



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:21 PM UTC

PDB ID : 1MC1 / pdb_00001mc1
Title : BETA-LACTAM SYNTHETASE WITH PRODUCT (DGPC), AMP AND PPI
Authors : Miller, M.T.; Bachmann, B.O.; Townsend, C.A.; Rosenzweig, A.C.
Deposited on : 2002-08-04
Resolution : 2.16 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

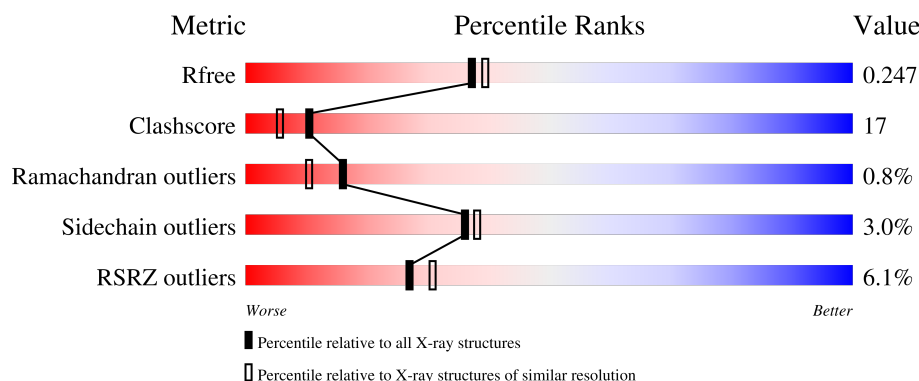
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.16 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	513	
1	B	513	

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	AMP	A	702	-	-	X	-
5	PCX	B	706	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7946 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.

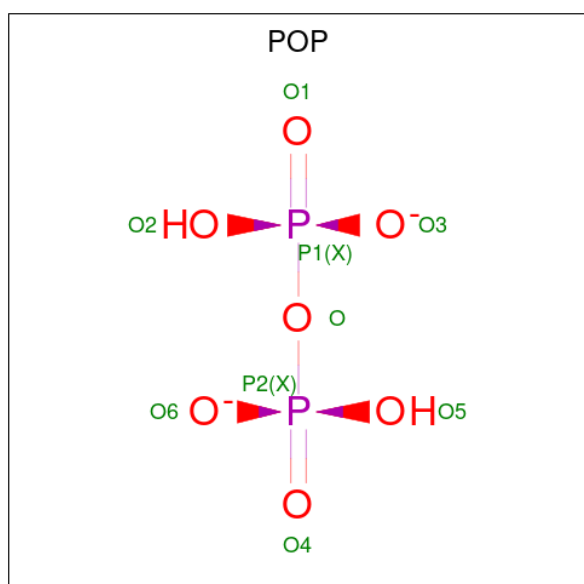
- Molecule 1 is a protein called BETA-LACTAM SYNTHETASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	491	Total	C	N	O	S	0	0	0
			3693	2305	678	702	8			
1	B	498	Total	C	N	O	S	0	0	0
			3748	2339	692	709	8			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

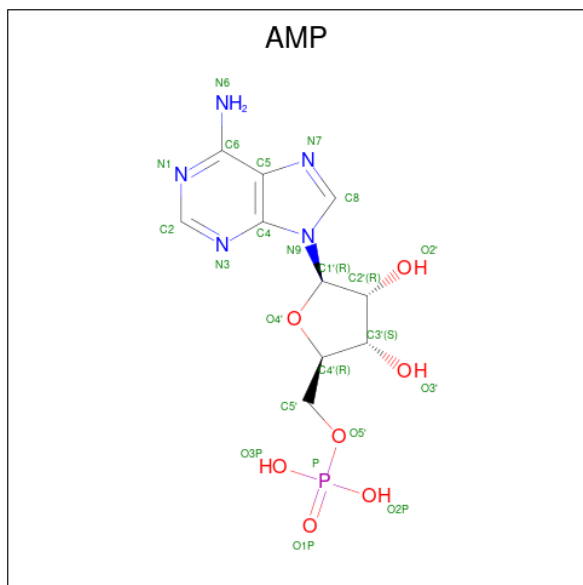
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mg	0	0
			2	2		
2	B	2	Total	Mg	0	0
			2	2		

- Molecule 3 is PYROPHOSPHATE 2- (CCD ID: POP) (formula: $\text{H}_2\text{O}_7\text{P}_2$).



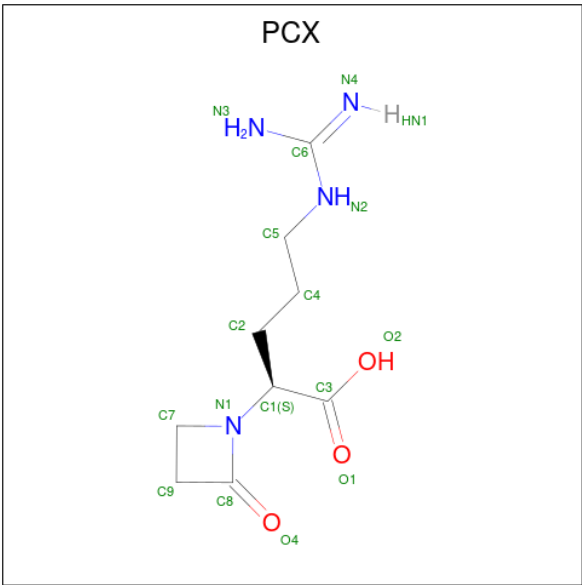
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			9	7	2		
3	B	1	Total	O	P	0	0
			9	7	2		

- Molecule 4 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: $C_{10}H_{14}N_5O_7P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
4	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 5 is DEOXYGUANIDINOPROCLAVAMINIC ACID (CCD ID: PCX) (formula: $C_9H_{16}N_4O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			16	9	4	3		
5	B	1	Total	C	N	O	0	0
			16	9	4	3		

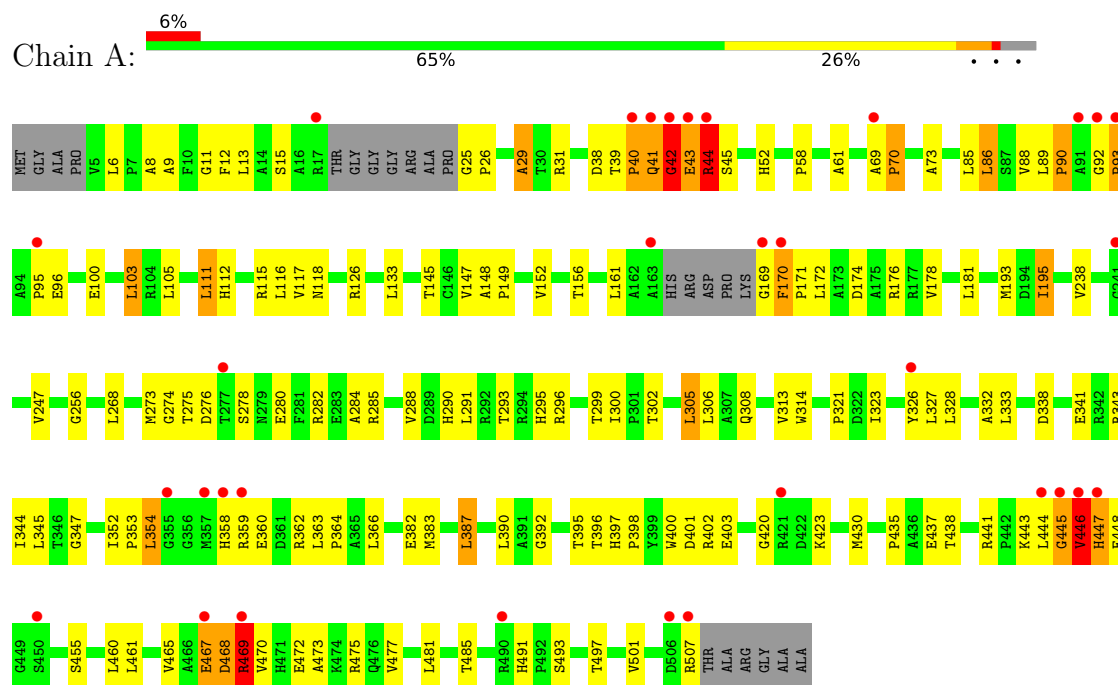
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	175	Total	O	0	0
			175	175		
6	B	230	Total	O	0	0
			230	230		

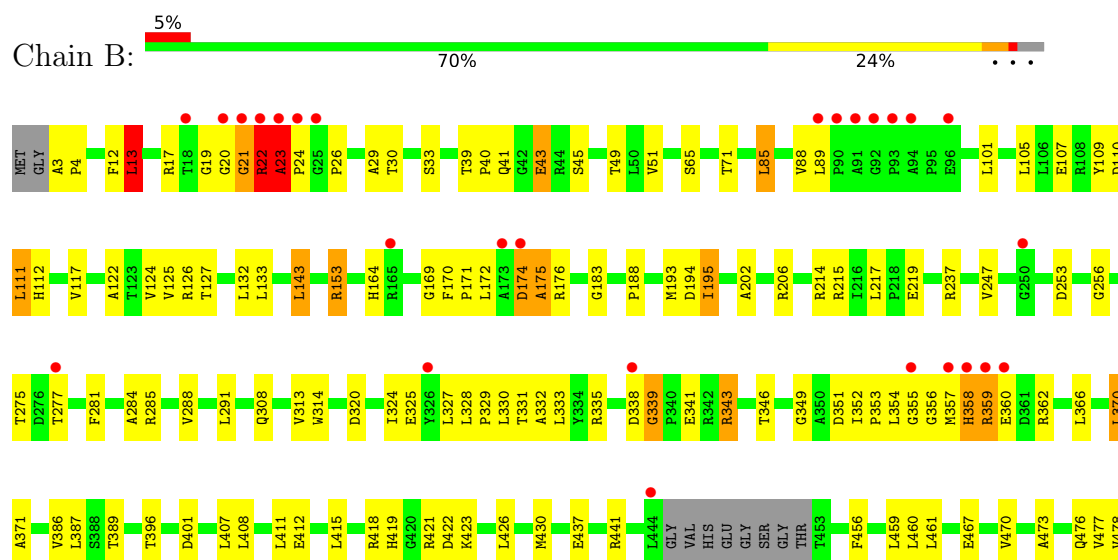
3 Residue-property plots

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: BETA-LACTAM SYNTHETASE



• Molecule 1: BETA-LACTAM SYNTHETASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.04Å 96.57Å 80.91Å 90.00° 90.54° 90.00°	Depositor
Resolution (Å)	29.10 – 2.16 29.10 – 2.16	Depositor EDS
% Data completeness (in resolution range)	94.7 (29.10-2.16) 94.8 (29.10-2.16)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.00 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.208 , 0.247 0.208 , 0.247	Depositor DCC
R_{free} test set	4350 reflections (6.86%)	wwPDB-VP
Wilson B-factor (Å ²)	21.6	Xtriage
Anisotropy	0.549	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.053 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7946	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.04% 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: PCX, MG, POP, AMP

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	0.66	12/3765 (0.3%)	1.18	41/5136 (0.8%)
1	B	0.53	3/3824 (0.1%)	1.05	24/5219 (0.5%)
All	All	0.60	15/7589 (0.2%)	1.12	65/10355 (0.6%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	9
1	B	1	5
All	All	1	14

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	468	ASP	N-CA	10.24	1.59	1.46
1	A	69	ALA	C-N	9.21	1.43	1.33
1	B	23	ALA	N-CA	7.61	1.56	1.46
1	A	170	PHE	C-N	7.11	1.42	1.33
1	A	445	GLY	C-O	7.00	1.33	1.23
1	A	467	GLU	C-O	6.85	1.32	1.24
1	B	22	ARG	C-O	6.85	1.31	1.23
1	A	169	GLY	CA-C	-6.12	1.41	1.52
1	A	169	GLY	N-CA	-5.99	1.35	1.45
1	B	175	ALA	N-CA	5.97	1.53	1.46
1	A	447	HIS	N-CA	-5.63	1.39	1.46
1	A	43	GLU	N-CA	-5.55	1.41	1.46
1	A	171	PRO	N-CA	5.50	1.53	1.47
1	A	170	PHE	N-CA	5.39	1.52	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	42	GLY	CA-C	-5.09	1.44	1.51

All (65) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	23	ALA	O-C-N	-19.13	99.32	121.32
1	A	469	ARG	O-C-N	-12.80	105.57	122.59
1	A	446	VAL	CA-C-N	12.26	140.36	120.63
1	A	446	VAL	C-N-CA	12.26	140.36	120.63
1	A	69	ALA	CA-C-N	-11.01	106.69	120.23
1	A	69	ALA	C-N-CA	-11.01	106.69	120.23
1	A	170	PHE	CA-C-N	-10.27	106.39	120.25
1	A	170	PHE	C-N-CA	-10.27	106.39	120.25
1	A	169	GLY	N-CA-C	10.19	142.85	113.30
1	A	445	GLY	O-C-N	-9.96	109.75	122.70
1	A	171	PRO	CA-N-CD	-9.58	98.59	112.00
1	B	22	ARG	CA-C-O	-9.20	109.79	121.11
1	A	70	PRO	CA-N-CD	-8.55	100.03	112.00
1	A	467	GLU	CA-C-O	-8.17	108.83	120.51
1	A	117	VAL	N-CA-C	8.12	119.61	108.89
1	B	22	ARG	N-CA-C	8.00	123.58	107.41
1	B	174	ASP	CA-CB-CG	7.52	120.12	112.60
1	B	195	ILE	N-CA-C	7.50	118.66	108.17
1	B	358	HIS	CA-C-N	7.42	132.66	122.34
1	B	358	HIS	C-N-CA	7.42	132.66	122.34
1	B	21	GLY	O-C-N	-7.42	113.06	122.70
1	A	170	PHE	C-N-CD	7.32	155.00	125.00
1	A	467	GLU	N-CA-C	7.19	126.11	110.80
1	A	447	HIS	N-CA-C	7.17	121.66	112.34
1	A	467	GLU	O-C-N	-7.10	113.14	122.59
1	B	401	ASP	N-CA-C	-7.09	100.50	110.50
1	A	69	ALA	C-N-CD	6.89	153.26	125.00
1	A	169	GLY	O-C-N	-6.80	112.12	123.00
1	A	29	ALA	N-CA-C	-6.72	103.76	112.23
1	B	22	ARG	O-C-N	-6.72	114.78	122.84
1	A	44	ARG	N-CA-C	-6.69	105.88	112.97
1	A	401	ASP	N-CA-C	-6.63	100.62	110.23
1	A	41	GLN	CA-C-N	6.34	133.84	121.41
1	A	41	GLN	C-N-CA	6.34	133.84	121.41
1	B	117	VAL	N-CA-C	6.27	117.82	108.54
1	A	70	PRO	CB-CA-C	6.25	118.37	111.56
1	A	170	PHE	N-CA-CB	-6.23	102.61	109.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	390	LEU	N-CA-C	-6.21	104.97	112.54
1	A	446	VAL	CA-C-O	-6.19	113.90	120.64
1	A	42	GLY	CA-C-N	-6.15	114.15	121.90
1	A	42	GLY	C-N-CA	-6.15	114.15	121.90
1	B	359	ARG	O-C-N	-6.05	115.32	122.58
1	B	29	ALA	N-CA-C	-6.01	105.03	112.90
1	B	20	GLY	CA-C-N	-6.00	109.65	121.41
1	B	20	GLY	C-N-CA	-6.00	109.65	121.41
1	A	278	SER	N-CA-C	5.99	118.18	109.07
1	B	21	GLY	CA-C-N	-5.87	114.02	123.23
1	B	21	GLY	C-N-CA	-5.87	114.02	123.23
1	A	468	ASP	N-CA-CB	-5.76	100.76	110.49
1	A	195	ILE	N-CA-C	5.72	116.12	108.11
1	A	70	PRO	N-CA-CB	5.67	107.83	103.30
1	A	343	ARG	N-CA-C	-5.61	98.61	108.20
1	A	170	PHE	N-CA-C	-5.54	103.66	110.31
1	A	467	GLU	CA-C-N	5.52	132.08	121.54
1	A	467	GLU	C-N-CA	5.52	132.08	121.54
1	B	39	THR	CA-C-N	5.49	125.16	119.56
1	B	39	THR	C-N-CA	5.49	125.16	119.56
1	B	22	ARG	CA-C-N	5.43	135.05	121.80
1	B	22	ARG	C-N-CA	5.43	135.05	121.80
1	A	25	GLY	CA-C-N	5.33	125.25	119.76
1	A	25	GLY	C-N-CA	5.33	125.25	119.76
1	A	43	GLU	O-C-N	5.24	126.92	121.79
1	B	13	LEU	N-CA-C	-5.17	98.85	107.80
1	B	22	ARG	CD-NE-CZ	5.15	131.61	124.40
1	B	343	ARG	N-CA-C	-5.01	99.63	108.20

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	22	ARG	CA

All (14) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	174	ASP	Sidechain
1	A	42	GLY	Mainchain,Peptide
1	A	44	ARG	Mainchain
1	A	445	GLY	Mainchain
1	A	446	VAL	Mainchain

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Mol	Chain	Res	Type	Group
1	A	467	GLU	Mainchain
1	A	469	ARG	Mainchain
1	A	70	PRO	Mainchain
1	B	21	GLY	Mainchain,Peptide
1	B	22	ARG	Mainchain,Peptide
1	B	23	ALA	Mainchain

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	3693	0	3681	141	2
1	B	3748	0	3743	110	1
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	9	0	0	1	0
3	B	9	0	0	1	0
4	A	23	0	12	14	0
4	B	23	0	12	3	0
5	A	16	0	14	5	0
5	B	16	0	14	7	0
6	A	175	0	0	3	0
6	B	230	0	0	7	0
All	All	7946	0	7476	253	2

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 17.

All (253) 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)
4:A:702:AMP:O3P	5:A:703:PCX:H9C1	1.14	1.32
1:A:149:PRO:CD	1:A:170:PHE:CZ	2.20	1.23
1:A:149:PRO:CD	1:A:170:PHE:CE1	2.30	1.15
1:A:149:PRO:HD3	1:A:170:PHE:CE1	1.80	1.15
1:A:326:TYR:CE2	1:A:447:HIS:CE1	2.44	1.05

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:702:AMP:O3P	5:A:703:PCX:C9	2.06	1.03
1:A:149:PRO:HD3	1:A:170:PHE:CZ	1.86	1.02
1:A:446:VAL:HG21	4:A:702:AMP:C8	1.97	1.00
1:A:446:VAL:CG2	4:A:702:AMP:C8	2.46	0.99
1:B:357:MET:HE1	5:B:706:PCX:O4	1.64	0.97
1:A:149:PRO:HD2	1:A:170:PHE:CE1	1.97	0.96
1:B:13:LEU:HD12	1:B:26:PRO:HB3	1.48	0.94
1:B:357:MET:CE	5:B:706:PCX:O4	2.15	0.94
1:A:149:PRO:HD2	1:A:170:PHE:CZ	2.02	0.94
1:A:41:GLN:O	1:A:43:GLU:N	2.01	0.93
5:B:706:PCX:H7C2	4:B:705:AMP:O3P	1.70	0.91
1:B:357:MET:HE1	5:B:706:PCX:C8	2.04	0.88
1:A:326:TYR:CE2	1:A:447:HIS:ND1	2.42	0.87
1:A:13:LEU:HD23	1:A:26:PRO:HG3	1.58	0.85
1:A:326:TYR:HE2	1:A:447:HIS:CE1	1.92	0.83
1:A:444:LEU:HD12	1:A:448:GLU:HB2	1.60	0.82
1:B:459:LEU:HD11	1:B:501:VAL:HG11	1.62	0.81
1:A:193:MET:HE2	1:A:195:ILE:HD11	1.62	0.81
1:A:446:VAL:HG22	4:A:702:AMP:N7	1.96	0.81
1:A:446:VAL:HG22	4:A:702:AMP:C8	2.16	0.80
1:A:444:LEU:CD1	1:A:448:GLU:HB2	2.13	0.79
1:A:446:VAL:CG2	4:A:702:AMP:H8	1.95	0.77
1:B:174:ASP:O	1:B:188:PRO:HG3	1.83	0.77
1:A:112:HIS:HD2	1:A:115:ARG:HE	1.31	0.76
1:B:17:ARG:HG2	1:B:19:GLY:H	1.52	0.74
1:A:149:PRO:CG	1:A:170:PHE:CZ	2.72	0.72
1:A:273:MET:HB2	1:A:333:LEU:HD13	1.72	0.71
1:B:22:ARG:CG	1:B:23:ALA:HB3	2.18	0.71
1:A:352:ILE:HB	1:A:353:PRO:HD3	1.73	0.70
1:A:300:ILE:HG13	1:A:300:ILE:O	1.91	0.70
1:B:22:ARG:HG2	1:B:23:ALA:HB3	1.72	0.69
1:B:328:LEU:HB2	1:B:329:PRO:HD3	1.73	0.69
1:B:174:ASP:O	1:B:188:PRO:CG	2.41	0.69
1:A:326:TYR:CD2	1:A:447:HIS:ND1	2.61	0.68
1:A:321:PRO:HG2	1:A:475:ARG:HG2	1.75	0.68
1:B:327:LEU:O	1:B:331:THR:HG23	1.93	0.68
1:B:153:ARG:HG2	1:B:153:ARG:HH21	1.61	0.66
1:A:446:VAL:HG21	4:A:702:AMP:H8	1.53	0.66
1:B:459:LEU:CD1	1:B:501:VAL:HG11	2.26	0.65
1:A:176:ARG:HG2	1:A:176:ARG:HH11	1.62	0.65
1:A:41:GLN:HG2	1:A:45:SER:HB2	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:172:LEU:HD13	1:B:193:MET:HE1	1.78	0.65
1:A:446:VAL:CG2	4:A:702:AMP:N7	2.59	0.64
1:A:44:ARG:O	1:A:126:ARG:CZ	2.46	0.64
1:A:327:LEU:HD22	1:A:387:LEU:HD21	1.79	0.63
1:A:444:LEU:HD12	1:A:448:GLU:OE2	1.99	0.62
1:A:280:GLU:CD	1:A:444:LEU:HD23	2.23	0.62
1:A:41:GLN:C	1:A:43:GLU:H	1.96	0.62
1:B:437:GLU:HG2	1:B:441:ARG:NH2	2.15	0.62
1:A:280:GLU:OE2	1:A:444:LEU:HD23	2.00	0.61
1:A:300:ILE:HD11	1:A:305:LEU:HG	1.83	0.61
1:B:360:GLU:CG	1:B:362:ARG:HG3	2.30	0.61
1:A:148:ALA:HA	1:A:170:PHE:CD1	2.35	0.61
1:A:443:LYS:HD3	1:A:444:LEU:N	2.15	0.60
1:A:73:ALA:HB2	1:A:103:LEU:HD13	1.83	0.60
1:B:41:GLN:HG2	1:B:45:SER:OG	2.01	0.60
1:A:360:GLU:HB2	6:A:820:HOH:O	2.01	0.60
1:B:132:LEU:HD23	1:B:194:ASP:HA	1.83	0.60
1:B:170:PHE:HB3	1:B:171:PRO:HD2	1.84	0.60
1:A:92:GLY:N	1:A:93:PRO:HD2	2.17	0.60
1:B:153:ARG:HH22	1:B:164:HIS:CD2	2.19	0.60
1:B:308:GLN:HG3	1:B:332:ALA:HB2	1.83	0.59
5:B:706:PCX:C7	4:B:705:AMP:O3P	2.47	0.59
1:B:3:ALA:HB3	1:B:4:PRO:HD3	1.85	0.59
1:B:386:VAL:HA	1:B:389:THR:OG1	2.03	0.59
1:A:314:TRP:CD1	1:A:485:THR:HB	2.38	0.58
1:B:359:ARG:HD3	6:B:927:HOH:O	2.03	0.58
1:B:85:LEU:HG	1:B:105:LEU:HD11	1.84	0.58
1:B:358:HIS:CE1	1:B:360:GLU:HB3	2.39	0.57
1:B:215:ARG:HH11	1:B:215:ARG:HG3	1.68	0.57
1:B:491:HIS:O	1:B:494:GLU:HG2	2.04	0.57
1:A:296:ARG:HH12	1:A:338:ASP:H	1.52	0.57
1:B:308:GLN:HE22	1:B:335:ARG:HH21	1.50	0.57
1:A:345:LEU:HD13	1:A:396:THR:HG23	1.84	0.57
1:B:85:LEU:O	1:B:88:VAL:HG22	2.05	0.57
1:B:360:GLU:HG3	1:B:362:ARG:HG3	1.87	0.57
1:A:282:ARG:HD3	1:A:282:ARG:C	2.29	0.57
1:B:355:GLY:HA3	6:B:849:HOH:O	2.04	0.57
1:B:314:TRP:CD1	1:B:485:THR:HB	2.40	0.57
1:A:176:ARG:HG2	1:A:176:ARG:NH1	2.20	0.56
1:A:354:LEU:HB3	1:A:423:LYS:HG2	1.88	0.56
1:B:169:GLY:HA3	1:B:183:GLY:O	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:418:ARG:HG2	1:B:419:HIS:CD2	2.40	0.56
1:B:360:GLU:HG3	1:B:362:ARG:H	1.70	0.56
1:B:65:SER:HB3	1:B:71:THR:HB	1.87	0.56
1:B:71:THR:OG1	1:B:127:THR:HG23	2.06	0.56
1:A:29:ALA:HA	1:A:314:TRP:CH2	2.41	0.55
1:A:306:LEU:HD21	1:A:455:SER:HB3	1.87	0.55
1:A:284:ALA:O	1:A:288:VAL:HG13	2.07	0.55
1:B:461:LEU:HD13	1:B:470:VAL:HG21	1.88	0.55
1:A:52:HIS:CE1	1:A:58:PRO:HG3	2.43	0.54
1:A:313:VAL:HG21	1:A:481:LEU:HB3	1.89	0.54
1:B:122:ALA:HB1	1:B:143:LEU:HG	1.89	0.54
1:B:351:ASP:OD2	1:B:356:GLY:HA2	2.07	0.54
1:A:112:HIS:CD2	1:A:115:ARG:HE	2.21	0.54
1:A:437:GLU:HG2	1:A:441:ARG:NH1	2.23	0.53
1:A:193:MET:CE	1:A:195:ILE:HD11	2.36	0.53
1:B:172:LEU:HB2	1:B:175:ALA:HB3	1.91	0.53
1:B:357:MET:CE	5:B:706:PCX:C8	2.77	0.53
1:B:370:LEU:HD13	1:B:370:LEU:C	2.34	0.53
1:B:437:GLU:HG2	1:B:441:ARG:HH22	1.72	0.53
1:B:459:LEU:HD12	1:B:460:LEU:HD22	1.91	0.52
1:A:472:GLU:OE2	1:A:475:ARG:HD2	2.10	0.52
1:A:52:HIS:NE2	1:A:58:PRO:HG3	2.25	0.52
1:A:149:PRO:HG2	1:A:170:PHE:CZ	2.45	0.52
1:A:285:ARG:O	1:A:288:VAL:HG22	2.10	0.52
1:A:507:ARG:HD3	1:A:507:ARG:N	2.25	0.51
1:A:148:ALA:HA	1:A:170:PHE:CE1	2.46	0.51
1:B:89:LEU:HD11	1:B:105:LEU:HD23	1.93	0.51
1:A:344:ILE:HB	1:A:395:THR:HG22	1.93	0.51
1:A:497:THR:O	1:A:501:VAL:HG23	2.10	0.51
1:B:153:ARG:HG2	1:B:153:ARG:NH2	2.26	0.51
1:B:247:VAL:HG23	4:B:705:AMP:O2'	2.10	0.51
1:B:111:LEU:HD12	1:B:112:HIS:N	2.26	0.51
1:B:125:VAL:CG1	1:B:132:LEU:HB2	2.41	0.51
1:B:22:ARG:HG3	1:B:23:ALA:HB3	1.91	0.50
1:B:253:ASP:HB2	3:B:704:POP:P1	2.51	0.50
1:B:327:LEU:HD13	1:B:387:LEU:HD21	1.92	0.50
1:A:41:GLN:C	1:A:43:GLU:N	2.62	0.50
1:A:435:PRO:HG2	1:A:438:THR:OG1	2.12	0.50
1:B:275:THR:OG1	1:B:277:THR:HG22	2.12	0.50
1:B:366:LEU:C	1:B:366:LEU:HD23	2.37	0.50
1:A:300:ILE:HD13	1:A:332:ALA:HB1	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:PRO:HG2	1:A:475:ARG:CG	2.40	0.49
1:B:256:GLY:HA2	1:B:430:MET:HE2	1.94	0.49
1:A:43:GLU:C	1:A:45:SER:H	2.19	0.49
1:A:358:HIS:HB3	6:A:776:HOH:O	2.11	0.49
1:B:85:LEU:O	1:B:89:LEU:HD13	2.13	0.49
1:A:446:VAL:CG2	4:A:702:AMP:H5'1	2.43	0.49
1:A:247:VAL:HG23	1:A:247:VAL:O	2.11	0.48
1:A:111:LEU:HD12	1:A:112:HIS:N	2.28	0.48
1:A:321:PRO:HG3	1:A:475:ARG:HA	1.95	0.48
1:B:176:ARG:NH1	6:B:853:HOH:O	2.43	0.48
1:A:85:LEU:O	1:A:88:VAL:HG22	2.14	0.48
1:A:147:VAL:HB	1:A:172:LEU:HD21	1.95	0.48
1:A:31:ARG:NH2	1:A:392:GLY:HA3	2.29	0.48
1:B:349:GLY:O	1:B:353:PRO:HD2	2.14	0.48
1:A:39:THR:O	1:A:42:GLY:HA2	2.13	0.48
1:B:421:ARG:HD2	6:B:761:HOH:O	2.14	0.48
1:A:275:THR:HG21	1:A:305:LEU:HD12	1.96	0.47
1:A:268:LEU:O	1:A:293:THR:HB	2.14	0.47
1:A:285:ARG:HA	1:A:288:VAL:HG22	1.96	0.47
1:A:323:ILE:O	1:A:327:LEU:HG	2.14	0.47
1:B:195:ILE:HG12	1:B:202:ALA:HB2	1.97	0.47
1:B:30:THR:HG23	1:B:389:THR:HB	1.97	0.47
1:B:215:ARG:HG3	1:B:215:ARG:NH1	2.30	0.47
1:A:302:THR:O	1:A:306:LEU:HD23	2.15	0.47
1:A:444:LEU:CD1	1:A:448:GLU:CB	2.89	0.47
1:A:491:HIS:HD2	1:A:493:SER:OG	1.96	0.47
1:B:281:PHE:O	1:B:285:ARG:HG3	2.15	0.47
1:A:314:TRP:HD1	1:A:485:THR:HB	1.80	0.47
1:A:444:LEU:HD11	1:A:448:GLU:HB2	1.91	0.46
1:B:237:ARG:O	1:B:343:ARG:HD3	2.15	0.46
1:B:473:ALA:O	1:B:477:VAL:HG23	2.15	0.46
1:A:92:GLY:H	1:A:93:PRO:HD2	1.80	0.46
1:A:256:GLY:HA2	1:A:430:MET:HE2	1.97	0.46
1:B:237:ARG:HD3	1:B:396:THR:HG23	1.97	0.46
1:A:86:LEU:HD11	1:A:95:PRO:HD2	1.98	0.46
1:A:362:ARG:NH2	1:B:107:GLU:O	2.49	0.46
1:A:469:ARG:O	1:A:470:VAL:C	2.59	0.45
1:B:22:ARG:HG2	1:B:23:ALA:O	2.16	0.45
1:B:40:PRO:HG2	1:B:41:GLN:NE2	2.31	0.45
1:A:15:SER:HA	1:A:152:VAL:O	2.16	0.45
1:A:133:LEU:HD12	1:A:193:MET:HE2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:422:ASP:OD2	1:B:423:LYS:HE2	2.16	0.45
1:B:247:VAL:HG23	1:B:247:VAL:O	2.17	0.45
1:A:178:VAL:HB	1:A:181:LEU:HD12	1.98	0.45
1:B:26:PRO:HG3	1:B:49:THR:HG21	1.99	0.45
1:A:288:VAL:HG11	1:A:295:HIS:CG	2.52	0.45
1:B:370:LEU:HD13	1:B:371:ALA:N	2.32	0.45
1:A:383:MET:HE1	1:A:400:TRP:CB	2.47	0.45
1:A:73:ALA:HB2	1:A:103:LEU:CD1	2.45	0.45
1:A:285:ARG:HD3	1:A:288:VAL:CG2	2.46	0.45
1:A:363:LEU:N	1:A:364:PRO:CD	2.80	0.45
1:B:338:ASP:O	1:B:339:GLY:O	2.35	0.45
1:A:460:LEU:HD13	1:A:473:ALA:HB3	1.99	0.44
1:B:125:VAL:HG13	1:B:132:LEU:HB2	2.00	0.44
1:A:274:GLY:O	1:A:299:THR:HA	2.17	0.44
1:A:6:LEU:HD11	1:A:341:GLU:HG3	1.98	0.44
1:B:71:THR:HA	1:B:126:ARG:O	2.18	0.44
1:A:85:LEU:HD13	1:A:105:LEU:HD11	2.00	0.44
1:B:89:LEU:HD11	1:B:105:LEU:CD2	2.48	0.44
1:A:382:GLU:OE1	5:A:703:PCX:H5C2	2.17	0.44
1:B:284:ALA:O	1:B:288:VAL:HG23	2.18	0.44
1:B:333:LEU:C	1:B:333:LEU:HD23	2.42	0.44
1:B:124:VAL:HG22	1:B:133:LEU:CD2	2.48	0.43
1:A:359:ARG:HH21	1:A:359:ARG:HG3	1.83	0.43
1:A:8:ALA:HB3	1:A:61:ALA:HB2	2.00	0.43
1:B:33:SER:O	1:B:51:VAL:HA	2.17	0.43
1:B:352:ILE:HB	1:B:353:PRO:CD	2.49	0.43
1:A:366:LEU:C	1:A:366:LEU:HD23	2.43	0.43
1:B:308:GLN:NE2	1:B:335:ARG:HH21	2.14	0.43
1:A:96:GLU:HG2	1:A:100:GLU:OE2	2.19	0.43
1:B:461:LEU:HD12	1:B:461:LEU:HA	1.95	0.43
1:A:116:LEU:HD23	1:B:214:ARG:HB2	2.01	0.43
1:B:407:LEU:O	1:B:411:LEU:HG	2.19	0.43
1:A:145:THR:HG21	1:A:193:MET:HE1	2.00	0.42
1:A:148:ALA:O	1:A:149:PRO:C	2.62	0.42
1:A:461:LEU:HD23	1:A:465:VAL:O	2.20	0.42
1:A:444:LEU:HD12	1:A:448:GLU:CB	2.41	0.42
1:B:41:GLN:HE21	1:B:45:SER:HB3	1.85	0.42
1:B:110:ASP:C	1:B:112:HIS:N	2.78	0.42
1:B:441:ARG:HH11	1:B:441:ARG:HG3	1.84	0.42
1:A:11:GLY:HA2	1:A:156:THR:HG21	2.01	0.42
1:B:174:ASP:O	1:B:188:PRO:HG2	2.16	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:217:LEU:HD11	6:B:875:HOH:O	2.19	0.42
1:A:12:PHE:CD1	1:A:12:PHE:C	2.98	0.42
1:A:382:GLU:HG2	6:A:751:HOH:O	2.19	0.42
1:B:12:PHE:CD1	1:B:12:PHE:C	2.97	0.42
1:B:110:ASP:O	1:B:112:HIS:N	2.53	0.42
1:A:443:LYS:HD3	1:A:443:LYS:C	2.45	0.42
1:A:460:LEU:HD22	1:A:465:VAL:HG21	2.01	0.42
1:A:9:ALA:HA	1:A:52:HIS:HB3	2.01	0.42
1:A:290:HIS:CG	1:A:435:PRO:HD3	2.54	0.42
1:A:291:LEU:N	1:A:291:LEU:HD12	2.35	0.42
4:A:702:AMP:H8	4:A:702:AMP:H5'1	1.85	0.42
1:A:149:PRO:HD3	1:A:170:PHE:CD1	2.46	0.42
1:A:397:HIS:HA	1:A:398:PRO:HD2	1.95	0.42
1:A:468:ASP:C	1:A:469:ARG:HG3	2.44	0.42
1:B:320:ASP:O	1:B:324:ILE:HG12	2.19	0.42
1:B:357:MET:N	1:B:357:MET:SD	2.93	0.41
1:A:247:VAL:HB	4:A:702:AMP:N3	2.35	0.41
1:A:473:ALA:O	1:A:477:VAL:HG23	2.21	0.41
1:B:354:LEU:HD21	1:B:408:LEU:HD22	2.03	0.41
1:B:24:PRO:HG2	1:B:164:HIS:CE1	2.55	0.41
1:B:206:ARG:NH2	6:B:802:HOH:O	2.53	0.41
1:A:446:VAL:HG21	4:A:702:AMP:H5'1	2.03	0.41
1:B:313:VAL:HG22	1:B:478:VAL:HG13	2.02	0.41
1:A:149:PRO:HD2	1:A:170:PHE:HE1	1.72	0.41
1:A:305:LEU:CD2	1:A:328:LEU:HB3	2.50	0.41
1:B:459:LEU:C	1:B:459:LEU:HD13	2.46	0.41
1:A:103:LEU:HD12	1:A:103:LEU:HA	1.81	0.41
1:A:238:VAL:O	1:A:238:VAL:HG23	2.21	0.41
1:B:219:GLU:CD	1:B:418:ARG:HH22	2.28	0.41
1:A:352:ILE:HD11	5:A:703:PCX:H5C1	2.03	0.41
1:B:43:GLU:H	1:B:43:GLU:HG2	1.54	0.41
1:B:325:GLU:OE1	1:B:456:PHE:HB3	2.21	0.41
1:B:476:GLN:HB3	6:B:835:HOH:O	2.20	0.41
1:A:31:ARG:HG2	1:A:31:ARG:HH11	1.85	0.40
1:A:38:ASP:C	1:A:40:PRO:HD3	2.46	0.40
1:A:352:ILE:HG13	5:A:703:PCX:O1	2.21	0.40
1:B:422:ASP:O	1:B:423:LYS:C	2.63	0.40
1:A:364:PRO:HG3	1:B:109:TYR:C	2.45	0.40
1:A:89:LEU:HA	1:A:90:PRO:HD3	1.84	0.40
1:A:118:ASN:HD21	1:A:402:ARG:H	1.68	0.40
1:A:347:GLY:C	4:A:702:AMP:O3'	2.65	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:443:LYS:NZ	3:A:701:POP:O3	2.52	0.40
1:B:357:MET:HE3	5:B:706:PCX:O4	2.12	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:ARG:NE	1:A:276:ASP:OD1[1_455]	2.10	0.10
1:A:420:GLY:O	1:B:359:ARG:NE[1_655]	2.10	0.10

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	485/513 (94%)	461 (95%)	18 (4%)	6 (1%)	10	5
1	B	494/513 (96%)	474 (96%)	18 (4%)	2 (0%)	30	26
All	All	979/1026 (95%)	935 (96%)	36 (4%)	8 (1%)	16	10

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	42	GLY
1	A	90	PRO
1	A	111	LEU
1	B	111	LEU
1	B	339	GLY
1	A	469	ARG
1	A	40	PRO
1	A	93	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	383/395 (97%)	375 (98%)	8 (2%)	47	52
1	B	388/395 (98%)	373 (96%)	15 (4%)	28	28
All	All	771/790 (98%)	748 (97%)	23 (3%)	36	38

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	86	LEU
1	A	103	LEU
1	A	161	LEU
1	A	305	LEU
1	A	308	GLN
1	A	354	LEU
1	A	387	LEU
1	A	403	GLU
1	B	13	LEU
1	B	43	GLU
1	B	85	LEU
1	B	101	LEU
1	B	143	LEU
1	B	153	ARG
1	B	291	LEU
1	B	330	LEU
1	B	341	GLU
1	B	346	THR
1	B	370	LEU
1	B	412	GLU
1	B	415	LEU
1	B	426	LEU
1	B	467	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	112	HIS
1	A	118	ASN
1	A	186	GLN
1	A	236	GLN
1	A	471	HIS
1	A	491	HIS
1	B	41	GLN
1	B	236	GLN
1	B	308	GLN
1	B	358	HIS
1	B	381	ASN
1	B	440	ASN
1	B	471	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 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	$\# Z > 2$
4	AMP	B	705	2	25,25,25	2.06	8 (32%)	37,38,38	1.41	7 (18%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	AMP	A	702	2	25,25,25	2.00	6 (24%)	37,38,38	1.49	7 (18%)
5	PCX	B	706	-	16,16,16	7.49	6 (37%)	14,21,21	7.27	5 (35%)
3	POP	A	701	2	6,8,8	1.15	0	12,13,13	1.09	0
5	PCX	A	703	-	16,16,16	10.33	7 (43%)	14,21,21	4.11	5 (35%)
3	POP	B	704	2	6,8,8	1.02	0	12,13,13	0.82	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	AMP	B	705	2	-	1/10/26/26	0/3/3/3
4	AMP	A	702	2	-	1/10/26/26	0/3/3/3
5	PCX	B	706	-	-	7/14/24/24	0/1/1/1
3	POP	A	701	2	-	0/6/6/6	-
5	PCX	A	703	-	-	9/14/24/24	0/1/1/1
3	POP	B	704	2	-	0/6/6/6	-

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	703	PCX	C1-N1	-36.87	1.02	1.46
5	B	706	PCX	C1-N1	-23.17	1.18	1.46
5	B	706	PCX	C9-C8	-15.35	1.41	1.52
5	A	703	PCX	C9-C8	-14.49	1.41	1.52
5	A	703	PCX	C7-C9	-8.20	1.42	1.54
5	B	706	PCX	C7-C9	-7.78	1.42	1.54
4	B	705	AMP	C3'-C4'	6.07	1.68	1.53
4	A	702	AMP	C3'-C4'	6.02	1.68	1.53
5	A	703	PCX	C7-N1	-4.59	1.40	1.48
5	B	706	PCX	C7-N1	-4.45	1.40	1.48
5	A	703	PCX	O4-C8	-4.44	1.13	1.23
5	B	706	PCX	C8-N1	4.42	1.43	1.36
5	A	703	PCX	O1-C3	4.14	1.34	1.22
5	B	706	PCX	O1-C3	3.99	1.33	1.22
4	B	705	AMP	O3'-C3'	3.29	1.51	1.43
4	A	702	AMP	O3'-C3'	3.07	1.50	1.43
4	B	705	AMP	C2'-C3'	2.77	1.60	1.53
4	B	705	AMP	C2'-C1'	2.73	1.62	1.53
4	A	702	AMP	C2'-C1'	2.64	1.61	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	705	AMP	C8-N9	2.63	1.42	1.37
4	A	702	AMP	C5-C4	2.59	1.43	1.39
4	A	702	AMP	C2'-C3'	2.55	1.60	1.53
4	B	705	AMP	C5-C4	2.54	1.43	1.39
5	A	703	PCX	C8-N1	2.52	1.40	1.36
4	B	705	AMP	C8-N7	2.11	1.35	1.31
4	A	702	AMP	C8-N9	2.08	1.41	1.37
4	B	705	AMP	C1'-N9	2.03	1.51	1.46

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	706	PCX	O4-C8-N1	24.99	145.44	131.14
5	A	703	PCX	O4-C8-N1	13.49	138.86	131.14
5	B	706	PCX	O4-C8-C9	-8.78	122.61	135.39
5	A	703	PCX	O4-C8-C9	-4.22	129.25	135.39
5	A	703	PCX	O1-C3-C1	-3.77	110.87	122.44
5	B	706	PCX	O1-C3-C1	-3.65	111.25	122.44
5	A	703	PCX	O2-C3-O1	3.26	131.49	124.08
4	A	702	AMP	N6-C6-N1	3.21	125.53	118.38
4	A	702	AMP	N9-C8-N7	3.12	118.36	113.94
4	A	702	AMP	C5-N7-C8	-3.08	98.60	103.45
5	B	706	PCX	O2-C3-O1	3.07	131.05	124.08
4	B	705	AMP	N6-C6-N1	3.04	125.14	118.38
4	A	702	AMP	C4-C5-N7	3.02	114.04	110.58
4	B	705	AMP	C4-C5-N7	2.85	113.84	110.58
4	B	705	AMP	C5-N7-C8	-2.70	99.21	103.45
4	A	702	AMP	C2-N1-C6	2.65	123.08	118.73
4	B	705	AMP	C2-N1-C6	2.60	123.00	118.73
4	B	705	AMP	N9-C8-N7	2.53	117.53	113.94
5	B	706	PCX	C7-C9-C8	2.46	87.87	85.40
4	A	702	AMP	N3-C2-N1	-2.40	124.95	128.58
4	B	705	AMP	N3-C2-N1	-2.25	125.17	128.58
5	A	703	PCX	C7-C9-C8	2.22	87.62	85.40
4	A	702	AMP	C4-N9-C8	-2.14	103.49	105.74
4	B	705	AMP	C3'-C2'-C1'	2.04	105.32	101.46

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	703	PCX	C2-C1-N1-C7

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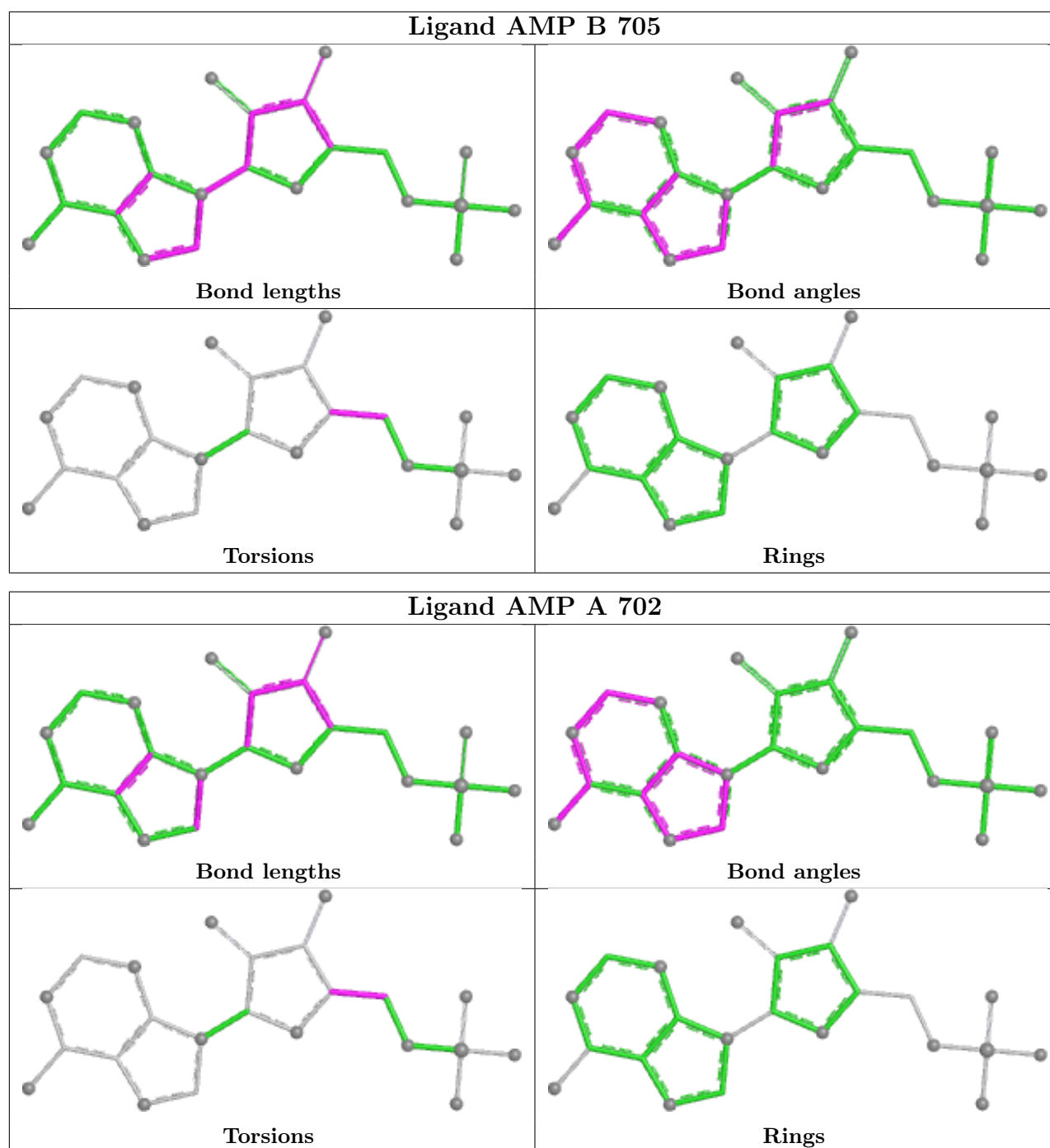
Mol	Chain	Res	Type	Atoms
5	A	703	PCX	N1-C1-C2-C4
5	A	703	PCX	C3-C1-C2-C4
5	A	703	PCX	C2-C1-C3-O1
5	B	706	PCX	C2-C1-N1-C7
5	B	706	PCX	N1-C1-C2-C4
5	B	706	PCX	C3-C1-C2-C4
5	B	706	PCX	C2-C1-C3-O1
5	B	706	PCX	C2-C4-C5-N2
5	A	703	PCX	C1-C2-C4-C5
5	A	703	PCX	C2-C1-C3-O2
5	B	706	PCX	C2-C1-C3-O2
5	A	703	PCX	C2-C4-C5-N2
5	A	703	PCX	N1-C1-C3-O1
5	A	703	PCX	N1-C1-C3-O2
5	B	706	PCX	C1-C2-C4-C5
4	A	702	AMP	O4'-C4'-C5'-O5'
4	B	705	AMP	O4'-C4'-C5'-O5'

There are no ring outliers.

6 monomers are involved in 27 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	705	AMP	3	0
4	A	702	AMP	14	0
5	B	706	PCX	7	0
3	A	701	POP	1	0
5	A	703	PCX	5	0
3	B	704	POP	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.



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	491/513 (95%)	0.35	32 (6%) 25 28	10, 25, 46, 59	0
1	B	498/513 (97%)	0.20	28 (5%) 30 34	13, 23, 42, 60	0
All	All	989/1026 (96%)	0.27	60 (6%) 27 31	10, 24, 45, 60	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	91	ALA	6.1
1	A	170	PHE	5.3
1	A	44	ARG	4.7
1	A	446	VAL	4.7
1	B	22	ARG	4.5
1	A	163	ALA	4.5
1	B	93	PRO	4.1
1	A	42	GLY	4.1
1	B	360	GLU	4.1
1	B	23	ALA	4.0
1	A	169	GLY	3.9
1	A	447	HIS	3.9
1	A	444	LEU	3.7
1	B	91	ALA	3.6
1	B	359	ARG	3.6
1	B	20	GLY	3.5
1	B	24	PRO	3.4
1	A	445	GLY	3.4
1	B	94	ALA	3.4
1	B	25	GLY	3.4
1	A	359	ARG	3.3
1	A	326	TYR	3.2
1	A	469	ARG	3.2
1	A	358	HIS	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	277	THR	3.0
1	A	92	GLY	3.0
1	A	93	PRO	3.0
1	A	357	MET	2.9
1	B	444	LEU	2.9
1	B	338	ASP	2.9
1	A	450	SER	2.8
1	B	355	GLY	2.8
1	A	490	ARG	2.8
1	B	357	MET	2.7
1	A	43	GLU	2.5
1	B	165	ARG	2.5
1	A	467	GLU	2.5
1	B	18	THR	2.5
1	B	358	HIS	2.5
1	A	41	GLN	2.4
1	B	174	ASP	2.4
1	A	40	PRO	2.4
1	B	173	ALA	2.4
1	A	507	ARG	2.3
1	A	69	ALA	2.3
1	A	506	ASP	2.3
1	B	326	TYR	2.3
1	A	17	ARG	2.2
1	B	21	GLY	2.2
1	B	92	GLY	2.2
1	B	96	GLU	2.2
1	B	89	LEU	2.1
1	A	421	ARG	2.1
1	B	250	GLY	2.1
1	A	355	GLY	2.1
1	B	277	THR	2.0
1	A	95	PRO	2.0
1	B	90	PRO	2.0
1	B	508	THR	2.0
1	A	241	GLY	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

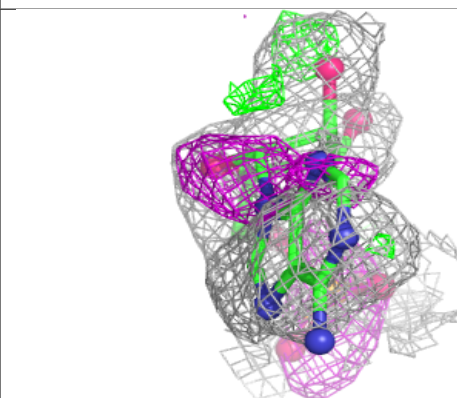
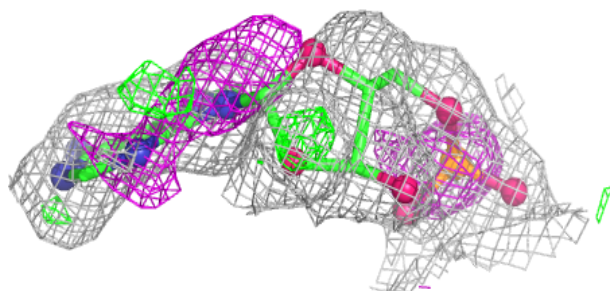
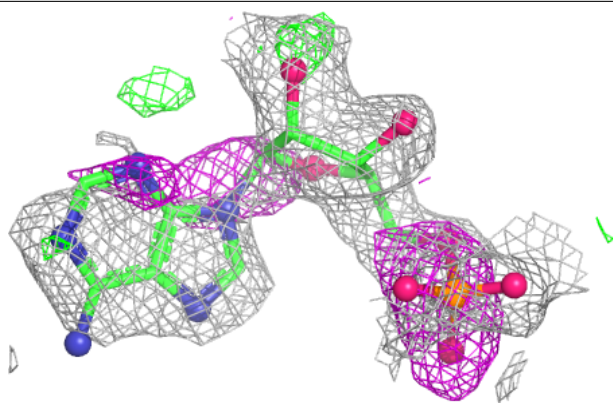
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	AMP	B	705	23/23	0.62	0.20	41,46,60,61	0
5	PCX	A	703	16/16	0.78	0.14	39,46,49,51	0
5	PCX	B	706	16/16	0.79	0.14	23,31,37,39	0
4	AMP	A	702	23/23	0.83	0.12	26,39,49,50	0
2	MG	A	603	1/1	0.94	0.05	35,35,35,35	0
3	POP	A	701	9/9	0.94	0.09	23,27,30,31	0
2	MG	B	605	1/1	0.95	0.06	29,29,29,29	0
2	MG	B	602	1/1	0.95	0.06	38,38,38,38	0
2	MG	A	601	1/1	0.96	0.06	32,32,32,32	0
3	POP	B	704	9/9	0.96	0.09	19,20,22,23	0

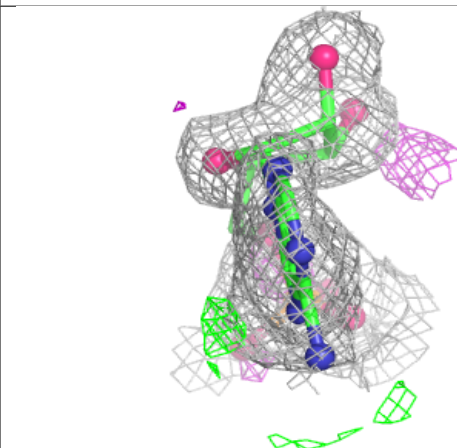
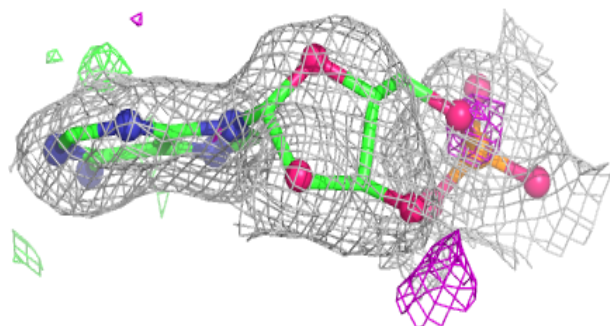
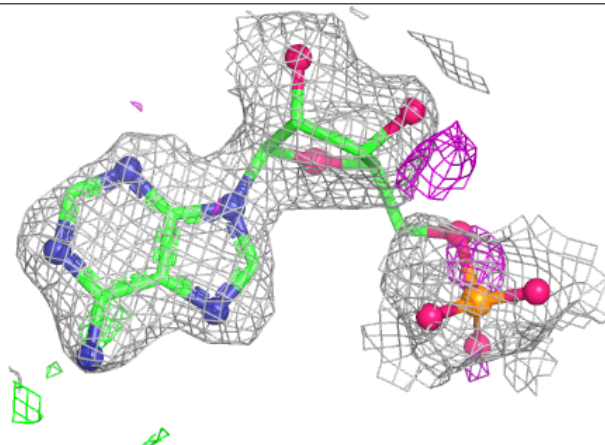
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around AMP B 705:

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

**Electron density around AMP A 702:**

$2mF_o-DF_c$ (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.