47	109.53	119.70
2	G	2	DA	P-O3'-C3'	-7.99	110.11	119.70
2	E	28	DA	P-O3'-C3'	-7.93	110.18	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	29	DA	P-O3'-C3'	-6.57	111.81	119.70
3	J	74	LEU	O-C-N	5.61	131.68	122.70
2	I	30	DA	P-O3'-C3'	-5.17	113.49	119.70

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	H	20	SER	Mainchain
3	H	70	ARG	Sidechain
3	J	48	ARG	Sidechain
3	J	69	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	1665	0	1719	52	0
1	B	1685	0	1750	51	0
1	D	1665	0	1719	41	0
1	F	1689	0	1753	53	0
2	C	574	0	320	13	0
2	E	615	0	343	12	0
2	G	615	0	343	14	0
2	I	615	0	343	9	0
3	H	703	0	681	35	0
3	J	701	0	681	46	0
3	K	701	0	681	22	0
4	A	10	0	4	1	0
4	B	10	0	4	2	0
4	D	10	0	4	3	0
4	F	10	0	4	0	0
All	All	11268	0	10349	296	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 14.

All (296) 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)
3:J:85:LYS:NZ	1:A:58:GLU:HG3	1.63	1.13
3:J:75:GLN:HA	3:J:78:ASN:HB2	1.31	1.10
3:J:75:GLN:HA	3:J:78:ASN:CB	1.85	1.05
3:H:68:LEU:HA	3:H:72:ALA:HB3	1.08	1.04
3:H:68:LEU:CA	3:H:72:ALA:HB3	1.91	1.01
1:A:118:ALA:HB1	1:B:118:ALA:HB1	1.43	0.97
3:J:85:LYS:NZ	1:A:58:GLU:CG	2.28	0.96
3:K:68:LEU:HA	3:K:72:ALA:HB3	1.49	0.94
3:H:68:LEU:HA	3:H:72:ALA:CB	1.97	0.94
3:J:85:LYS:HZ2	1:A:58:GLU:HG3	1.21	0.92
1:B:82:GLY:HA2	1:B:104:ILE:CD1	2.00	0.92
1:F:11:MET:HB3	1:F:117:LEU:HD13	1.50	0.91
3:H:43:ASP:HB3	3:H:46:GLY:HA3	1.51	0.89
1:B:78:SER:O	1:B:104:ILE:HD12	1.74	0.88
1:D:157:ARG:HD3	3:H:32:TYR:HB2	1.57	0.86
1:F:95:THR:O	1:F:97:VAL:HG23	1.73	0.86
3:J:85:LYS:HZ2	1:A:58:GLU:CG	1.87	0.85
3:H:68:LEU:HD23	3:H:72:ALA:HB1	1.58	0.85
1:A:190:THR:HG23	2:C:10:DG:OP2	1.77	0.85
1:A:125:LEU:HG	1:B:122:LEU:HD23	1.60	0.84
1:D:157:ARG:HG3	3:H:31:LEU:HB3	1.60	0.84
1:F:171:LEU:HD11	1:F:173:LEU:HD21	1.60	0.83
1:F:78:SER:HB2	1:F:104:ILE:HG21	1.62	0.82
3:J:68:LEU:HA	3:J:72:ALA:HB3	1.61	0.79
3:J:85:LYS:HD2	1:A:58:GLU:HG2	1.66	0.77
3:J:85:LYS:HZ1	1:A:58:GLU:HG3	1.47	0.77
1:D:157:ARG:HA	3:H:31:LEU:HD23	1.69	0.74
3:J:75:GLN:HA	3:J:78:ASN:CG	2.08	0.73
3:H:77:TYR:O	3:H:81:GLN:HG3	1.89	0.73
1:A:125:LEU:HG	1:B:122:LEU:CD2	2.20	0.72
3:J:74:LEU:O	3:J:78:ASN:N	2.23	0.72
1:F:191:ARG:NH2	2:G:19:DG:N7	2.38	0.71
3:H:68:LEU:HD23	3:H:72:ALA:CB	2.20	0.71
3:J:75:GLN:O	3:J:79:GLN:N	2.20	0.71
3:K:68:LEU:HD23	3:K:72:ALA:HB1	1.73	0.71
1:A:109:LYS:O	1:A:113:GLU:HG3	1.90	0.71
3:J:85:LYS:CD	1:A:58:GLU:HG2	2.22	0.70
1:F:94:PHE:HE2	1:F:171:LEU:HD13	1.57	0.69
1:F:78:SER:HB2	1:F:104:ILE:CG2	2.22	0.68
1:D:11:MET:CB	1:D:117:LEU:HD13	2.24	0.68
1:F:159:PHE:CB	1:F:171:LEU:HD23	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:77:TYR:CE2	3:H:81:GLN:NE2	2.62	0.68
1:B:57:TYR:HE2	1:B:63:ILE:HD12	1.59	0.68
1:D:76:VAL:HG11	1:F:129:ILE:HD13	1.75	0.67
3:J:74:LEU:O	3:J:77:TYR:HB3	1.94	0.67
1:A:118:ALA:CB	1:B:118:ALA:HB1	2.22	0.67
1:F:159:PHE:CG	1:F:171:LEU:HD23	2.29	0.67
1:B:82:GLY:HA2	1:B:104:ILE:HD13	1.74	0.67
1:B:219:GLN:NE2	3:K:86:GLN:O	2.28	0.66
1:F:169:ILE:HG22	1:F:171:LEU:HG	1.77	0.66
1:B:221:PHE:O	1:B:222:SER:C	2.31	0.66
1:B:15:LEU:O	1:B:22:LEU:HB2	1.96	0.65
1:D:140:ALA:O	1:F:183:GLY:HA2	1.95	0.65
3:K:68:LEU:CA	3:K:72:ALA:HB3	2.26	0.65
1:F:174:SER:OG	1:F:177:ALA:HB2	1.97	0.65
1:A:169:ILE:HG22	1:A:171:LEU:HD23	1.77	0.65
3:J:75:GLN:CA	3:J:78:ASN:HB2	2.19	0.65
1:F:69:ARG:NH1	1:F:158:ASP:O	2.29	0.64
2:E:4:DT:H2''	2:E:5:DT:O5'	1.97	0.64
1:D:11:MET:HB3	1:D:117:LEU:HD13	1.80	0.64
1:A:132:THR:HG22	1:B:132:THR:HG22	1.79	0.64
3:K:68:LEU:HD23	3:K:72:ALA:CB	2.28	0.64
2:G:2:DA:H2'	2:G:3:DT:C6	2.32	0.64
1:A:112:ILE:HA	1:B:115:PRO:HB2	1.79	0.64
1:A:11:MET:HB3	1:A:117:LEU:HD13	1.79	0.64
3:J:85:LYS:HZ1	1:A:58:GLU:CG	2.08	0.63
1:F:197:ARG:HH21	1:F:204:ILE:HD12	1.64	0.63
1:F:171:LEU:CD1	1:F:173:LEU:HD21	2.27	0.63
1:D:75:GLY:N	4:D:301:AKG:O1	2.32	0.63
1:B:82:GLY:HA2	1:B:104:ILE:HD12	1.80	0.62
3:J:68:LEU:CA	3:J:72:ALA:HB3	2.30	0.62
1:B:131:GLN:HG2	1:B:157:ARG:HH22	1.65	0.62
1:A:69:ARG:NH1	1:A:158:ASP:O	2.32	0.61
1:A:69:ARG:HH11	1:A:69:ARG:HB2	1.66	0.60
1:F:94:PHE:CE2	1:F:171:LEU:HD13	2.37	0.60
1:B:76:VAL:HG23	4:B:501:AKG:O1	2.00	0.60
1:D:76:VAL:HG23	4:D:301:AKG:O1	2.00	0.60
1:B:212:PHE:O	1:B:213:ASN:C	2.40	0.60
1:F:174:SER:OG	1:F:177:ALA:CB	2.49	0.60
1:F:82:GLY:HA2	1:F:104:ILE:CD1	2.32	0.59
1:B:214:PRO:O	1:B:215:VAL:C	2.41	0.59
3:J:68:LEU:O	3:J:72:ALA:N	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:76:VAL:HG11	1:F:129:ILE:CD1	2.33	0.59
1:A:169:ILE:HD12	1:A:209:ILE:HB	1.85	0.58
2:C:5:DT:H1'	2:C:6:DT:H5''	1.85	0.58
3:J:85:LYS:CE	1:A:58:GLU:HG2	2.34	0.58
3:H:71:ASP:OD2	3:H:71:ASP:N	2.33	0.58
1:A:118:ALA:HB1	1:B:118:ALA:CB	2.27	0.58
1:F:16:GLY:HA2	1:F:22:LEU:HB2	1.85	0.57
1:B:214:PRO:O	1:B:217:LEU:N	2.38	0.57
1:D:84:ARG:NH1	3:J:51:PRO:O	2.37	0.57
1:D:139:LEU:HA	1:D:147:ARG:HD3	1.86	0.57
1:A:126:SER:HB2	1:B:80:LEU:HD21	1.86	0.57
2:E:13:DT:H2''	2:E:14:DC:O5'	2.04	0.57
3:H:33:ALA:O	3:H:34:GLN:C	2.44	0.56
2:E:14:DC:H4'	2:E:15:DA:OP1	2.06	0.56
3:K:74:LEU:O	3:K:75:GLN:C	2.42	0.56
3:J:33:ALA:O	3:J:34:GLN:C	2.43	0.56
1:D:151:PHE:CE2	1:D:155:LEU:HD11	2.41	0.55
3:H:43:ASP:CB	3:H:46:GLY:HA3	2.31	0.55
2:C:4:DT:H2''	2:C:5:DT:O5'	2.07	0.55
3:J:43:ASP:C	3:J:45:ARG:N	2.58	0.55
3:K:68:LEU:HA	3:K:72:ALA:CB	2.28	0.55
1:A:186:ARG:NH2	2:C:10:DG:N7	2.55	0.55
1:F:46:LEU:HD22	1:F:51:VAL:HG23	1.88	0.54
2:I:14:DC:H4'	2:I:15:DA:OP1	2.06	0.54
3:J:43:ASP:HB2	3:J:48:ARG:HD2	1.89	0.54
1:F:142:ARG:H	1:F:142:ARG:HE	1.55	0.54
2:G:13:DT:H2''	2:G:14:DC:O5'	2.06	0.54
3:J:65:LEU:HB3	3:J:77:TYR:HE1	1.73	0.54
1:B:137:GLU:OE2	3:K:35:ARG:NH2	2.40	0.54
1:B:151:PHE:CE2	1:B:155:LEU:HD11	2.42	0.54
1:B:167:ILE:HD11	3:K:6:TYR:CD2	2.42	0.54
1:A:133:GLU:OE2	1:B:128:ARG:NH2	2.41	0.54
3:H:44:ALA:C	3:H:46:GLY:H	2.10	0.54
2:E:30:DA:H2''	2:E:31:DT:H5'	1.89	0.54
1:F:159:PHE:CD2	1:F:171:LEU:CD2	2.92	0.53
1:A:57:TYR:HE2	1:A:63:ILE:HD13	1.72	0.53
3:J:43:ASP:O	3:J:44:ALA:C	2.44	0.53
3:K:74:LEU:HD12	3:K:75:GLN:H	1.74	0.53
1:F:171:LEU:HD12	1:F:171:LEU:C	2.29	0.53
1:F:214:PRO:O	1:F:215:VAL:C	2.46	0.53
1:A:115:PRO:HG3	1:B:115:PRO:HG3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:VAL:HA	1:A:152:LEU:HD12	1.91	0.52
2:I:13:DT:H2''	2:I:14:DC:O5'	2.10	0.52
1:A:85:SER:OG	3:K:36:LEU:CD2	2.58	0.52
1:D:15:LEU:HD22	1:D:110:ALA:HB1	1.92	0.52
3:J:68:LEU:HD23	3:J:72:ALA:HB1	1.92	0.52
2:C:8:DA:C2	2:E:26:DA:C2	2.98	0.52
1:D:188:THR:OG1	2:I:19:DG:OP2	2.26	0.51
3:H:26:GLU:HB2	3:H:28:PHE:HE1	1.75	0.51
2:C:14:DC:H2'	2:C:15:DA:C8	2.45	0.51
3:J:19:CYS:HB3	3:J:80:LEU:HD13	1.93	0.51
1:F:161:ILE:HG12	1:F:170:ASP:HB2	1.93	0.51
2:G:4:DT:H2''	2:G:5:DT:O5'	2.11	0.51
1:F:144:MET:HB2	2:G:18:DT:OP1	2.11	0.51
1:A:85:SER:OG	3:K:36:LEU:HD22	2.10	0.51
1:D:189:VAL:O	1:D:193:LEU:HG	2.11	0.51
1:A:163:SER:OG	1:A:164:PRO:HD2	2.11	0.51
1:F:82:GLY:CA	1:F:104:ILE:CD1	2.89	0.51
3:J:68:LEU:HD23	3:J:72:ALA:CB	2.40	0.51
1:D:139:LEU:HA	1:D:147:ARG:CD	2.41	0.50
3:K:78:ASN:O	3:K:81:GLN:HB3	2.11	0.50
3:J:77:TYR:O	3:J:81:GLN:HB2	2.11	0.50
1:A:38:ASP:O	1:A:87:ARG:HG2	2.11	0.50
1:F:126:SER:HA	1:F:129:ILE:HD12	1.92	0.50
1:A:161:ILE:HG12	1:A:170:ASP:HB2	1.94	0.50
1:F:93:ALA:O	1:F:94:PHE:C	2.50	0.50
1:B:172:LYS:HE3	1:B:207:LYS:HD2	1.94	0.50
1:D:46:LEU:HD22	1:D:51:VAL:HG23	1.93	0.50
3:H:40:VAL:HA	3:H:48:ARG:O	2.12	0.50
3:H:44:ALA:C	3:H:46:GLY:N	2.63	0.50
3:J:85:LYS:CE	1:A:58:GLU:CG	2.90	0.49
1:B:43:VAL:HG11	1:B:79:LEU:HD11	1.94	0.49
3:J:68:LEU:HA	3:J:72:ALA:CB	2.36	0.49
3:J:85:LYS:NZ	1:A:58:GLU:HG2	2.20	0.49
1:D:185:THR:HG21	2:I:19:DG:H2'	1.94	0.49
1:B:77:LEU:HB2	4:B:501:AKG:O4	2.12	0.49
3:H:20:SER:O	3:H:21:PHE:HB2	2.12	0.49
1:F:82:GLY:HA2	1:F:104:ILE:HD12	1.94	0.48
1:B:16:GLY:HA3	1:B:20:LEU:O	2.13	0.48
1:B:11:MET:HB3	1:B:117:LEU:HG	1.95	0.48
1:B:46:LEU:HD22	1:B:51:VAL:HG23	1.96	0.48
1:A:136:ILE:HD13	1:B:136:ILE:HD13	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:VAL:HG23	4:A:500:AKG:O1	2.13	0.48
2:C:6:DT:H1'	2:C:7:DT:H5'	1.96	0.48
3:H:43:ASP:C	3:H:46:GLY:H	2.17	0.48
3:J:58:ARG:HG3	3:J:59:MET:N	2.28	0.48
2:G:26:DA:C2	2:I:8:DA:C2	3.01	0.48
1:B:95:THR:O	1:B:97:VAL:HG23	2.14	0.48
3:J:75:GLN:CA	3:J:78:ASN:H	2.27	0.48
1:D:147:ARG:HH12	1:F:147:ARG:NH2	2.12	0.48
1:D:71:ASN:HB2	3:H:10:PRO:O	2.14	0.47
2:G:14:DC:H2'	2:G:15:DA:C8	2.49	0.47
1:A:144:MET:HB2	2:E:18:DT:OP1	2.14	0.47
1:F:18:GLY:HA3	1:F:20:LEU:HD13	1.96	0.47
3:H:39:LEU:HB2	3:H:50:GLU:HB3	1.96	0.47
3:K:82:GLN:O	3:K:86:GLN:HG3	2.15	0.47
2:G:3:DT:C6	2:G:4:DT:H72	2.48	0.47
3:J:32:TYR:O	3:J:33:ALA:HB3	2.14	0.47
3:J:57:ALA:O	3:J:61:VAL:HG23	2.14	0.47
3:H:44:ALA:O	3:H:46:GLY:N	2.48	0.47
3:H:79:GLN:O	3:H:82:GLN:HB3	2.14	0.47
1:D:179:ALA:HA	1:D:182:ILE:HG12	1.97	0.47
1:F:46:LEU:CD1	1:F:97:VAL:HG13	2.45	0.47
1:F:151:PHE:CZ	1:F:155:LEU:HD21	2.50	0.47
1:A:170:ASP:O	1:A:171:LEU:HB3	2.15	0.46
1:D:188:THR:HA	1:D:191:ARG:HE	1.80	0.46
2:G:4:DT:H4'	2:G:5:DT:OP1	2.14	0.46
3:H:66:ARG:O	3:H:69:ARG:HB3	2.15	0.46
1:B:46:LEU:CD1	1:B:97:VAL:HG13	2.45	0.46
2:C:5:DT:H1'	2:C:6:DT:C5'	2.45	0.46
1:D:160:GLY:HA2	1:D:169:ILE:HA	1.97	0.46
1:B:213:ASN:O	1:B:214:PRO:C	2.53	0.46
2:E:16:DG:H2''	2:E:17:DC:C5	2.50	0.46
1:F:94:PHE:CD2	1:F:95:THR:HG23	2.51	0.46
3:J:19:CYS:HB3	3:J:80:LEU:CD1	2.46	0.46
3:J:74:LEU:HD23	3:J:74:LEU:HA	1.81	0.46
3:H:30:THR:O	3:H:34:GLN:HA	2.16	0.46
1:B:22:LEU:HD13	1:B:102:VAL:HG12	1.96	0.46
1:F:95:THR:HB	1:F:96:PRO:HD2	1.98	0.46
2:G:2:DA:H2'	2:G:3:DT:H6	1.78	0.46
3:H:32:TYR:O	3:H:33:ALA:HB3	2.16	0.46
1:D:169:ILE:HG22	1:D:171:LEU:CD2	2.46	0.46
1:F:151:PHE:CE2	1:F:155:LEU:HD11	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:30:DA:H2''	2:E:31:DT:H3'	1.98	0.46
3:H:21:PHE:HB2	3:H:25:LYS:O	2.15	0.46
1:A:38:ASP:HB2	1:A:87:ARG:HG3	1.97	0.46
1:B:196:LEU:HB3	1:B:202:ILE:HG12	1.98	0.46
1:D:186:ARG:O	1:D:189:VAL:HB	2.16	0.46
1:F:16:GLY:O	1:F:18:GLY:O	2.33	0.46
1:A:56:VAL:HG22	1:A:62:GLU:HG2	1.98	0.46
1:A:27:PHE:HB2	1:A:97:VAL:HG13	1.97	0.45
1:D:76:VAL:N	4:D:301:AKG:O3	2.49	0.45
3:J:75:GLN:HG2	3:J:79:GLN:HG3	1.98	0.45
3:J:79:GLN:O	3:J:82:GLN:HB3	2.16	0.45
3:K:68:LEU:O	3:K:72:ALA:HB3	2.17	0.45
1:F:171:LEU:CD1	1:F:173:LEU:CD2	2.93	0.45
1:D:215:VAL:HG11	3:H:86:GLN:NE2	2.31	0.45
2:G:16:DG:H2''	2:G:17:DC:C5	2.51	0.45
1:A:214:PRO:HA	1:A:217:LEU:HD12	1.98	0.45
3:K:74:LEU:HD12	3:K:75:GLN:N	2.30	0.45
2:C:5:DT:C1'	2:C:6:DT:H5''	2.46	0.45
3:H:74:LEU:O	3:H:75:GLN:C	2.53	0.45
1:B:178:ILE:O	1:B:182:ILE:HG12	2.17	0.45
1:D:69:ARG:HH11	1:D:158:ASP:HA	1.81	0.45
1:D:169:ILE:HD12	1:D:209:ILE:HB	1.99	0.45
1:A:69:ARG:NH1	1:A:69:ARG:HB2	2.29	0.45
1:B:82:GLY:CA	1:B:104:ILE:HD13	2.43	0.45
2:C:4:DT:C4	2:C:5:DT:C4	3.05	0.44
3:K:83:VAL:HA	3:K:86:GLN:HE21	1.82	0.44
2:I:4:DT:H4'	2:I:5:DT:OP1	2.16	0.44
1:A:88:PHE:CG	3:K:35:ARG:HD3	2.52	0.44
1:D:182:ILE:HD11	1:D:189:VAL:HG22	1.98	0.44
1:F:175:HIS:HB3	1:F:186:ARG:HG2	1.98	0.44
3:J:36:LEU:HD23	3:J:36:LEU:HA	1.67	0.44
1:B:15:LEU:HD12	1:B:102:VAL:HG11	2.00	0.44
1:B:95:THR:HB	1:B:96:PRO:HD2	2.00	0.44
1:D:147:ARG:HH22	1:F:147:ARG:HH22	1.66	0.44
1:B:219:GLN:C	1:B:221:PHE:N	2.71	0.44
2:C:13:DT:H2''	2:C:14:DC:O5'	2.18	0.44
3:H:36:LEU:HD23	3:H:36:LEU:HA	1.72	0.44
1:F:82:GLY:HA2	1:F:104:ILE:HD13	1.99	0.43
2:C:4:DT:H1'	2:C:5:DT:OP1	2.18	0.43
1:F:11:MET:CB	1:F:117:LEU:HD13	2.35	0.43
1:F:213:ASN:O	1:F:214:PRO:C	2.57	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:140:ALA:O	1:F:183:GLY:CA	2.65	0.43
2:I:13:DT:H4'	2:I:14:DC:OP1	2.18	0.43
1:D:179:ALA:O	1:D:183:GLY:N	2.52	0.43
1:A:169:ILE:HG22	1:A:171:LEU:CD2	2.47	0.43
3:K:26:GLU:HB2	3:K:28:PHE:HE1	1.84	0.43
3:K:69:ARG:HG3	3:K:70:ARG:N	2.34	0.43
2:E:30:DA:C2'	2:E:31:DT:H5'	2.48	0.42
1:B:102:VAL:HA	1:B:103:PRO:HD3	1.92	0.42
3:J:74:LEU:C	3:J:78:ASN:H	2.22	0.42
1:F:154:ILE:O	1:F:155:LEU:C	2.58	0.42
2:C:18:DT:H3	2:E:15:DA:H61	1.68	0.42
3:J:45:ARG:O	3:J:47:THR:HG23	2.20	0.42
1:A:42:ARG:HG2	1:A:103:PRO:HA	2.02	0.42
3:J:22:GLY:HA2	3:J:24:SER:HA	2.02	0.42
3:J:43:ASP:HB2	3:J:48:ARG:CD	2.49	0.42
1:B:161:ILE:HG12	1:B:170:ASP:HB2	2.02	0.42
1:B:175:HIS:CE1	1:B:193:LEU:HD12	2.54	0.42
1:D:11:MET:HB2	1:D:117:LEU:HD13	2.02	0.41
1:D:22:LEU:HD23	1:D:100:PHE:HD2	1.84	0.41
1:D:170:ASP:O	1:D:171:LEU:HB3	2.20	0.41
3:H:68:LEU:CD2	3:H:72:ALA:HB1	2.40	0.41
1:B:219:GLN:NE2	1:B:219:GLN:HA	2.34	0.41
1:F:93:ALA:O	1:F:95:THR:N	2.53	0.41
1:A:151:PHE:CE2	1:A:155:LEU:HD11	2.54	0.41
2:G:16:DG:H2''	2:G:17:DC:C6	2.55	0.41
3:H:36:LEU:HD22	3:H:37:PHE:H	1.84	0.41
3:H:65:LEU:O	3:H:77:TYR:HE1	2.03	0.41
2:E:4:DT:H4'	2:E:5:DT:OP1	2.21	0.41
1:D:179:ALA:HB1	1:D:184:SER:O	2.20	0.41
1:B:219:GLN:C	1:B:221:PHE:H	2.22	0.41
1:F:145:GLY:O	1:F:149:VAL:HG23	2.20	0.41
2:G:17:DC:H42	2:I:16:DG:H1	1.67	0.41
2:E:16:DG:H2''	2:E:17:DC:C6	2.56	0.41
3:K:75:GLN:HG3	3:K:76:GLU:N	2.35	0.41
1:A:201:LEU:O	1:A:212:PHE:HB2	2.21	0.41
1:B:211:VAL:HG23	1:B:214:PRO:HD3	2.02	0.41
3:K:41:ALA:HB3	3:K:48:ARG:CB	2.51	0.41
3:J:43:ASP:O	3:J:45:ARG:N	2.53	0.41
1:A:167:ILE:O	1:A:211:VAL:HG22	2.21	0.41
1:B:78:SER:O	1:B:104:ILE:CD1	2.57	0.41
1:F:155:LEU:HB3	1:F:169:ILE:HD13	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:215:VAL:O	1:D:219:GLN:HG2	2.21	0.40
3:H:41:ALA:O	3:H:47:THR:HA	2.21	0.40
1:F:214:PRO:O	1:F:217:LEU:N	2.52	0.40
2:G:6:DT:O2	2:I:28:DA:C2	2.74	0.40
1:D:149:VAL:HA	1:D:152:LEU:HD12	2.04	0.40
1:F:146:SER:HB3	1:F:221:PHE:HE1	1.86	0.40
1:B:160:GLY:HA2	1:B:169:ILE:HA	2.02	0.40
1:D:161:ILE:HG12	1:D:170:ASP:HB2	2.03	0.40
1:A:166:GLY:HA3	1:A:211:VAL:O	2.22	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	209/222 (94%)	200 (96%)	9 (4%)	0	100	100
1	B	211/222 (95%)	202 (96%)	8 (4%)	1 (0%)	25	63
1	D	209/222 (94%)	201 (96%)	8 (4%)	0	100	100
1	F	212/222 (96%)	201 (95%)	11 (5%)	0	100	100
3	H	81/89 (91%)	72 (89%)	9 (11%)	0	100	100
3	J	80/89 (90%)	73 (91%)	7 (9%)	0	100	100
3	K	80/89 (90%)	75 (94%)	5 (6%)	0	100	100
All	All	1082/1155 (94%)	1024 (95%)	57 (5%)	1 (0%)	48	83

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	215	VAL

### 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	183/194 (94%)	174 (95%)	9 (5%)	21	43
1	B	187/194 (96%)	181 (97%)	6 (3%)	34	55
1	D	183/194 (94%)	177 (97%)	6 (3%)	33	55
1	F	187/194 (96%)	181 (97%)	6 (3%)	34	55
3	H	72/78 (92%)	67 (93%)	5 (7%)	13	34
3	J	72/78 (92%)	64 (89%)	8 (11%)	5	19
3	K	72/78 (92%)	64 (89%)	8 (11%)	5	19
All	All	956/1010 (95%)	908 (95%)	48 (5%)	20	43

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	9	LEU
1	D	38	ASP
1	D	48	LYS
1	D	63	ILE
1	D	142	ARG
1	D	186	ARG
1	F	9	LEU
1	F	13	ARG
1	F	67	LEU
1	F	106	PHE
1	F	142	ARG
1	F	163	SER
3	J	36	LEU
3	J	60	LEU
3	J	62	ASP
3	J	68	LEU
3	J	69	ARG
3	J	71	ASP
3	J	74	LEU
3	J	75	GLN

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Mol	Chain	Res	Type
3	H	19	CYS
3	H	36	LEU
3	H	60	LEU
3	H	62	ASP
3	H	71	ASP
1	A	38	ASP
1	A	55	ARG
1	A	63	ILE
1	A	69	ARG
1	A	125	LEU
1	A	142	ARG
1	A	165	ASP
1	A	186	ARG
1	A	211	VAL
1	B	13	ARG
1	B	106	PHE
1	B	117	LEU
1	B	142	ARG
1	B	163	SER
1	B	165	ASP
3	K	25	LYS
3	K	45	ARG
3	K	58	ARG
3	K	62	ASP
3	K	64	ARG
3	K	71	ASP
3	K	74	LEU
3	K	81	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
3	H	86	GLN
1	A	213	ASN
1	B	23	GLN
1	B	119	ASN
1	B	123	GLN
3	K	86	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 ⓘ

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	$\# Z  > 2$
4	AKG	F	301	-	9,9,9	6.30	3 (33%)	11,11,11	2.16	4 (36%)
4	AKG	A	500	-	9,9,9	6.19	2 (22%)	11,11,11	1.85	3 (27%)
4	AKG	D	301	-	9,9,9	6.24	2 (22%)	11,11,11	1.95	3 (27%)
4	AKG	B	501	-	9,9,9	6.33	2 (22%)	11,11,11	1.96	3 (27%)

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
4	AKG	F	301	-	-	5/9/9/9	-
4	AKG	A	500	-	-	4/9/9/9	-
4	AKG	D	301	-	-	5/9/9/9	-
4	AKG	B	501	-	-	4/9/9/9	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	501	AKG	C2-C1	-18.67	1.28	1.53
4	F	301	AKG	C2-C1	-18.56	1.28	1.53
4	D	301	AKG	C2-C1	-18.40	1.28	1.53
4	A	500	AKG	C2-C1	-18.27	1.28	1.53
4	D	301	AKG	O2-C1	-2.66	1.22	1.30
4	B	501	AKG	O2-C1	-2.61	1.23	1.30
4	A	500	AKG	O2-C1	-2.57	1.23	1.30
4	F	301	AKG	O2-C1	-2.45	1.23	1.30
4	F	301	AKG	O3-C5	2.05	1.28	1.22

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	301	AKG	O1-C1-C2	-4.95	115.11	121.72
4	B	501	AKG	O1-C1-C2	-4.05	116.31	121.72
4	A	500	AKG	C3-C2-C1	4.00	123.40	115.97
4	D	301	AKG	C3-C2-C1	3.95	123.31	115.97
4	B	501	AKG	C3-C2-C1	3.66	122.78	115.97
4	D	301	AKG	O1-C1-C2	-3.52	117.03	121.72
4	A	500	AKG	O1-C1-C2	-3.21	117.44	121.72
4	F	301	AKG	C3-C2-C1	2.85	121.26	115.97
4	B	501	AKG	O2-C1-C2	2.34	120.38	113.97
4	F	301	AKG	C4-C3-C2	2.31	117.40	113.03
4	D	301	AKG	O2-C1-C2	2.19	119.97	113.97
4	F	301	AKG	O2-C1-C2	2.12	119.78	113.97
4	A	500	AKG	O2-C1-C2	2.01	119.46	113.97

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	301	AKG	O1-C1-C2-O5
4	D	301	AKG	O1-C1-C2-C3
4	D	301	AKG	O2-C1-C2-O5
4	D	301	AKG	O2-C1-C2-C3
4	D	301	AKG	C1-C2-C3-C4
4	F	301	AKG	O2-C1-C2-C3
4	A	500	AKG	O2-C1-C2-C3
4	B	501	AKG	O2-C1-C2-C3
4	F	301	AKG	C2-C3-C4-C5
4	F	301	AKG	O1-C1-C2-O5

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Mol	Chain	Res	Type	Atoms
4	A	500	AKG	O1-C1-C2-O5
4	F	301	AKG	O1-C1-C2-C3
4	A	500	AKG	O1-C1-C2-C3
4	B	501	AKG	O1-C1-C2-C3
4	F	301	AKG	O2-C1-C2-O5
4	A	500	AKG	O2-C1-C2-O5
4	B	501	AKG	O2-C1-C2-O5
4	B	501	AKG	O1-C1-C2-O5

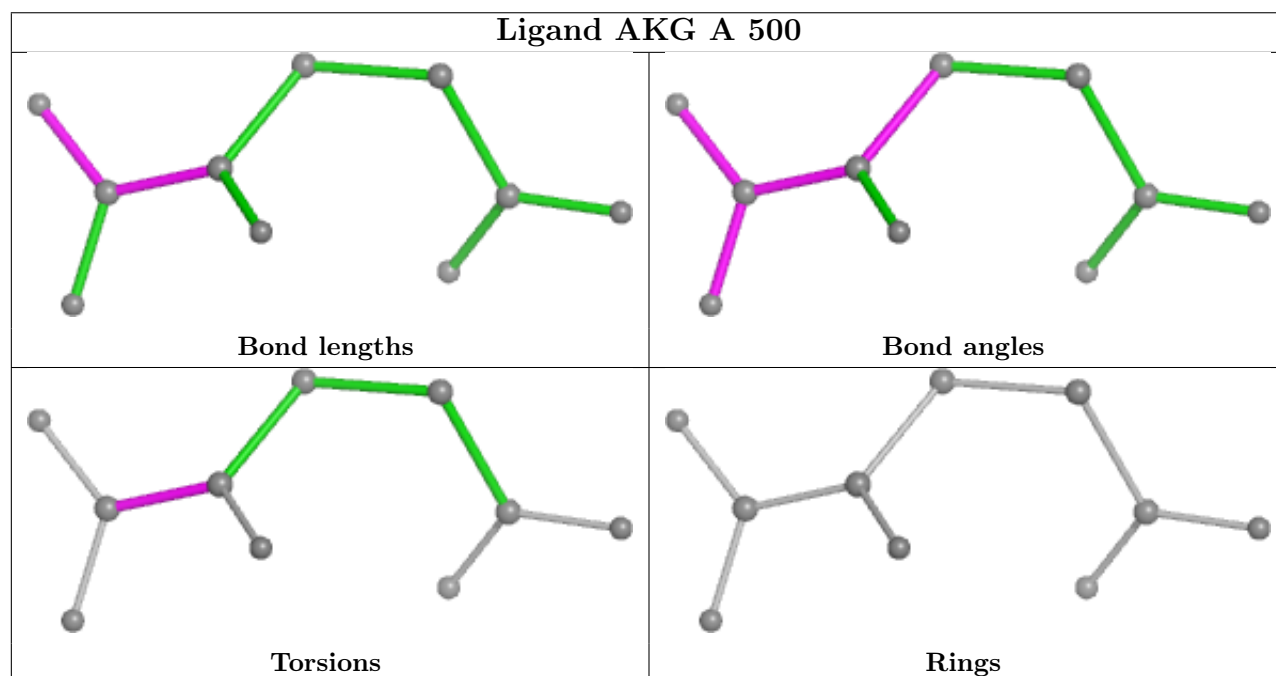
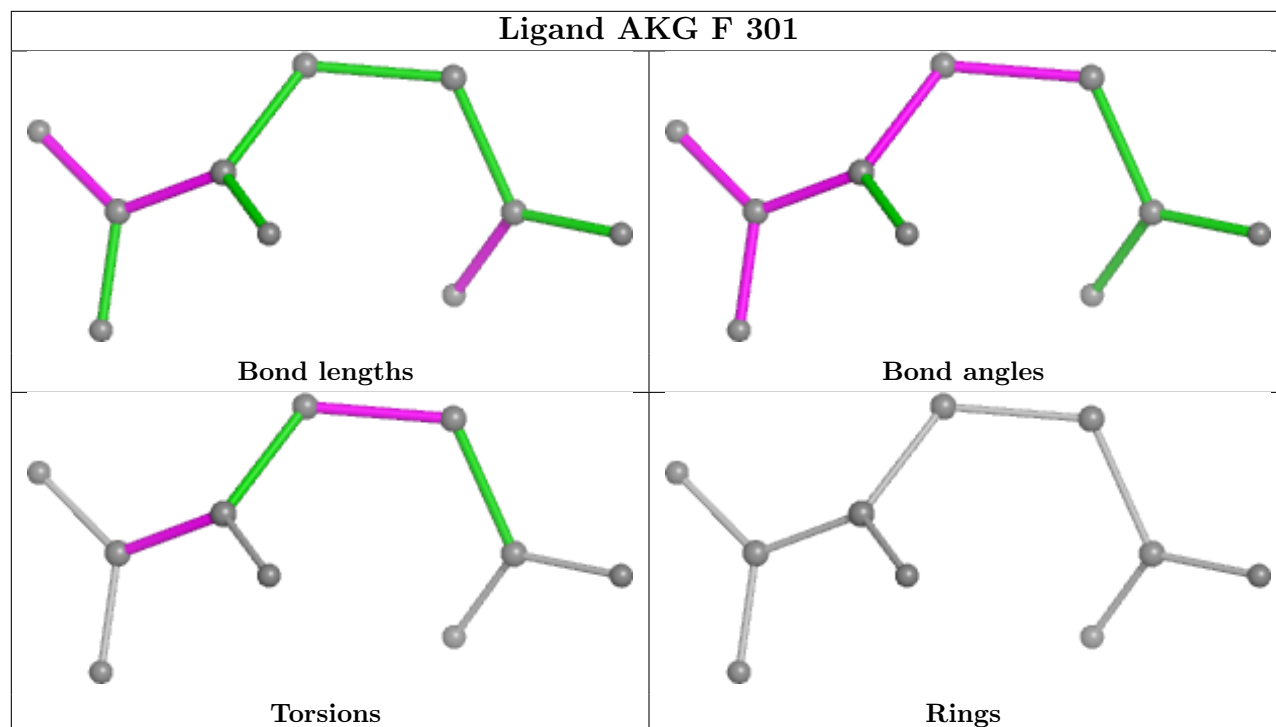
There are no ring outliers.

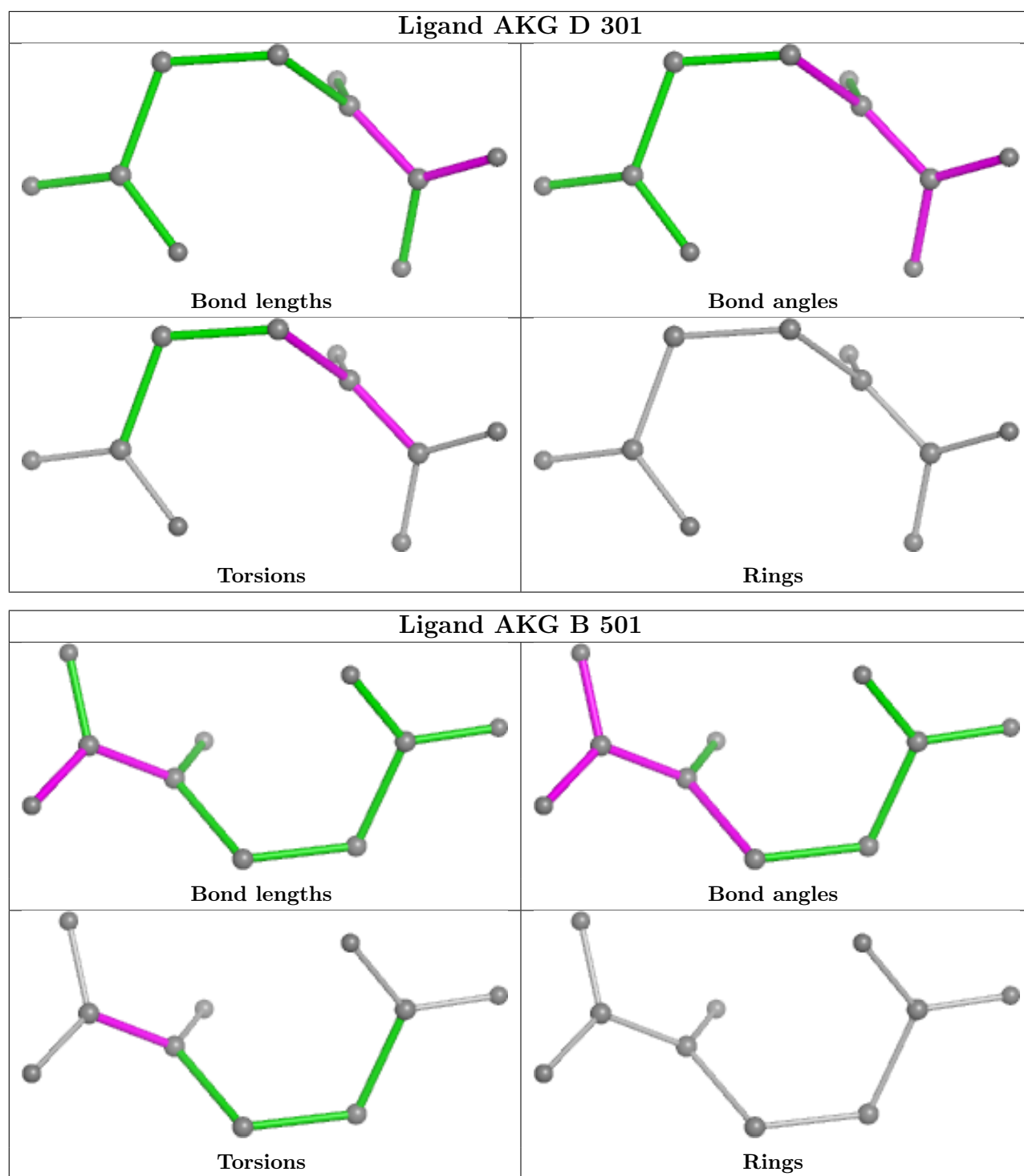
3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	500	AKG	1	0
4	D	301	AKG	3	0
4	B	501	AKG	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.







## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	213/222 (95%)	-0.61	0 100 100	67, 88, 114, 129	0
1	B	215/222 (96%)	-0.60	1 (0%) 87 76	56, 79, 105, 121	0
1	D	213/222 (95%)	-0.63	0 100 100	69, 91, 114, 135	0
1	F	216/222 (97%)	-0.56	1 (0%) 87 76	59, 85, 116, 140	0
2	C	28/30 (93%)	-0.43	0 100 100	112, 127, 151, 163	0
2	E	30/30 (100%)	-0.26	0 100 100	118, 131, 150, 160	0
2	G	30/30 (100%)	-0.36	0 100 100	93, 108, 121, 132	0
2	I	30/30 (100%)	-0.26	0 100 100	116, 139, 152, 154	0
3	H	85/89 (95%)	-0.64	0 100 100	77, 87, 109, 118	0
3	J	84/89 (94%)	-0.51	1 (1%) 76 61	84, 95, 123, 137	0
3	K	84/89 (94%)	-0.41	2 (2%) 59 45	71, 83, 111, 131	0
All	All	1228/1275 (96%)	-0.56	5 (0%) 89 78	56, 89, 129, 163	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	K	34	GLN	2.3
1	B	49	GLY	2.2
3	K	33	ALA	2.1
3	J	24	SER	2.1
1	F	81	THR	2.0

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

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

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 6.4 Ligands [i](#)

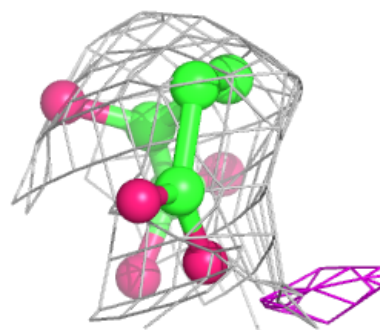
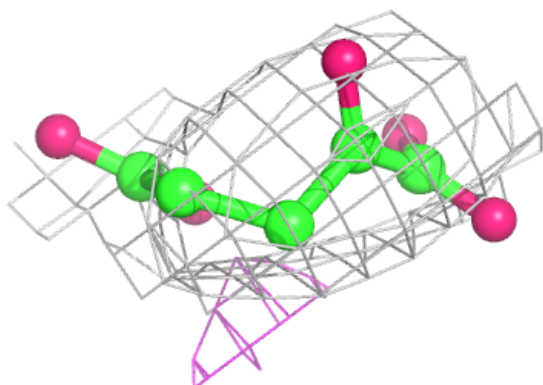
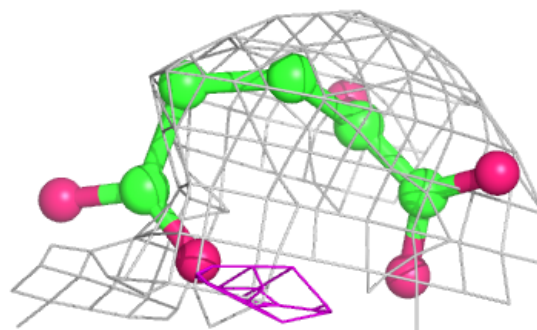
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	AKG	D	301	10/10	0.99	0.08	81,81,81,81	0
4	AKG	F	301	10/10	0.99	0.23	81,81,81,81	0
4	AKG	A	500	10/10	0.99	0.10	81,81,81,81	0
4	AKG	B	501	10/10	0.99	0.15	81,81,81,81	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.

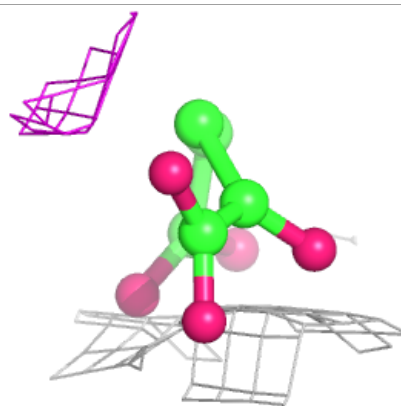
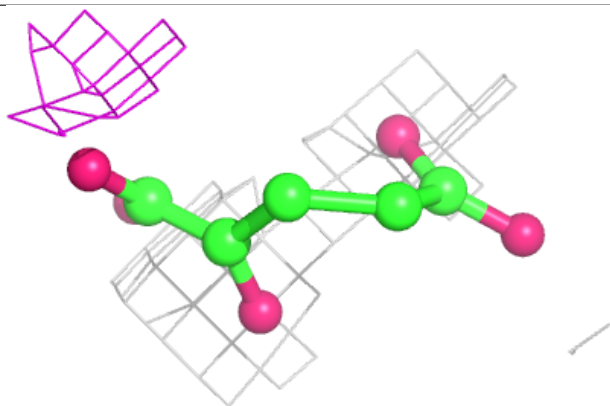
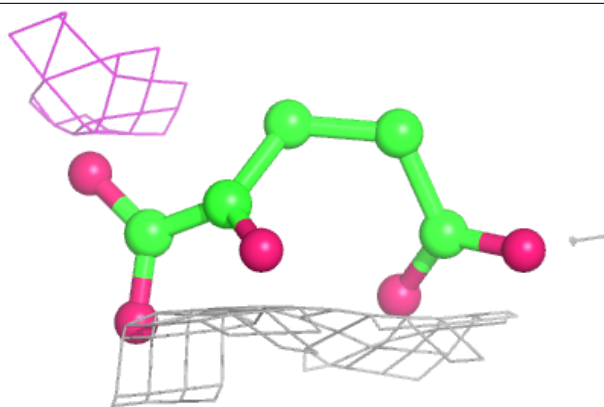
#### Electron density around AKG D 301:

2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)

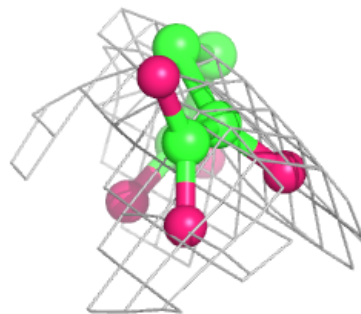
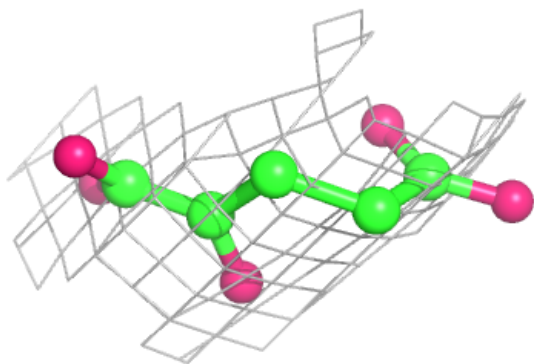
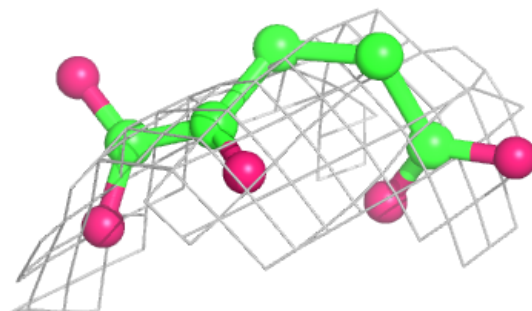


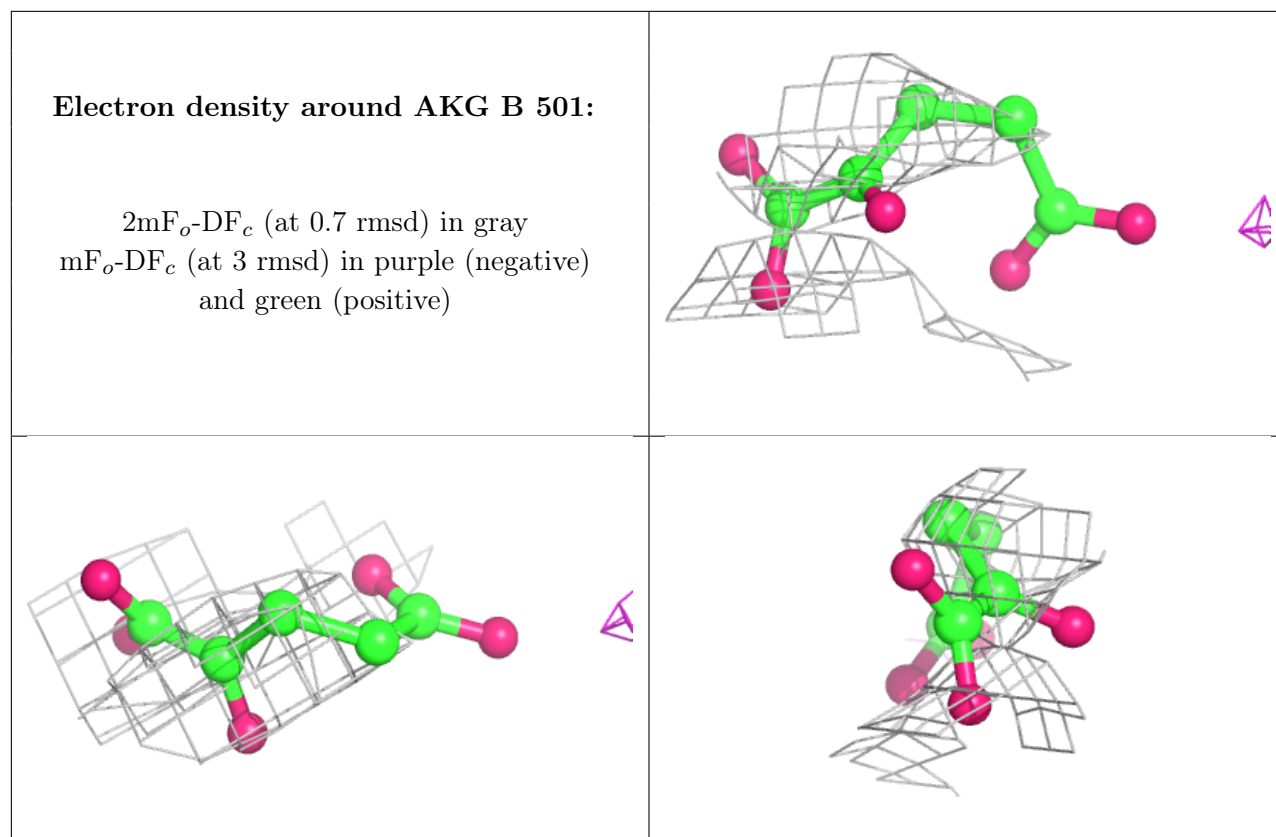
**Electron density around AKG F 301:**

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

**Electron density around AKG A 500:**

$2mF_o-DF_c$  (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.