



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 08:21 am BST

PDB ID : 1U0H
Title : STRUCTURAL BASIS FOR THE INHIBITION OF MAMMALIAN ADENYLYL CYCLASE BY MANT-GTP
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Deposited on : 2004-07-13
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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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

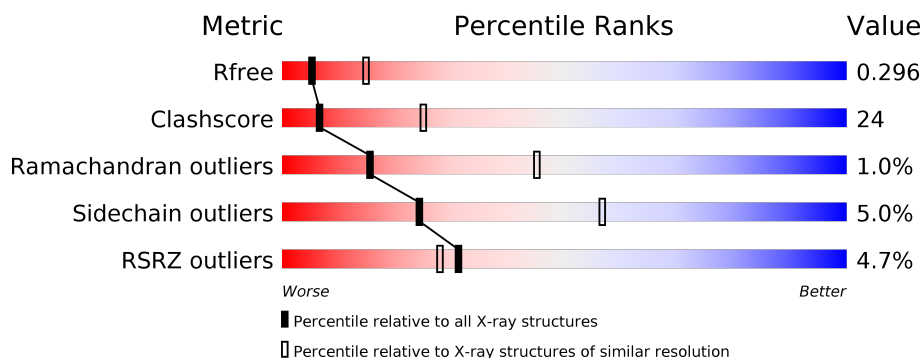
1 Overall quality at a glance

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 ≥ 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	225	<div> <div>7%</div> <div> <div>40%</div> <div>40%</div> <div>•</div> <div>16%</div> </div> </div>
2	B	212	<div> <div>6%</div> <div> <div>51%</div> <div>35%</div> <div>•</div> <div>11%</div> </div> </div>
3	C	394	<div> <div>%</div> <div> <div>49%</div> <div>32%</div> <div>•</div> <div>16%</div> </div> </div>

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 5769 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 Adenylate cyclase, type V.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	189	Total	C	N	O	S	0	0	0
			1476	929	259	271	17			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	356	MET	-	INITIATING METHIONINE	UNP P30803
A	357	HIS	-	EXPRESSION TAG	UNP P30803
A	358	HIS	-	EXPRESSION TAG	UNP P30803
A	359	HIS	-	EXPRESSION TAG	UNP P30803
A	360	HIS	-	EXPRESSION TAG	UNP P30803
A	362	HIS	VAL	CONFLICT	UNP P30803
A	476	MET	VAL	CONFLICT	UNP P30803

- Molecule 2 is a protein called Adenylate cyclase, type II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	189	Total	C	N	O	S	0	0	0
			1467	936	242	279	10			

- Molecule 3 is a protein called Guanine nucleotide-binding protein G(s), alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	330	Total	C	N	O	S	0	0	0
			2702	1714	470	505	13			

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

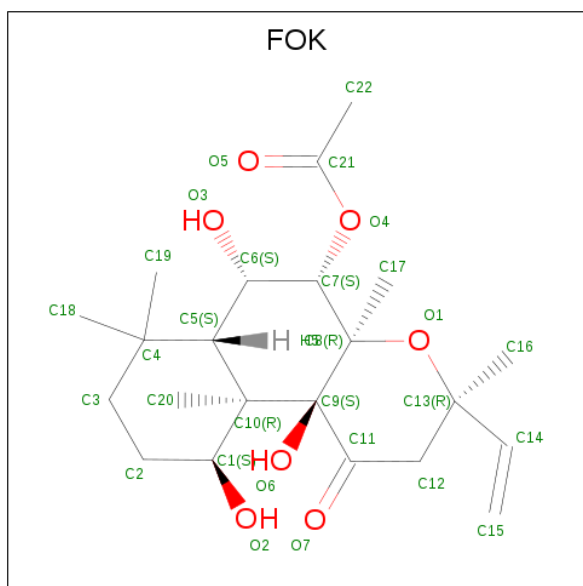
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mg	0	0
			2	2		

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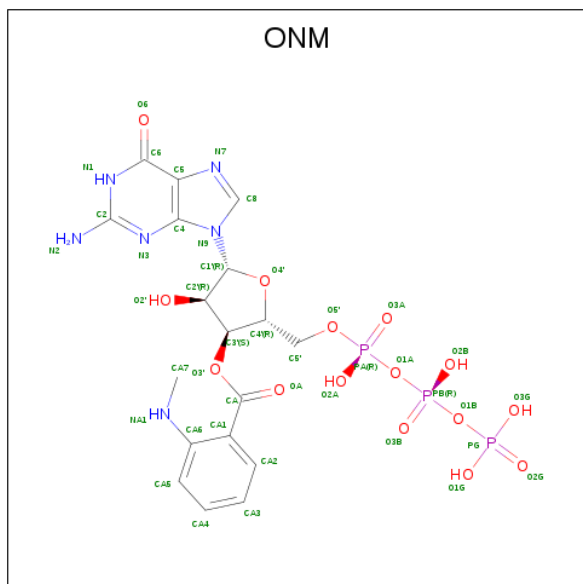
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	1	Total	Mg	0	0
			1	1		

- Molecule 5 is FORSKOLIN (three-letter code: FOK) (formula: $C_{22}H_{34}O_7$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	C O	0	0
			29	22 7		

- Molecule 6 is 3'-O-(N-METHYLANTHRANILOYL)-GUANOSINE-5'-TRIPHOSPHATE (three-letter code: ONM) (formula: $C_{18}H_{23}N_6O_{15}P_3$).

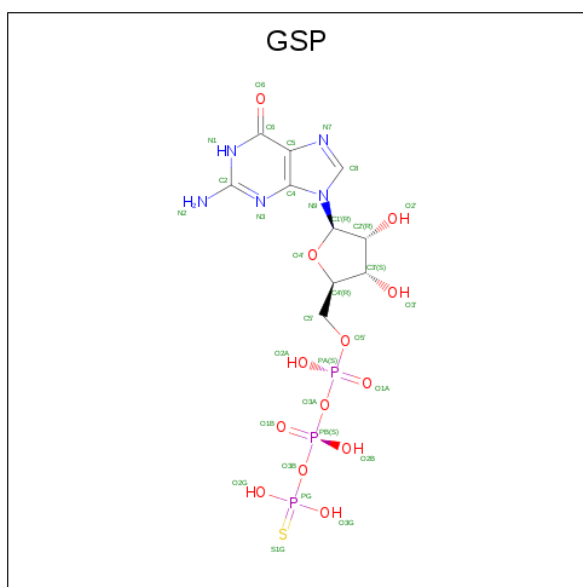


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			42	18	6	15	3		

- Molecule 7 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	C	1	Total	Cl	0	0
			1	1		

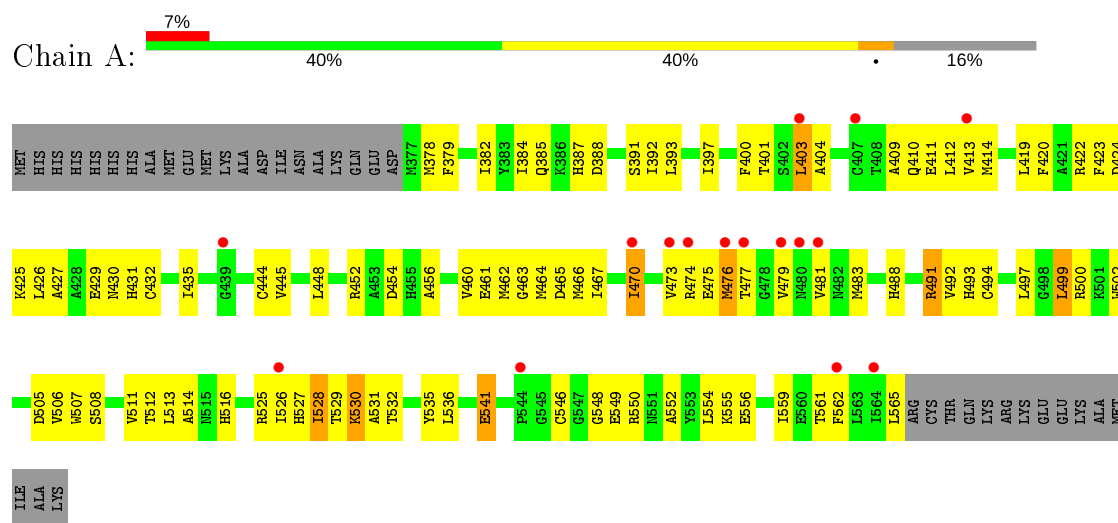
- Molecule 8 is 5'-GUANOSINE-DIPHOSPHATE-MONOTHIOPHOSPHATE (three-letter code: GSP) (formula: C₁₀H₁₆N₅O₁₃P₃S).



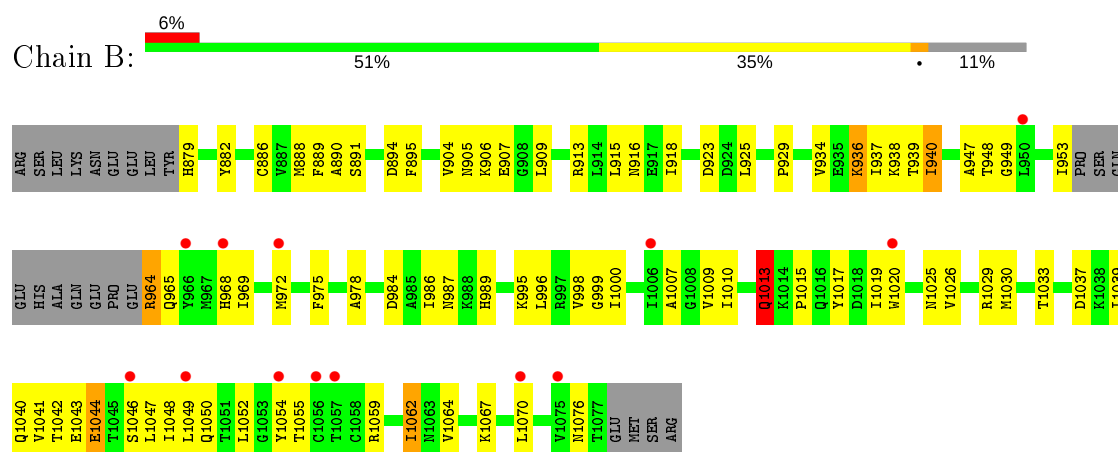
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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.

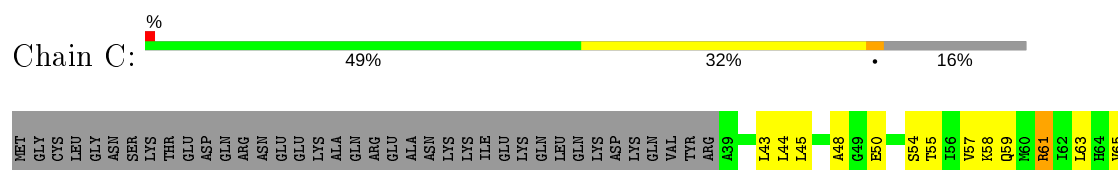
- Molecule 1: Adenylate cyclase, type V

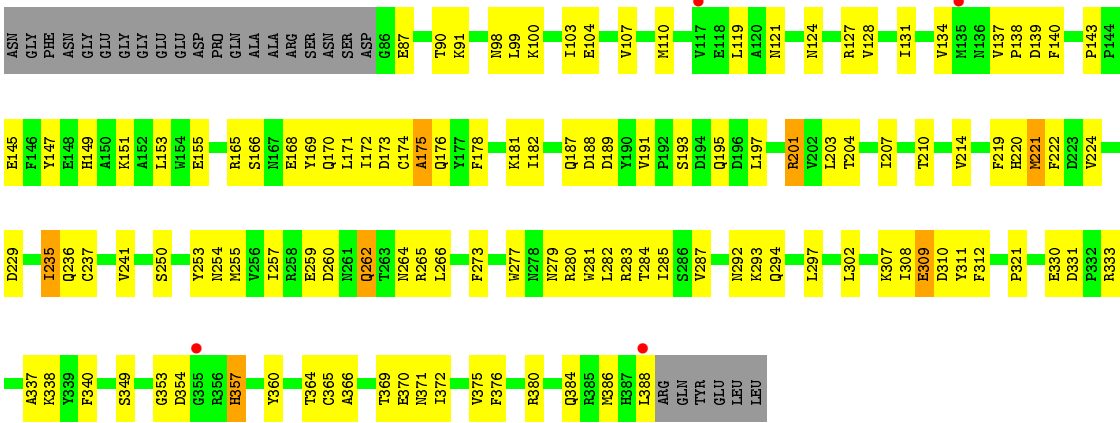


- Molecule 2: Adenylate cyclase, type II



- Molecule 3: Guanine nucleotide-binding protein G(s), alpha subunit





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	118.40 Å 133.00 Å 70.40 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.90 45.31 – 2.91	Depositor EDS
% Data completeness (in resolution range)	88.1 (15.00-2.90) 88.2 (45.31-2.91)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.91 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.245 , 0.280 0.267 , 0.296	Depositor DCC
R_{free} test set	1064 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	40.9	Xtriage
Anisotropy	1.226	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 69.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	5769	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GSP, MG, CL, FOK, ONM

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	0.38	0/1504	0.58	0/2027
2	B	0.45	0/1492	0.62	0/2014
3	C	0.40	0/2759	0.60	0/3733
All	All	0.41	0/5755	0.60	0/7774

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1476	0	1450	98	0
2	B	1467	0	1470	70	0
3	C	2702	0	2651	115	0
4	A	2	0	0	0	0
4	C	1	0	0	0	0
5	A	29	0	34	2	0
6	A	42	0	19	2	0
7	C	1	0	0	0	0
8	C	32	0	12	0	0
9	A	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	B	7	0	0	2	0
9	C	6	0	0	5	0
All	All	5769	0	5636	269	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 24.

All (269) 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:C:119:LEU:HD12	3:C:119:LEU:H	1.40	0.87
1:A:554:LEU:HA	1:A:559:ILE:HD12	1.59	0.84
3:C:207:ILE:HG12	3:C:224:VAL:HG12	1.61	0.82
1:A:530:LYS:HD2	1:A:531:ALA:N	1.93	0.81
1:A:456:ALA:HB2	1:A:535:TYR:HB3	1.61	0.81
1:A:508:SER:OG	1:A:511:VAL:HG23	1.81	0.80
2:B:1064:VAL:HG21	2:B:1070:LEU:HD12	1.61	0.79
1:A:456:ALA:HB1	1:A:536:LEU:HD12	1.63	0.79
1:A:530:LYS:HD2	1:A:531:ALA:H	1.45	0.78
3:C:99:LEU:HD11	3:C:182:ILE:HD13	1.65	0.77
3:C:127:ARG:HH11	3:C:149:HIS:HA	1.48	0.77
3:C:100:LYS:O	3:C:104:GLU:HG2	1.84	0.77
3:C:364:THR:HG22	3:C:375:VAL:HG11	1.66	0.76
3:C:257:ILE:HD12	3:C:259:GLU:HB2	1.67	0.75
2:B:953:ILE:HD12	2:B:953:ILE:H	1.51	0.74
3:C:255:MET:HB2	3:C:265:ARG:HD2	1.69	0.74
2:B:1043:GLU:O	2:B:1046:SER:HB3	1.88	0.74
1:A:404:ALA:HA	1:A:412:LEU:HD11	1.72	0.72
1:A:424:ASP:HB3	2:B:1013:GLN:HG2	1.70	0.72
5:A:101:FOK:H202	5:A:101:FOK:H193	1.71	0.71
1:A:526:ILE:HD12	1:A:526:ILE:N	2.05	0.71
1:A:391:SER:O	1:A:445:VAL:HG23	1.90	0.70
1:A:401:THR:HA	1:A:404:ALA:HB3	1.73	0.70
1:A:528:ILE:HD11	1:A:562:PHE:HB2	1.73	0.70
2:B:1007:ALA:HA	2:B:1019:ILE:HG22	1.74	0.69
1:A:529:THR:HG22	1:A:530:LYS:HD2	1.73	0.69
1:A:470:ILE:O	1:A:470:ILE:HD13	1.93	0.69
3:C:87:GLU:O	3:C:90:THR:HG22	1.92	0.69
1:A:546:CYS:O	1:A:549:GLU:HG2	1.93	0.69
1:A:378:MET:HE1	3:C:284:THR:N	2.09	0.68
1:A:473:VAL:O	1:A:477:THR:HG23	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:107:VAL:HA	3:C:110:MET:HG3	1.75	0.67
5:A:101:FOK:H201	5:A:101:FOK:H173	1.77	0.66
1:A:452:ARG:HG2	1:A:454:ASP:H	1.60	0.66
3:C:43:LEU:HD11	3:C:219:PHE:HD2	1.61	0.66
2:B:940:ILE:H	2:B:940:ILE:HD12	1.61	0.65
1:A:397:ILE:HD13	1:A:483:MET:SD	2.36	0.65
2:B:1044:GLU:HA	2:B:1047:LEU:HD12	1.78	0.64
3:C:380:ARG:O	3:C:384:GLN:HG2	1.98	0.64
3:C:58:LYS:HA	3:C:61:ARG:HH12	1.60	0.64
2:B:1009:VAL:HG11	2:B:1015:PRO:HB3	1.80	0.63
2:B:953:ILE:HD12	2:B:953:ILE:N	2.13	0.63
3:C:365:CYS:H	3:C:371:ASN:ND2	1.97	0.62
2:B:915:LEU:HA	2:B:918:ILE:HD12	1.79	0.62
1:A:378:MET:HE1	3:C:284:THR:H	1.62	0.62
1:A:500:ARG:HG3	2:B:923:ASP:HB3	1.80	0.62
3:C:121:ASN:HB2	3:C:124:ASN:HD22	1.65	0.62
1:A:411:GLU:HA	1:A:414:MET:HE2	1.81	0.61
3:C:273:PHE:HE1	3:C:287:VAL:HG11	1.66	0.60
3:C:63:LEU:HD21	3:C:369:THR:HG22	1.83	0.60
3:C:266:LEU:HD23	3:C:312:PHE:CZ	2.36	0.60
1:A:462:MET:HE2	1:A:466:MET:HG3	1.83	0.60
1:A:403:LEU:HD22	1:A:481:VAL:CG1	2.31	0.60
2:B:891:SER:HB2	9:B:8:HOH:O	2.01	0.60
1:A:463:GLY:O	1:A:467:ILE:HG12	2.00	0.60
3:C:166:SER:HA	3:C:169:TYR:CE2	2.37	0.60
1:A:384:ILE:HG22	1:A:385:GLN:N	2.17	0.59
3:C:124:ASN:OD1	3:C:127:ARG:HD2	2.03	0.59
3:C:119:LEU:HD12	3:C:119:LEU:N	2.14	0.59
2:B:1009:VAL:HG12	2:B:1010:ILE:N	2.17	0.58
1:A:552:ALA:O	1:A:556:GLU:HG3	2.03	0.58
3:C:170:GLN:HG3	9:C:400:HOH:O	2.02	0.58
1:A:387:HIS:CG	1:A:448:LEU:HD21	2.39	0.57
1:A:452:ARG:HD3	1:A:454:ASP:HB3	1.85	0.57
3:C:121:ASN:HD22	3:C:124:ASN:ND2	2.03	0.57
3:C:61:ARG:HH11	3:C:61:ARG:HB2	1.70	0.57
2:B:888:MET:HG2	2:B:889:PHE:N	2.18	0.57
2:B:906:LYS:HB2	2:B:909:LEU:HB3	1.86	0.57
1:A:474:ARG:HG3	1:A:481:VAL:HG22	1.85	0.56
3:C:307:LYS:HB2	3:C:310:ASP:OD2	2.05	0.56
1:A:528:ILE:CD1	1:A:562:PHE:HB2	2.35	0.56
1:A:460:VAL:O	1:A:464:MET:HG2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:203:LEU:HD23	3:C:204:THR:N	2.20	0.56
3:C:302:LEU:HA	3:C:333:ARG:HH21	1.69	0.56
2:B:1050:GLN:C	2:B:1052:LEU:H	2.06	0.56
2:B:1009:VAL:HG12	2:B:1010:ILE:H	1.70	0.56
1:A:423:PHE:HA	1:A:426:LEU:HD12	1.88	0.55
1:A:435:ILE:HD11	1:A:445:VAL:HG12	1.88	0.55
1:A:532:THR:O	1:A:536:LEU:HD13	2.07	0.55
3:C:45:LEU:HD11	3:C:221:MET:HG3	1.89	0.55
2:B:948:THR:HG23	2:B:972:MET:SD	2.46	0.55
3:C:99:LEU:HD21	3:C:182:ILE:HG23	1.88	0.55
3:C:219:PHE:O	3:C:220:HIS:HD2	1.89	0.55
3:C:151:LYS:O	3:C:155:GLU:HG2	2.07	0.55
3:C:371:ASN:O	3:C:375:VAL:HG22	2.07	0.55
1:A:411:GLU:HA	1:A:414:MET:CE	2.36	0.54
1:A:497:LEU:HD22	2:B:916:ASN:HA	1.89	0.54
2:B:886:CYS:HB2	2:B:969:ILE:HD13	1.89	0.54
3:C:61:ARG:HA	3:C:65:VAL:HB	1.89	0.54
3:C:143:PRO:HB2	3:C:145:GLU:OE2	2.08	0.54
1:A:384:ILE:HG22	1:A:385:GLN:H	1.73	0.54
2:B:975:PHE:O	2:B:978:ALA:HB3	2.07	0.54
3:C:337:ALA:O	3:C:340:PHE:HB3	2.08	0.54
1:A:378:MET:HE3	3:C:283:ARG:N	2.23	0.54
3:C:166:SER:HA	3:C:169:TYR:CZ	2.43	0.54
3:C:360:TYR:OH	3:C:386:MET:HG3	2.07	0.54
1:A:420:PHE:HA	1:A:423:PHE:HB2	1.90	0.53
1:A:476:MET:SD	1:A:476:MET:N	2.82	0.53
3:C:59:GLN:OE1	3:C:59:GLN:HA	2.09	0.53
3:C:124:ASN:O	3:C:128:VAL:HG23	2.08	0.53
3:C:279:ASN:O	3:C:283:ARG:HG3	2.08	0.53
1:A:529:THR:HG22	1:A:530:LYS:N	2.24	0.52
6:A:100:ONM:H3'	2:B:1025:ASN:HD22	1.74	0.52
2:B:953:ILE:CD1	2:B:953:ILE:H	2.21	0.52
1:A:425:LYS:O	1:A:429:GLU:HG3	2.09	0.52
1:A:499:LEU:HD12	1:A:500:ARG:HG2	1.92	0.52
2:B:1059:ARG:NH1	2:B:1062:ILE:HG13	2.25	0.52
1:A:435:ILE:HD11	1:A:445:VAL:CG1	2.40	0.52
1:A:527:HIS:NE2	1:A:561:THR:HB	2.25	0.52
3:C:127:ARG:O	3:C:131:ILE:HG12	2.10	0.52
3:C:207:ILE:HD13	3:C:237:CYS:SG	2.50	0.52
3:C:176:GLN:HB3	9:C:403:HOH:O	2.09	0.52
1:A:461:GLU:O	1:A:464:MET:HB2	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:493:HIS:HB2	1:A:507:TRP:O	2.10	0.51
1:A:525:ARG:NH2	1:A:565:LEU:HD23	2.25	0.51
1:A:378:MET:CE	3:C:283:ARG:N	2.74	0.51
2:B:1064:VAL:CG2	2:B:1070:LEU:HD12	2.36	0.51
6:A:100:ONM:HA3	2:B:1020:TRP:HB3	1.92	0.51
1:A:505:ASP:HB3	1:A:507:TRP:CZ2	2.45	0.51
2:B:1030:MET:HE1	2:B:1042:THR:HG23	1.92	0.51
3:C:104:GLU:HA	3:C:107:VAL:HG22	1.93	0.51
3:C:386:MET:HB2	3:C:388:LEU:HG	1.91	0.51
1:A:513:LEU:HD13	1:A:559:ILE:HD13	1.91	0.51
3:C:44:LEU:HD12	3:C:222:PHE:HB3	1.91	0.51
2:B:888:MET:HG3	2:B:1000:ILE:HG12	1.93	0.51
2:B:1026:VAL:HG22	2:B:1064:VAL:HG11	1.93	0.50
3:C:58:LYS:HA	3:C:61:ARG:NH1	2.26	0.50
2:B:1048:ILE:O	2:B:1052:LEU:HG	2.12	0.50
3:C:165:ARG:HB3	3:C:168:GLU:OE2	2.11	0.50
3:C:294:GLN:O	3:C:297:LEU:HB3	2.12	0.50
1:A:525:ARG:HH21	1:A:565:LEU:HD23	1.77	0.49
1:A:529:THR:HG22	1:A:530:LYS:CD	2.40	0.49
3:C:241:VAL:CG1	3:C:285:ILE:HD13	2.43	0.49
1:A:431:HIS:HB3	1:A:452:ARG:CZ	2.41	0.49
1:A:427:ALA:HA	1:A:462:MET:SD	2.52	0.49
2:B:949:GLY:HA3	2:B:968:HIS:HD2	1.77	0.49
1:A:382:ILE:HD12	1:A:382:ILE:O	2.13	0.49
1:A:403:LEU:HD22	1:A:481:VAL:HG12	1.93	0.49
2:B:1049:LEU:O	2:B:1054:TYR:HB2	2.13	0.49
2:B:964:ARG:NE	2:B:964:ARG:HA	2.27	0.48
3:C:308:ILE:C	3:C:310:ASP:H	2.16	0.48
2:B:906:LYS:O	2:B:907:GLU:HB2	2.13	0.48
1:A:491:ARG:NH2	2:B:907:GLU:HG3	2.29	0.48
2:B:886:CYS:O	2:B:947:ALA:HA	2.13	0.48
1:A:403:LEU:HB2	1:A:479:VAL:HG11	1.95	0.48
2:B:918:ILE:HG12	2:B:986:ILE:HD13	1.96	0.48
3:C:99:LEU:HD11	3:C:182:ILE:CD1	2.42	0.48
3:C:45:LEU:N	3:C:45:LEU:HD12	2.29	0.48
3:C:174:CYS:HB2	3:C:178:PHE:CD2	2.49	0.48
3:C:43:LEU:HD12	3:C:221:MET:CE	2.44	0.48
1:A:474:ARG:HG3	1:A:481:VAL:CG2	2.44	0.48
2:B:1067:LYS:HG2	2:B:1070:LEU:HD21	1.95	0.47
2:B:998:VAL:HG12	2:B:999:GLY:N	2.29	0.47
2:B:890:ALA:HB1	2:B:996:LEU:HD21	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:191:VAL:O	3:C:191:VAL:HG13	2.14	0.47
2:B:984:ASP:HA	2:B:987:ASN:HD22	1.79	0.47
3:C:254:ASN:HA	3:C:311:TYR:CE2	2.48	0.47
1:A:541:GLU:CG	1:A:565:LEU:HD12	2.44	0.47
2:B:929:PRO:HD2	9:B:26:HOH:O	2.14	0.47
3:C:307:LYS:HB3	3:C:309:GLU:OE2	2.14	0.47
3:C:44:LEU:HD12	3:C:222:PHE:CB	2.44	0.47
3:C:43:LEU:HD11	3:C:219:PHE:CD2	2.45	0.47
1:A:470:ILE:HD11	1:A:481:VAL:HG23	1.96	0.47
3:C:119:LEU:H	3:C:119:LEU:CD1	2.16	0.47
2:B:1007:ALA:CA	2:B:1019:ILE:HG22	2.44	0.47
3:C:166:SER:HB2	3:C:171:LEU:HD22	1.96	0.47
2:B:1043:GLU:O	2:B:1047:LEU:HG	2.15	0.47
2:B:1055:THR:OG1	2:B:1076:ASN:HB2	2.15	0.47
2:B:1049:LEU:HA	2:B:1052:LEU:HD12	1.97	0.47
3:C:166:SER:HB2	3:C:171:LEU:CD2	2.45	0.47
3:C:250:SER:HB2	3:C:292:ASN:O	2.14	0.47
1:A:378:MET:HE3	3:C:283:ARG:H	1.80	0.46
2:B:895:PHE:CZ	2:B:915:LEU:HB2	2.51	0.46
1:A:548:GLY:O	1:A:555:LYS:HB2	2.15	0.46
2:B:995:LYS:HD2	2:B:1037:ASP:OD2	2.16	0.46
3:C:121:ASN:HB2	3:C:124:ASN:ND2	2.28	0.46
3:C:203:LEU:HD23	3:C:204:THR:O	2.16	0.46
3:C:284:THR:C	3:C:357:HIS:HB3	2.36	0.46
3:C:376:PHE:O	3:C:380:ARG:HG3	2.15	0.46
2:B:1007:ALA:HB1	2:B:1017:TYR:OH	2.16	0.46
3:C:176:GLN:N	9:C:403:HOH:O	2.49	0.46
3:C:264:ASN:OD1	3:C:266:LEU:N	2.49	0.46
2:B:882:TYR:CD1	2:B:882:TYR:N	2.84	0.46
3:C:57:VAL:HG22	3:C:221:MET:HG2	1.98	0.46
2:B:1029:ARG:HH12	2:B:1033:THR:HG23	1.82	0.45
3:C:103:ILE:HG23	3:C:104:GLU:N	2.31	0.45
1:A:526:ILE:CD1	1:A:526:ILE:N	2.75	0.45
3:C:131:ILE:HG13	3:C:153:LEU:CD1	2.46	0.45
1:A:470:ILE:HD13	1:A:470:ILE:C	2.36	0.45
3:C:173:ASP:C	3:C:175:ALA:H	2.19	0.45
1:A:382:ILE:HG12	2:B:913:ARG:NH1	2.31	0.45
1:A:494:CYS:HB3	1:A:506:VAL:HG12	1.97	0.45
2:B:1030:MET:CE	2:B:1041:VAL:C	2.85	0.45
2:B:949:GLY:HA3	2:B:968:HIS:CD2	2.52	0.45
3:C:330:GLU:OE2	3:C:338:LYS:HE2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:55:THR:HG22	3:C:366:ALA:HB1	1.99	0.45
2:B:923:ASP:C	2:B:925:LEU:H	2.19	0.44
2:B:1000:ILE:HD12	2:B:1039:ILE:HG21	1.98	0.44
1:A:550:ARG:NH1	1:A:550:ARG:HB2	2.33	0.44
3:C:187:GLN:HB3	3:C:189:ASP:OD1	2.18	0.44
3:C:292:ASN:HA	3:C:364:THR:O	2.17	0.44
1:A:392:ILE:HD11	1:A:492:VAL:CG1	2.48	0.44
1:A:382:ILE:HG22	2:B:916:ASN:ND2	2.32	0.44
1:A:397:ILE:HD11	1:A:419:LEU:CD2	2.48	0.44
2:B:934:VAL:HA	2:B:948:THR:HG22	2.00	0.44
2:B:936:LYS:HE3	2:B:939:THR:HG23	2.00	0.44
3:C:195:GLN:HA	3:C:195:GLN:NE2	2.33	0.43
1:A:400:PHE:HA	1:A:403:LEU:HD23	1.99	0.43
3:C:193:SER:O	3:C:197:LEU:HG	2.17	0.43
3:C:255:MET:CB	3:C:265:ARG:HD2	2.45	0.43
3:C:59:GLN:HG3	3:C:372:ILE:HG13	2.00	0.43
2:B:904:VAL:HG11	3:C:207:ILE:HB	2.00	0.43
3:C:277:TRP:NE1	3:C:349:SER:HA	2.33	0.43
2:B:989:HIS:CG	3:C:280:ARG:HG2	2.54	0.43
3:C:210:THR:HB	3:C:221:MET:HB3	2.01	0.43
1:A:488:HIS:HB3	1:A:514:ALA:HB2	2.01	0.43
1:A:530:LYS:CD	1:A:531:ALA:N	2.75	0.43
2:B:1050:GLN:C	2:B:1052:LEU:N	2.72	0.43
3:C:214:VAL:HG11	3:C:380:ARG:CZ	2.49	0.42
1:A:378:MET:HE3	3:C:282:LEU:CA	2.49	0.42
1:A:422:ARG:O	1:A:426:LEU:HG	2.19	0.42
1:A:409:ALA:O	1:A:413:VAL:HG23	2.18	0.42
1:A:529:THR:HG22	1:A:530:LYS:H	1.84	0.42
2:B:1040:GLN:OE1	2:B:1059:ARG:HD2	2.20	0.42
1:A:422:ARG:HH21	1:A:425:LYS:CD	2.32	0.42
1:A:530:LYS:N	1:A:530:LYS:HD2	2.34	0.42
3:C:254:ASN:HB2	3:C:311:TYR:CZ	2.53	0.42
3:C:365:CYS:HB2	9:C:399:HOH:O	2.18	0.42
1:A:456:ALA:O	1:A:460:VAL:HG23	2.19	0.42
1:A:493:HIS:O	1:A:507:TRP:HE3	2.02	0.42
2:B:1030:MET:HE3	2:B:1041:VAL:C	2.40	0.42
2:B:1033:THR:HB	2:B:1059:ARG:HH12	1.85	0.42
1:A:475:GLU:HB2	1:A:476:MET:SD	2.60	0.42
2:B:1030:MET:HE1	2:B:1042:THR:CG2	2.49	0.42
1:A:499:LEU:HA	1:A:502:TRP:HE1	1.84	0.42
3:C:253:TYR:HB3	9:C:401:HOH:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:331:ASP:OD1	3:C:333:ARG:HG3	2.19	0.42
1:A:432:CYS:SG	1:A:462:MET:HB2	2.60	0.42
3:C:292:ASN:CG	3:C:293:LYS:H	2.23	0.42
1:A:512:THR:HG22	1:A:516:HIS:HD2	1.85	0.41
1:A:513:LEU:CD1	1:A:559:ILE:HD13	2.50	0.41
3:C:172:ILE:HD12	3:C:174:CYS:SG	2.60	0.41
3:C:235:ILE:HD13	3:C:235:ILE:O	2.20	0.41
1:A:426:LEU:HD13	1:A:465:ASP:HB3	2.01	0.41
1:A:456:ALA:HB1	1:A:536:LEU:CD1	2.43	0.41
3:C:260:ASP:OD1	3:C:262:GLN:N	2.53	0.41
3:C:50:GLU:HG2	3:C:201:ARG:NH2	2.36	0.41
3:C:147:TYR:HD1	3:C:182:ILE:HG22	1.86	0.41
1:A:393:LEU:HD23	1:A:393:LEU:C	2.40	0.41
2:B:998:VAL:CG1	2:B:999:GLY:N	2.84	0.41
3:C:281:TRP:O	3:C:282:LEU:HD23	2.19	0.41
2:B:905:ASN:HB3	3:C:236:GLN:OE1	2.20	0.41
3:C:134:VAL:HA	3:C:137:VAL:HG23	2.02	0.41
3:C:134:VAL:HB	3:C:140:PHE:CE1	2.55	0.41
3:C:254:ASN:HA	3:C:311:TYR:CD2	2.56	0.41
1:A:435:ILE:HG13	1:A:444:CYS:HA	2.03	0.41
3:C:90:THR:HG23	3:C:91:LYS:HG3	2.02	0.41
1:A:410:GLN:O	1:A:414:MET:HG3	2.21	0.41
1:A:507:TRP:O	1:A:508:SER:HB3	2.21	0.41
1:A:414:MET:HB2	1:A:414:MET:HE2	1.88	0.40
1:A:379:PHE:CE1	3:C:281:TRP:HB3	2.56	0.40
3:C:54:SER:O	3:C:58:LYS:HG3	2.22	0.40
1:A:530:LYS:HA	1:A:562:PHE:HE1	1.85	0.40
2:B:937:ILE:O	2:B:938:LYS:HB3	2.21	0.40
3:C:119:LEU:HD21	3:C:128:VAL:HG21	2.03	0.40
3:C:121:ASN:HD22	3:C:124:ASN:HD22	1.68	0.40
3:C:353:GLY:HA2	3:C:357:HIS:NE2	2.36	0.40
2:B:889:PHE:O	2:B:998:VAL:HA	2.21	0.40
3:C:280:ARG:HA	3:C:283:ARG:NE	2.36	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	187/225 (83%)	171 (91%)	15 (8%)	1 (0%)	29	61
2	B	185/212 (87%)	171 (92%)	13 (7%)	1 (0%)	29	61
3	C	326/394 (83%)	292 (90%)	29 (9%)	5 (2%)	10	34
All	All	698/831 (84%)	634 (91%)	57 (8%)	7 (1%)	15	45

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	1013	GLN
3	C	138	PRO
3	C	321	PRO
3	C	175	ALA
1	A	430	ASN
3	C	48	ALA
3	C	309	GLU

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	158/189 (84%)	149 (94%)	9 (6%)	20	51
2	B	162/184 (88%)	153 (94%)	9 (6%)	21	52
3	C	297/351 (85%)	284 (96%)	13 (4%)	28	61
All	All	617/724 (85%)	586 (95%)	31 (5%)	24	57

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	388	ASP
1	A	403	LEU
1	A	470	ILE
1	A	476	MET
1	A	491	ARG
1	A	499	LEU
1	A	528	ILE
1	A	530	LYS
1	A	541	GLU
2	B	879	HIS
2	B	894	ASP
2	B	936	LYS
2	B	940	ILE
2	B	964	ARG
2	B	965	GLN
2	B	1013	GLN
2	B	1044	GLU
2	B	1062	ILE
3	C	61	ARG
3	C	98	ASN
3	C	139	ASP
3	C	181	LYS
3	C	188	ASP
3	C	201	ARG
3	C	221	MET
3	C	229	ASP
3	C	235	ILE
3	C	262	GLN
3	C	354	ASP
3	C	357	HIS
3	C	370	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	385	GLN
1	A	387	HIS
1	A	455	HIS
1	A	515	ASN
2	B	989	HIS
2	B	1001	ASN

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Mol	Chain	Res	Type
2	B	1013	GLN
2	B	1025	ASN
2	B	1050	GLN
2	B	1076	ASN
3	C	97	ASN
3	C	121	ASN
3	C	220	HIS
3	C	362	HIS
3	C	371	ASN
3	C	377	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 for Mogul analysis.

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	$\# Z > 2$
6	ONM	A	100	4	37,45,45	3.64	16 (43%)	48,69,69	3.20	23 (47%)
5	FOK	A	101	-	28,31,31	3.42	9 (32%)	35,54,54	1.63	6 (17%)
8	GSP	C	395	4	26,34,34	1.74	6 (23%)	28,54,54	2.74	9 (32%)

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
6	ONM	A	100	4	-	4/28/48/48	0/4/4/4
5	FOK	A	101	-	-	4/7/80/80	0/3/3/3
8	GSP	C	395	4	-	2/17/38/38	0/3/3/3

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	101	FOK	O7-C11	13.75	1.43	1.21
6	A	100	ONM	CA4-CA5	10.27	1.60	1.38
6	A	100	ONM	OA-CA	9.06	1.47	1.22
6	A	100	ONM	CA3-CA2	8.72	1.57	1.38
6	A	100	ONM	CA5-CA6	7.86	1.53	1.39
5	A	101	FOK	C10-C5	6.02	1.67	1.56
5	A	101	FOK	C4-C5	-5.91	1.48	1.56
6	A	100	ONM	C6-N1	5.82	1.43	1.33
8	C	395	GSP	C6-N1	4.36	1.40	1.33
5	A	101	FOK	C9-C11	4.19	1.58	1.53
6	A	100	ONM	PB-O3B	4.00	1.65	1.50
6	A	100	ONM	PA-O3A	3.80	1.64	1.50
6	A	100	ONM	O4'-C1'	3.80	1.46	1.41
6	A	100	ONM	C2-N1	3.72	1.42	1.35
8	C	395	GSP	C2-N1	3.11	1.40	1.35
6	A	100	ONM	PG-O2G	3.11	1.60	1.50
8	C	395	GSP	O4'-C1'	2.99	1.45	1.41
6	A	100	ONM	CA1-CA	2.87	1.56	1.50
8	C	395	GSP	C8-N7	-2.59	1.30	1.34
5	A	101	FOK	C17-C8	2.53	1.56	1.51
6	A	100	ONM	CA3-CA4	2.47	1.44	1.38
6	A	100	ONM	O3'-C3'	2.45	1.48	1.44
6	A	100	ONM	CA7-NA1	2.43	1.49	1.45
5	A	101	FOK	C10-C1	2.39	1.59	1.55
6	A	100	ONM	CA2-CA1	2.37	1.43	1.39
5	A	101	FOK	C20-C10	2.32	1.58	1.53
5	A	101	FOK	O1-C13	2.27	1.49	1.45
8	C	395	GSP	C6-C5	2.26	1.45	1.41
6	A	100	ONM	C4-N3	2.24	1.39	1.35
5	A	101	FOK	C3-C4	2.18	1.58	1.54
8	C	395	GSP	C2'-C3'	2.18	1.59	1.53

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	C	395	GSP	C5-C6-N1	-9.96	109.81	123.43
6	A	100	ONM	O1G-PG-O3G	-8.54	74.98	107.64
6	A	100	ONM	PB-O1B-PG	-8.25	104.50	132.83
8	C	395	GSP	C6-N1-C2	6.75	126.66	115.93
6	A	100	ONM	C2-N3-C4	6.36	122.63	115.36
6	A	100	ONM	N3-C2-N1	-6.31	118.80	127.22
6	A	100	ONM	O3G-PG-O1B	5.51	123.12	104.64
6	A	100	ONM	C3'-O3'-CA	5.25	125.63	117.21
6	A	100	ONM	O1G-PG-O1B	-5.03	87.78	104.64
6	A	100	ONM	O3'-CA-CA1	4.88	119.29	111.69
5	A	101	FOK	C20-C10-C1	-4.36	101.21	107.68
6	A	100	ONM	C6-C5-C4	-4.31	116.68	120.80
5	A	101	FOK	C10-C5-C4	-4.14	113.01	116.46
6	A	100	ONM	CA3-CA4-CA5	-4.13	113.90	120.19
6	A	100	ONM	C5'-C4'-C3'	3.83	127.09	114.40
6	A	100	ONM	CA7-NA1-CA6	3.59	128.06	122.44
8	C	395	GSP	N3-C2-N1	-3.43	122.65	127.22
5	A	101	FOK	C2-C3-C4	3.42	118.80	113.38
5	A	101	FOK	C13-O1-C8	3.42	125.27	119.84
6	A	100	ONM	O5'-C5'-C4'	3.37	120.61	108.99
6	A	100	ONM	C1'-N9-C4	-3.31	120.83	126.64
6	A	100	ONM	PB-O1A-PA	-3.15	122.03	132.83
8	C	395	GSP	O4'-C1'-C2'	-2.77	102.88	106.93
8	C	395	GSP	C2-N3-C4	-2.71	112.26	115.36
5	A	101	FOK	C19-C4-C18	-2.46	104.26	107.89
6	A	100	ONM	O4'-C1'-C2'	2.43	110.48	106.93
8	C	395	GSP	PA-O3A-PB	-2.42	124.53	132.83
6	A	100	ONM	O2B-PB-O3B	-2.40	100.36	112.24
8	C	395	GSP	C1'-N9-C4	-2.40	122.42	126.64
6	A	100	ONM	N2-C2-N1	2.36	120.93	117.25
6	A	100	ONM	O3'-CA-OA	-2.33	119.73	123.53
6	A	100	ONM	O5'-PA-O3A	-2.30	100.08	109.07
6	A	100	ONM	C6-N1-C2	2.20	119.42	115.93
8	C	395	GSP	O3G-PG-O3B	2.14	111.79	104.64
8	C	395	GSP	O2A-PA-O1A	2.10	122.64	112.24
6	A	100	ONM	CA4-CA5-CA6	2.08	122.94	118.62
5	A	101	FOK	C19-C4-C5	2.06	119.06	111.86
6	A	100	ONM	C4-C5-N7	-2.06	107.26	109.40

There are no chirality outliers.

All (10) torsion outliers are listed below:

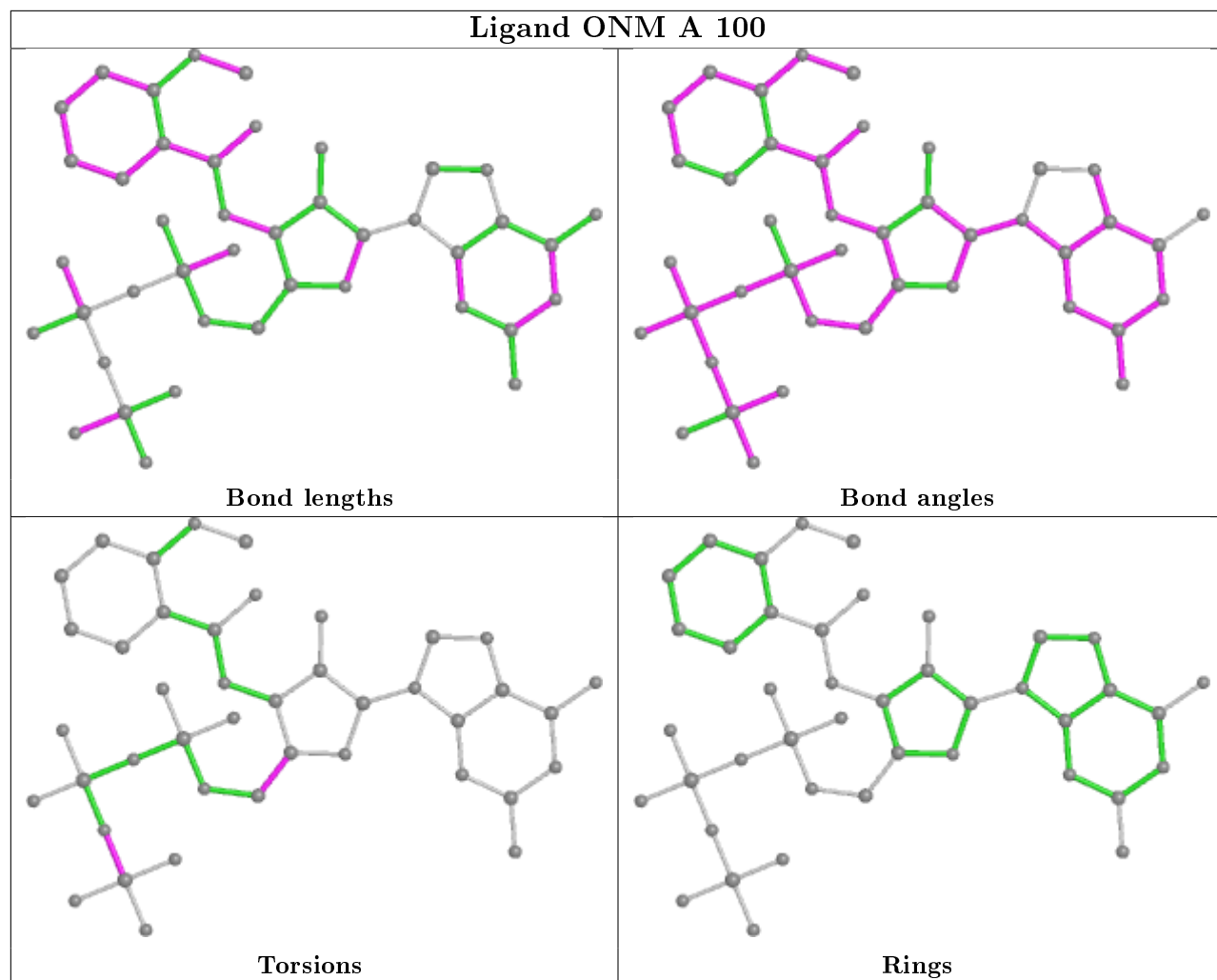
Mol	Chain	Res	Type	Atoms
6	A	100	ONM	O4'-C4'-C5'-O5'
5	A	101	FOK	C16-C13-C14-C15
5	A	101	FOK	C22-C21-O4-C7
5	A	101	FOK	O5-C21-O4-C7
6	A	100	ONM	C3'-C4'-C5'-O5'
5	A	101	FOK	O1-C13-C14-C15
8	C	395	GSP	PA-O3A-PB-O1B
6	A	100	ONM	PB-O1B-PG-O2G
6	A	100	ONM	PB-O1B-PG-O1G
8	C	395	GSP	PA-O3A-PB-O2B

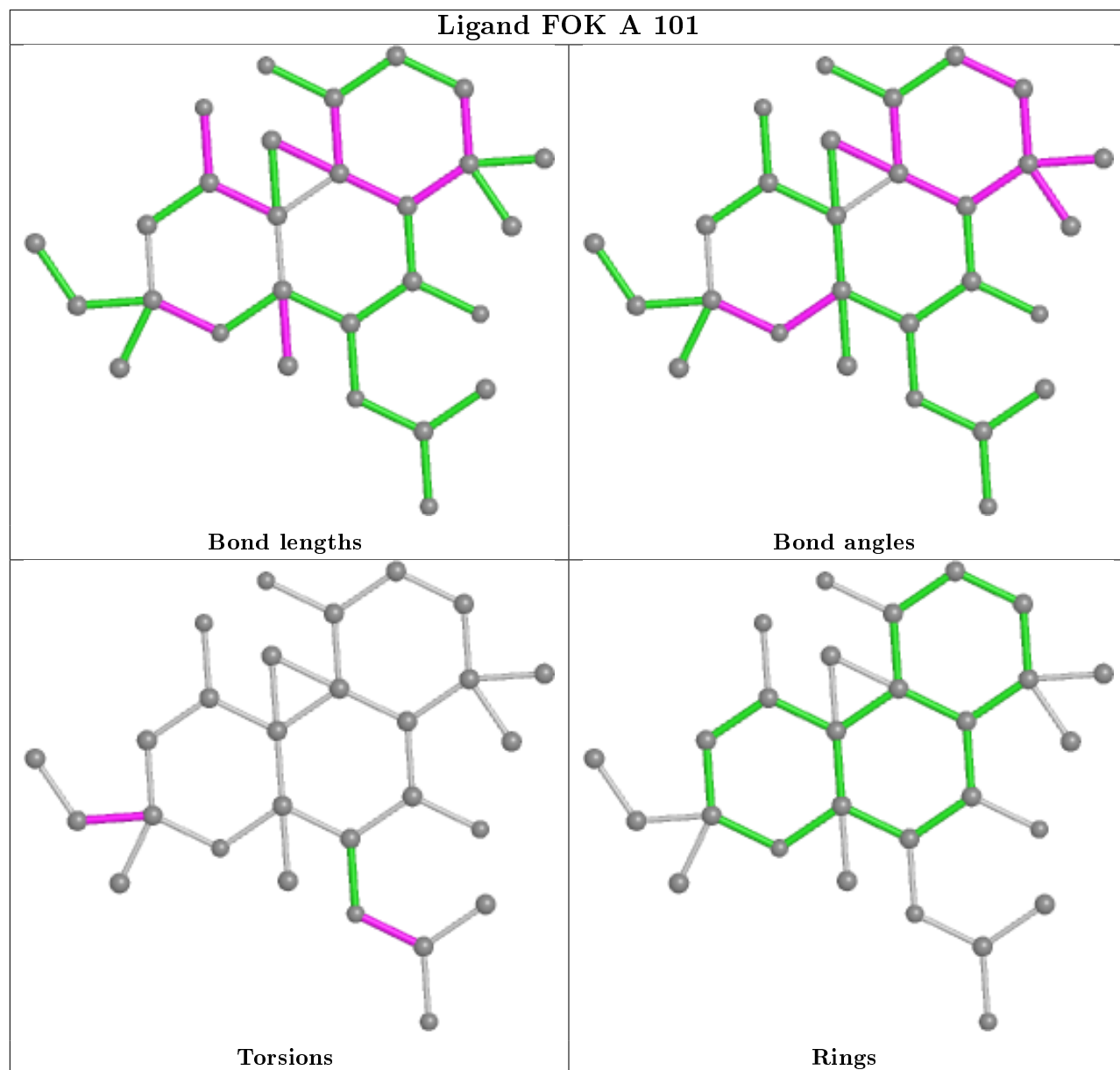
There are no ring outliers.

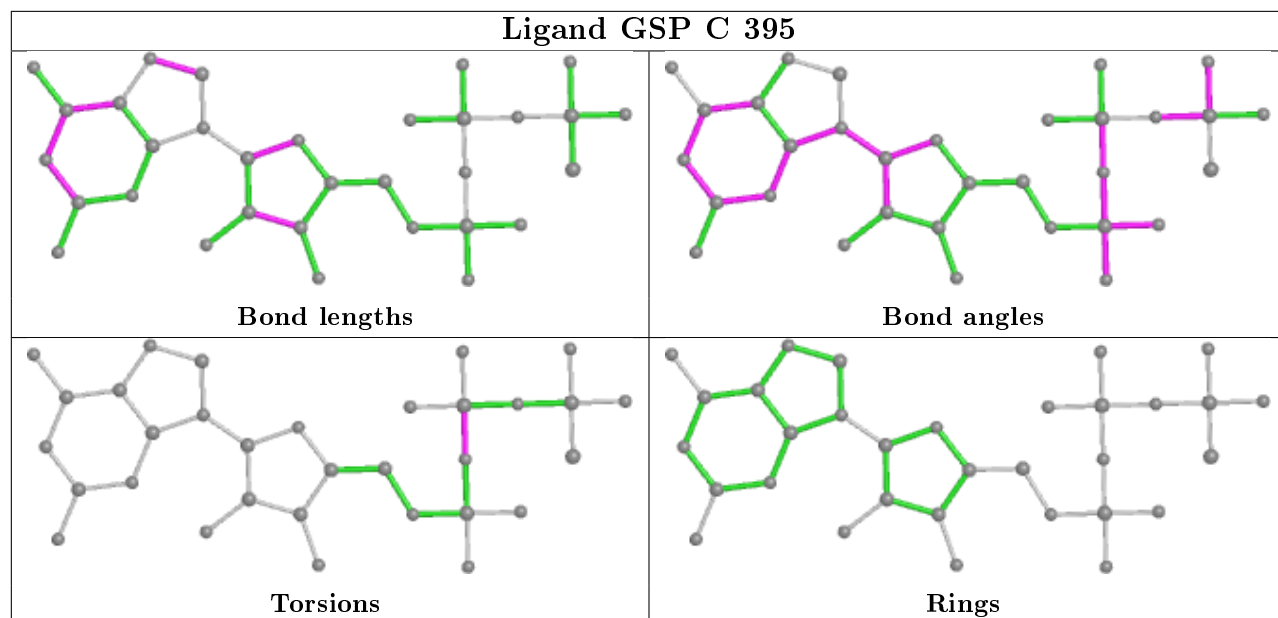
2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	100	ONM	2	0
5	A	101	FOK	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 [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	189/225 (84%)	0.80	16 (8%) 10 8	27, 64, 97, 116	0
2	B	189/212 (89%)	0.72	13 (6%) 16 13	20, 46, 80, 103	0
3	C	330/394 (83%)	0.28	4 (1%) 79 79	26, 49, 75, 103	0
All	All	708/831 (85%)	0.54	33 (4%) 31 28	20, 52, 84, 116	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	480	ASN	6.1
1	A	477	THR	5.3
1	A	479	VAL	4.9
2	B	1056	CYS	3.3
1	A	403	LEU	3.3
2	B	966	TYR	3.3
1	A	564	ILE	3.2
1	A	473	VAL	3.1
3	C	388	LEU	3.1
2	B	1006	ILE	3.0
2	B	1057	THR	2.9
2	B	1075	VAL	2.8
1	A	407	CYS	2.8
1	A	476	MET	2.7
1	A	481	VAL	2.7
1	A	526	ILE	2.7
1	A	439	GLY	2.6
2	B	1046	SER	2.6
1	A	470	ILE	2.5
2	B	1020	TRP	2.4
2	B	1070	LEU	2.4
2	B	1054	TYR	2.3
3	C	135	MET	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	544	PRO	2.3
2	B	968	HIS	2.1
1	A	562	PHE	2.1
2	B	950	LEU	2.1
3	C	117	VAL	2.1
1	A	474	ARG	2.1
1	A	413	VAL	2.1
2	B	972	MET	2.0
3	C	355	GLY	2.0
2	B	1049	LEU	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 carbohydrates in this entry.

6.4 Ligands [i](#)

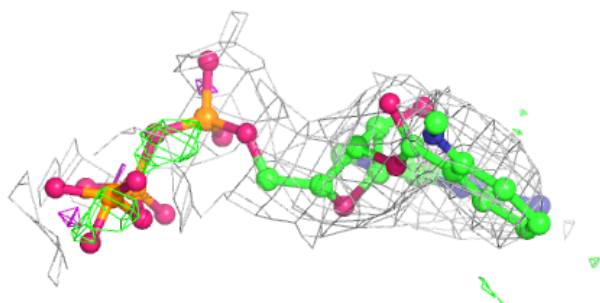
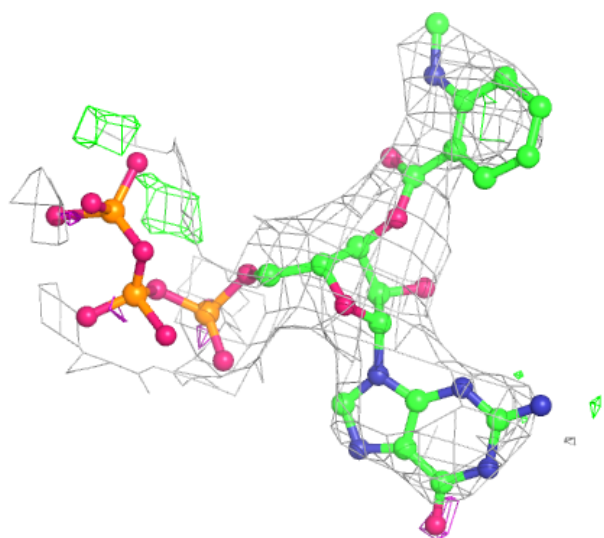
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
6	ONM	A	100	42/42	0.81	0.28	54,74,82,89	0
4	MG	A	581	1/1	0.85	0.15	38,38,38,38	0
7	CL	C	397	1/1	0.87	0.18	49,49,49,49	0
5	FOK	A	101	29/29	0.90	0.29	30,36,44,44	0
4	MG	C	396	1/1	0.90	0.18	21,21,21,21	0
4	MG	A	582	1/1	0.91	0.09	38,38,38,38	0
8	GSP	C	395	32/32	0.92	0.20	23,38,66,67	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.

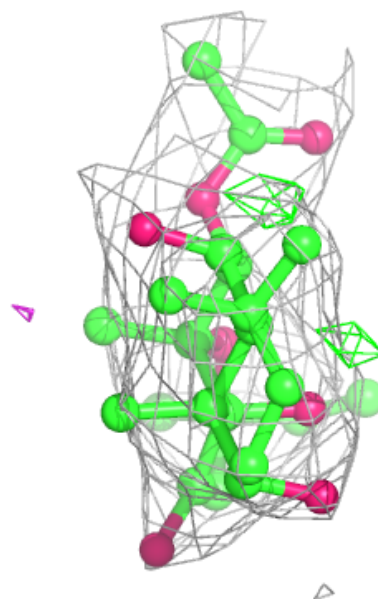
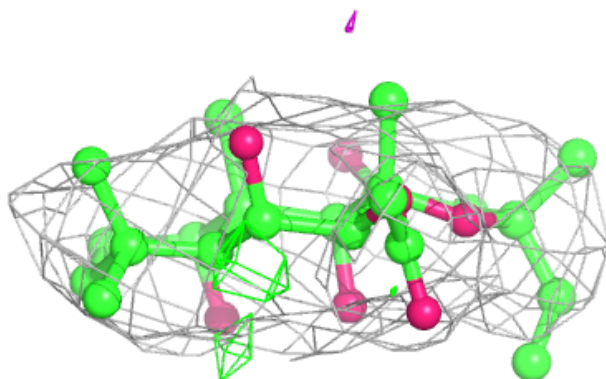
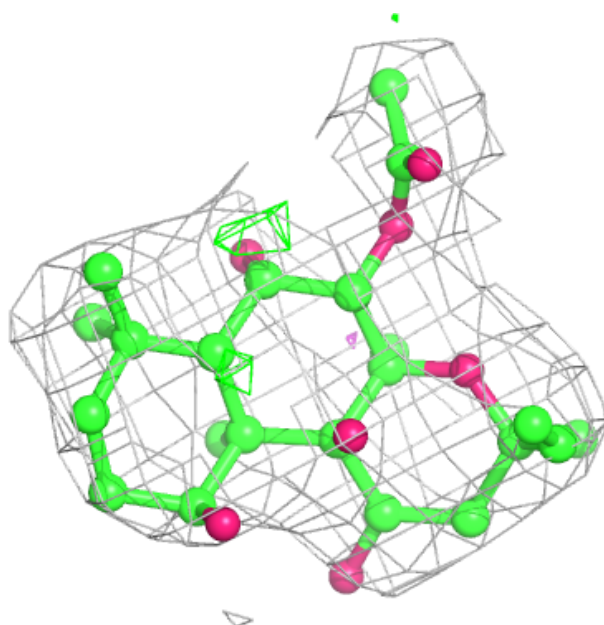
Electron density around ONM A 100:

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



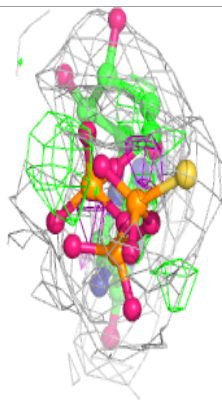
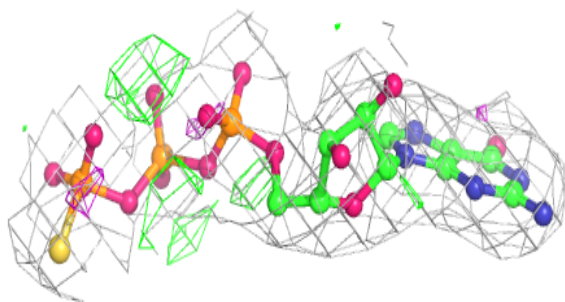
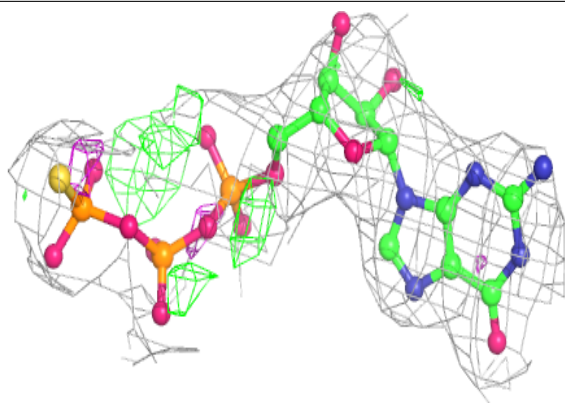
Electron density around FOK A 101:

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



Electron density around GSP C 395:

$2mF_o-DF_c$ (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.