



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 23, 2026 – 07:20 AM EDT

PDB ID : 1C9K / pdb_00001c9k
Title : THE THREE DIMENSIONAL STRUCTURE OF ADENOSYLCOBINAMIDE KINASE/ ADENOSYLCOBINAMIDE PHOSPHATE GUALYLTRANSFERASE (COBU) COMPLEXED WITH GMP: EVIDENCE FOR A SUBSTRATE INDUCED TRANSFERASE ACTIVE SITE
Authors : Thompson, T.B.; Thomas, M.G.; Esclante-Semerena, J.C.; Rayment, I.
Deposited on : 1999-08-02
Resolution : 2.20 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](#) ①) 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

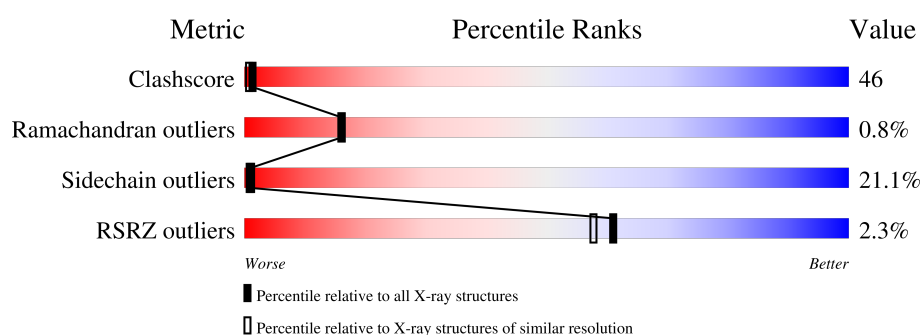
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	180	44% 36% 13% 6%
1	B	180	44% 40% 13% •
1	C	180	36% 37% 17% 8%

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
2	PO4	A	600	-	-	X	-
2	PO4	B	601	-	-	X	-

2 Entry composition [i](#)

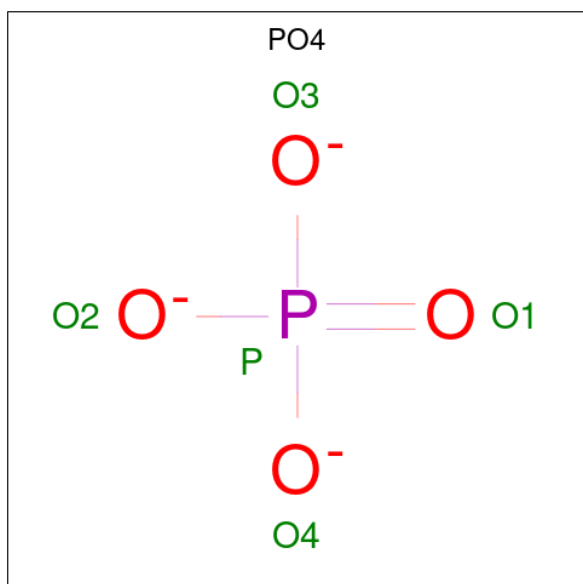
There are 6 unique types of molecules in this entry. The entry contains 4253 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ADENOSYLCOBINAMIDE KINASE.

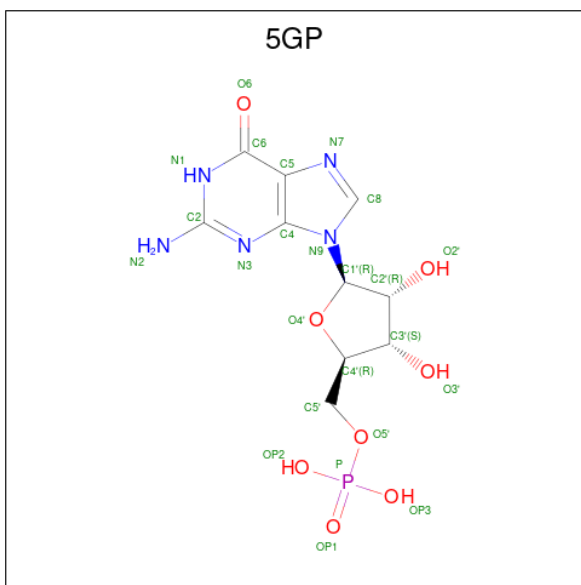
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	170	Total	C	N	O	S	0	0	0
			1287	812	232	235	8			
1	B	180	Total	C	N	O	S	0	0	0
			1328	837	240	242	9			
1	C	165	Total	C	N	O	S	0	0	0
			1214	768	215	224	7			

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is GUANOSINE-5'-MONOPHOSPHATE (CCD ID: 5GP) (formula: C₁₀H₁₄N₅O₈P).

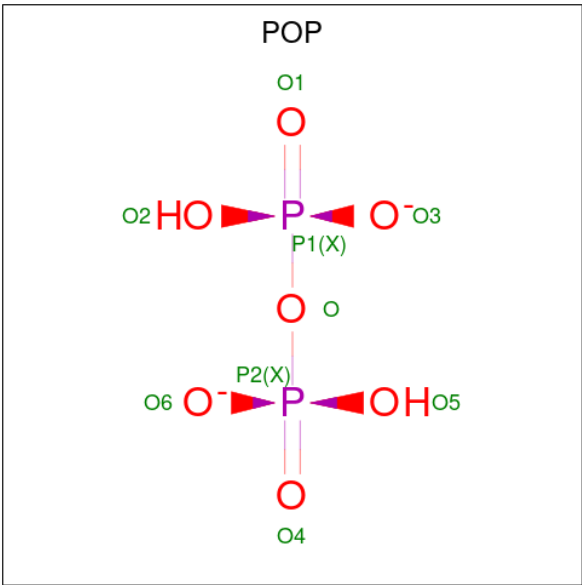


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			24	10	5	8	1		
3	B	1	Total	C	N	O	P	0	0
			24	10	5	8	1		
3	C	1	Total	C	N	O		0	0
			20	10	5	5			

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 5 is PYROPHOSPHATE 2- (CCD ID: POP) (formula: H₂O₇P₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	O	P	0	0
			9	7	2		

- Molecule 6 is water.

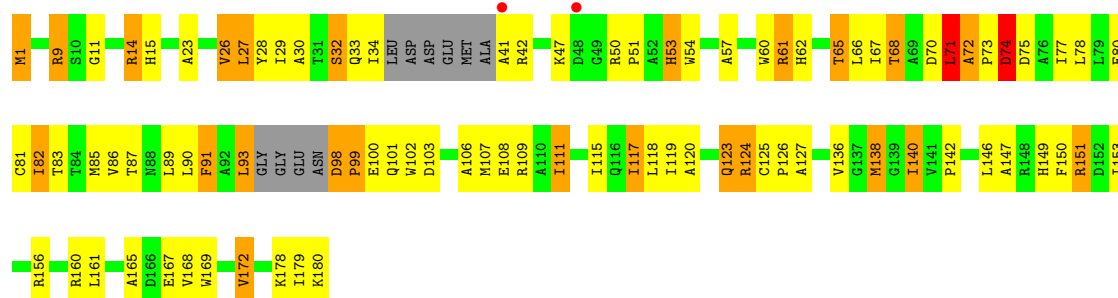
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	134	Total	O	0	0
			134	134		
6	B	116	Total	O	0	0
			116	116		
6	C	86	Total	O	0	0
			86	86		

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.

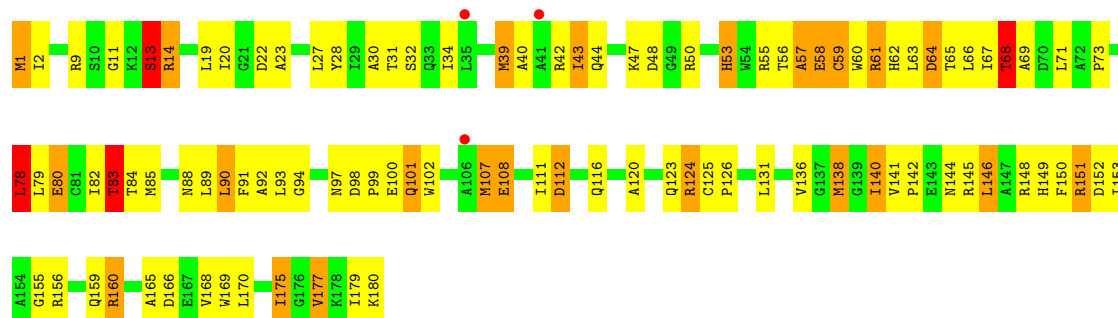
• Molecule 1: ADENOSYLCOBINAMIDE KINASE

Chain A: 

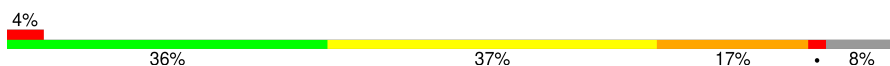


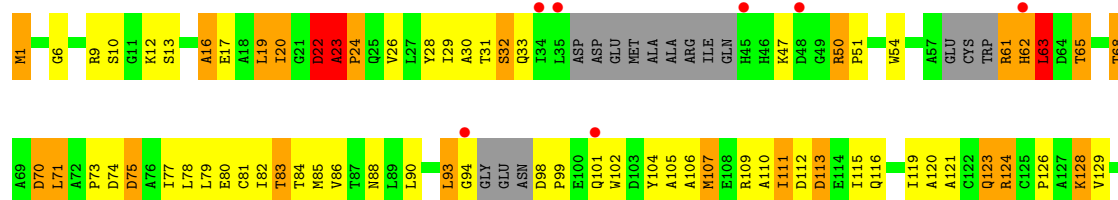
• Molecule 1: ADENOSYLCOBINAMIDE KINASE

Chain B: 



• Molecule 1: ADENOSYLCOBINAMIDE KINASE

Chain C: 



T133	M134		G137	M138	G139	I140		E143	N144	R145	L146	A147	R148	H149	F150	R151		G155	R156		Q159	R160	L161		A164	A165	D166	E167		L170	V171	V172		I175		K178	I179	K180
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.38Å 87.77Å 101.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.20 30.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.3 (30.00-2.20) 97.2 (30.00-2.20)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.53 (at 2.10Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.191 , (Not available) 0.194 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	26.2	Xtriage
Anisotropy	0.500	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 169.3	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	0.94	EDS
Total number of atoms	4253	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.02% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, POP, MG, 5GP

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	0.99	0/1306	1.61	20/1774 (1.1%)
1	B	1.17	3/1349 (0.2%)	1.74	22/1837 (1.2%)
1	C	1.11	2/1229 (0.2%)	1.87	28/1670 (1.7%)
All	All	1.10	5/3884 (0.1%)	1.74	70/5281 (1.3%)

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	1	0
1	B	0	1
1	C	1	0
All	All	2	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	58	GLU	N-CA	-13.60	1.29	1.46
1	C	23	ALA	CA-C	-13.00	1.36	1.52
1	C	23	ALA	N-CA	-6.58	1.36	1.46
1	B	57	ALA	C-N	-5.29	1.27	1.33
1	B	58	GLU	CA-C	-5.25	1.47	1.53

All (70) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	57	ALA	O-C-N	13.19	139.62	123.26
1	C	149	HIS	CA-CB-CG	-12.59	101.21	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	61	ARG	CA-C-N	11.79	139.10	122.36
1	C	61	ARG	C-N-CA	11.79	139.10	122.36
1	C	23	ALA	O-C-N	11.77	134.86	121.32
1	B	57	ALA	CA-C-N	10.36	138.19	122.23
1	B	57	ALA	C-N-CA	10.36	138.19	122.23
1	B	59	CYS	N-CA-C	9.70	123.88	109.24
1	C	62	HIS	CA-C-N	8.93	132.62	120.38
1	C	62	HIS	C-N-CA	8.93	132.62	120.38
1	B	166	ASP	N-CA-C	-8.00	103.31	113.23
1	C	22	ASP	N-CA-CB	-7.96	97.04	110.49
1	A	74	ASP	CA-CB-CG	7.90	120.50	112.60
1	C	31	THR	N-CA-C	7.86	122.13	112.54
1	A	172	VAL	N-CA-C	7.81	119.15	107.75
1	A	172	VAL	CB-CA-C	7.57	121.51	110.77
1	C	61	ARG	O-C-N	7.49	134.98	123.00
1	B	68	THR	N-CA-C	7.19	121.28	110.14
1	A	68	THR	N-CA-C	7.17	121.71	110.17
1	A	72	ALA	CA-C-N	6.89	126.40	119.24
1	A	72	ALA	C-N-CA	6.89	126.40	119.24
1	C	68	THR	CA-CB-OG1	6.79	119.79	109.60
1	A	82	ILE	N-CA-C	6.76	117.52	110.62
1	C	62	HIS	O-C-N	6.42	130.52	122.55
1	C	150	PHE	CA-CB-CG	-6.42	107.38	113.80
1	B	169	TRP	CB-CA-C	-6.38	98.49	109.65
1	B	64	ASP	CA-CB-CG	6.33	118.93	112.60
1	C	75	ASP	CA-CB-CG	6.33	118.93	112.60
1	C	63	LEU	N-CA-C	6.20	118.83	111.33
1	B	92	ALA	N-CA-C	-6.20	104.21	110.97
1	C	19	LEU	CA-C-N	6.19	128.36	120.56
1	C	19	LEU	C-N-CA	6.19	128.36	120.56
1	C	51	PRO	N-CA-CB	6.14	108.70	103.35
1	A	53	HIS	CA-CB-CG	-6.12	107.68	113.80
1	A	23	ALA	O-C-N	6.11	128.35	121.32
1	C	65	THR	CA-CB-OG1	6.07	118.71	109.60
1	C	88	ASN	CA-CB-CG	-5.92	106.67	112.60
1	A	98	ASP	CA-C-N	5.92	127.23	119.84
1	A	98	ASP	C-N-CA	5.92	127.23	119.84
1	A	32	SER	N-CA-CB	5.90	118.64	109.85
1	B	59	CYS	N-CA-CB	-5.81	101.58	111.69
1	B	23	ALA	N-CA-C	5.77	117.90	109.71
1	B	53	HIS	CA-CB-CG	-5.73	108.07	113.80
1	B	69	ALA	N-CA-C	5.72	119.98	112.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	14	ARG	N-CA-CB	-5.71	101.43	110.28
1	C	166	ASP	CA-CB-CG	5.71	118.31	112.60
1	A	23	ALA	CA-C-N	5.61	126.85	119.84
1	A	23	ALA	C-N-CA	5.61	126.85	119.84
1	C	68	THR	CA-CB-CG2	5.60	120.02	110.50
1	B	83	THR	N-CA-C	5.58	118.08	111.33
1	A	71	LEU	CA-C-N	5.52	128.07	120.67
1	A	71	LEU	C-N-CA	5.52	128.07	120.67
1	B	78	LEU	CA-C-N	-5.50	115.41	123.00
1	B	78	LEU	C-N-CA	-5.50	115.41	123.00
1	C	65	THR	CA-CB-CG2	5.50	119.85	110.50
1	C	33	GLN	N-CA-C	-5.46	100.81	109.76
1	A	91	PHE	CA-CB-CG	-5.43	108.37	113.80
1	C	68	THR	OG1-CB-CG2	5.41	120.12	109.30
1	B	22	ASP	N-CA-C	5.40	119.62	113.19
1	C	172	VAL	N-CA-CB	5.39	119.96	111.99
1	B	82	ILE	N-CA-C	5.37	116.00	110.36
1	A	150	PHE	CA-CB-CG	-5.36	108.44	113.80
1	B	146	LEU	O-C-N	5.34	129.70	122.59
1	B	13	SER	N-CA-CB	5.34	117.75	110.01
1	C	113	ASP	CA-CB-CG	-5.21	107.39	112.60
1	B	97	ASN	CA-CB-CG	5.21	117.81	112.60
1	C	149	HIS	O-C-N	5.17	127.40	122.07
1	A	26	VAL	N-CA-C	5.11	115.48	108.12
1	C	16	ALA	N-CA-C	-5.02	105.50	110.97
1	A	172	VAL	N-CA-CB	5.01	118.16	111.64

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	172	VAL	CA
1	C	68	THR	CB

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	58	GLU	Mainchain

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	1287	0	1297	140	0
1	B	1328	0	1318	133	0
1	C	1214	0	1215	128	0
2	A	5	0	0	2	0
2	B	5	0	0	3	0
3	A	24	0	12	4	0
3	B	24	0	12	1	0
3	C	20	0	12	2	0
4	C	1	0	0	0	0
5	C	9	0	0	0	0
6	A	134	0	0	9	0
6	B	116	0	0	6	0
6	C	86	0	0	5	0
All	All	4253	0	3866	358	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 46.

All (358) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:GLU:HG3	1:A:153:ILE:HG12	1.24	1.08
1:A:124:ARG:HH11	1:A:124:ARG:HB2	1.14	1.06
1:A:65:THR:HG22	1:A:66:LEU:HG	1.37	1.06
1:C:121:ALA:HA	1:C:124:ARG:HD3	1.38	1.05
1:B:68:THR:HG22	1:B:71:LEU:HG	1.35	1.04
1:B:34:ILE:HG21	1:B:40:ALA:HB2	1.40	0.98
1:A:9:ARG:NH2	1:C:159:GLN:HE22	1.61	0.97
1:A:9:ARG:HE	1:A:9:ARG:HA	1.30	0.94
1:B:102:TRP:CD2	1:B:107:MET:HE1	2.03	0.93
1:B:34:ILE:HG23	1:B:39:MET:HG2	1.48	0.93
1:C:145:ARG:HG3	1:C:149:HIS:CE1	2.03	0.93
1:A:156:ARG:HD3	1:B:9:ARG:HH21	1.33	0.92
1:A:138:MET:HE2	1:C:138:MET:HG2	1.51	0.90
1:B:179:ILE:HD12	1:C:175:ILE:HB	1.54	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:ARG:HD3	1:C:156:ARG:HD2	1.55	0.88
1:B:43:ILE:HG22	1:B:47:LYS:HE2	1.55	0.88
1:C:156:ARG:HH11	1:C:156:ARG:HG2	1.40	0.87
1:A:9:ARG:HD3	1:C:156:ARG:CD	2.05	0.86
1:A:102:TRP:CD2	1:A:107:MET:HE1	2.11	0.86
1:A:82:ILE:HD13	1:A:85:MET:HE1	1.59	0.85
1:A:138:MET:HB3	1:C:138:MET:HB3	1.60	0.83
1:B:1:MET:HE3	1:B:165:ALA:CA	2.09	0.83
1:B:141:VAL:HG22	1:B:151:ARG:NH2	1.93	0.83
1:C:121:ALA:CA	1:C:124:ARG:HD3	2.09	0.82
1:B:43:ILE:HG22	1:B:47:LYS:CE	2.09	0.82
1:B:34:ILE:HG21	1:B:40:ALA:CB	2.08	0.81
1:A:124:ARG:HB2	1:A:124:ARG:NH1	1.94	0.81
1:A:9:ARG:HH11	1:C:156:ARG:NH1	1.78	0.81
1:B:68:THR:CG2	1:B:71:LEU:HG	2.10	0.81
1:A:167:GLU:HG2	1:A:178:LYS:HE2	1.61	0.81
1:C:99:PRO:HA	1:C:102:TRP:CD2	2.15	0.81
1:A:61:ARG:HH11	1:A:61:ARG:HG3	1.45	0.80
1:A:82:ILE:HD13	1:A:85:MET:CE	2.11	0.80
1:A:136:VAL:HB	1:C:140:ILE:CD1	2.10	0.80
1:A:50:ARG:NH2	3:A:604:5GP:N7	2.30	0.80
1:B:108:GLU:HG3	1:B:153:ILE:HG12	1.64	0.80
1:B:141:VAL:HG22	1:B:151:ARG:HH22	1.47	0.79
1:B:90:LEU:HD23	1:B:111:ILE:HD11	1.64	0.78
1:A:106:ALA:HA	1:A:109:ARG:NH1	1.98	0.78
1:A:138:MET:HB3	1:B:138:MET:HB3	1.66	0.78
1:B:102:TRP:CE3	1:B:107:MET:HE1	2.19	0.78
1:B:138:MET:HB3	1:C:138:MET:HB3	1.63	0.77
1:A:138:MET:HB3	1:B:138:MET:CB	2.14	0.77
1:C:145:ARG:HG3	1:C:149:HIS:ND1	1.99	0.77
1:B:43:ILE:HG22	1:B:47:LYS:NZ	1.98	0.77
1:C:90:LEU:HD23	1:C:111:ILE:HD11	1.67	0.77
1:C:9:ARG:HA	1:C:9:ARG:NE	2.01	0.76
1:B:68:THR:HG22	1:B:71:LEU:CG	2.15	0.75
1:C:68:THR:HG22	1:C:71:LEU:HD13	1.69	0.75
1:A:140:ILE:HD11	1:B:151:ARG:HB2	1.68	0.74
1:C:99:PRO:HA	1:C:102:TRP:CG	2.22	0.73
1:B:1:MET:HE3	1:B:165:ALA:HA	1.70	0.73
1:B:138:MET:HB3	1:C:138:MET:CB	2.19	0.73
1:A:26:VAL:HG21	1:A:54:TRP:NE1	2.04	0.73
1:B:138:MET:CG	1:C:138:MET:HE2	2.18	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:SER:OG	3:A:604:5GP:H1'	1.89	0.72
1:C:28:TYR:CE1	1:C:30:ALA:HB2	2.23	0.72
1:C:13:SER:O	1:C:17:GLU:HG3	1.89	0.72
1:B:138:MET:HG2	1:C:138:MET:HE2	1.71	0.72
1:B:152:ASP:HB3	1:B:156:ARG:NH2	2.05	0.72
1:B:1:MET:HE3	1:B:165:ALA:HB2	1.70	0.71
1:A:68:THR:HB	1:A:71:LEU:HD22	1.71	0.71
1:A:140:ILE:O	1:A:142:PRO:HD3	1.90	0.71
1:A:53:HIS:H	1:A:53:HIS:CD2	2.09	0.71
1:A:99:PRO:HA	1:A:102:TRP:CE2	2.26	0.70
1:C:61:ARG:C	1:C:63:LEU:HD23	2.16	0.70
1:B:1:MET:HE3	1:B:165:ALA:CB	2.22	0.70
1:C:62:HIS:O	1:C:65:THR:HB	1.90	0.70
1:A:138:MET:CB	1:C:138:MET:HB3	2.21	0.70
1:B:99:PRO:HA	1:B:102:TRP:CD2	2.27	0.70
1:C:144:ASN:O	1:C:147:ALA:HB3	1.91	0.70
1:A:72:ALA:HB3	1:A:75:ASP:OD1	1.91	0.69
1:B:179:ILE:HD12	1:C:175:ILE:CB	2.22	0.69
1:C:9:ARG:HA	1:C:9:ARG:HE	1.57	0.69
1:C:99:PRO:HA	1:C:102:TRP:CE2	2.27	0.69
1:C:73:PRO:HA	1:C:126:PRO:O	1.93	0.69
1:C:1:MET:HB3	1:C:129:VAL:HB	1.75	0.69
1:A:62:HIS:O	1:A:65:THR:HB	1.92	0.69
1:A:42:ARG:N	1:A:42:ARG:HD3	2.07	0.68
1:B:20:ILE:HD12	1:B:78:LEU:HG	1.75	0.68
1:C:121:ALA:HA	1:C:124:ARG:CD	2.20	0.68
1:B:47:LYS:HE3	6:B:812:HOH:O	1.92	0.68
1:A:99:PRO:HA	1:A:102:TRP:CD2	2.29	0.68
1:A:156:ARG:CD	1:B:9:ARG:HH21	2.07	0.68
1:B:156:ARG:NH1	1:C:9:ARG:HD3	2.10	0.67
1:B:98:ASP:HB3	1:B:101:GLN:NE2	2.09	0.67
1:A:9:ARG:NH1	1:C:156:ARG:NH1	2.41	0.67
1:C:115:ILE:O	1:C:119:ILE:HG13	1.96	0.66
1:A:9:ARG:HE	1:A:9:ARG:CA	2.04	0.66
1:B:34:ILE:HG21	1:B:40:ALA:CA	2.26	0.66
1:A:61:ARG:HG3	1:A:61:ARG:NH1	2.10	0.66
1:B:34:ILE:HA	1:B:39:MET:HG2	1.77	0.66
1:B:11:GLY:HA2	6:B:749:HOH:O	1.96	0.66
1:A:9:ARG:HD3	1:C:156:ARG:NE	2.11	0.65
1:A:68:THR:H	1:A:71:LEU:CD2	2.10	0.65
1:B:9:ARG:HE	1:B:9:ARG:HA	1.62	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:119:ILE:HD13	1:C:164:ALA:HB2	1.77	0.65
1:C:68:THR:CG2	1:C:71:LEU:HD13	2.26	0.65
1:B:112:ASP:OD2	1:B:160:ARG:NH1	2.30	0.65
1:B:146:LEU:O	1:B:149:HIS:N	2.30	0.64
1:A:9:ARG:NH2	1:C:159:GLN:NE2	2.42	0.64
1:B:155:GLY:O	1:B:159:GLN:HG3	1.97	0.64
1:B:1:MET:CE	1:B:165:ALA:HB2	2.26	0.64
1:C:137:GLY:HA2	6:C:707:HOH:O	1.96	0.64
1:B:71:LEU:O	1:B:126:PRO:HG2	1.96	0.64
1:C:83:THR:O	1:C:86:VAL:HB	1.96	0.64
1:C:90:LEU:CD2	1:C:111:ILE:HD11	2.27	0.64
1:A:82:ILE:O	1:A:86:VAL:HG23	1.98	0.64
1:A:136:VAL:HB	1:C:140:ILE:HD12	1.79	0.64
1:C:28:TYR:HE1	1:C:30:ALA:HB2	1.62	0.64
1:A:153:ILE:HG23	6:A:862:HOH:O	1.97	0.64
1:C:102:TRP:CD2	1:C:107:MET:HE1	2.33	0.64
1:A:27:LEU:HD11	1:A:57:ALA:HB2	1.80	0.63
1:C:68:THR:HG22	1:C:71:LEU:HD22	1.80	0.63
1:A:9:ARG:HD2	6:A:763:HOH:O	1.98	0.63
1:A:93:LEU:N	1:A:93:LEU:HD23	2.13	0.63
1:A:98:ASP:O	1:A:100:GLU:N	2.30	0.63
1:A:29:ILE:HD12	1:A:77:ILE:HG23	1.80	0.63
1:A:167:GLU:HG2	1:A:178:LYS:CE	2.28	0.62
1:A:11:GLY:N	2:A:600:PO4:O2	2.30	0.62
1:B:63:LEU:HD23	1:B:66:LEU:HD12	1.79	0.62
1:C:99:PRO:HA	1:C:102:TRP:CD1	2.34	0.62
1:C:104:TYR:HE1	1:C:145:ARG:HE	1.47	0.62
1:A:124:ARG:HH11	1:A:124:ARG:CB	2.02	0.61
1:C:29:ILE:HD12	1:C:77:ILE:CG2	2.30	0.61
1:C:68:THR:HG22	1:C:71:LEU:CD1	2.30	0.61
1:C:106:ALA:HA	1:C:109:ARG:NH1	2.16	0.61
1:A:140:ILE:HD12	1:B:136:VAL:O	2.01	0.60
1:A:156:ARG:NH1	1:B:9:ARG:HD2	2.16	0.60
1:B:94:GLY:HA3	1:B:102:TRP:CZ2	2.36	0.60
1:C:145:ARG:O	1:C:149:HIS:ND1	2.34	0.60
1:B:98:ASP:HB3	1:B:101:GLN:HE21	1.65	0.60
1:C:155:GLY:O	1:C:159:GLN:HG3	2.01	0.60
1:B:98:ASP:O	1:B:101:GLN:HG2	2.01	0.60
1:A:61:ARG:HH11	1:A:61:ARG:CG	2.15	0.59
1:B:142:PRO:HG2	1:B:148:ARG:HA	1.83	0.59
1:A:60:TRP:HZ3	1:A:61:ARG:HH12	1.49	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:28:TYR:CE1	1:B:30:ALA:HB2	2.38	0.59
1:C:145:ARG:HG3	1:C:149:HIS:HD1	1.68	0.59
1:C:156:ARG:HH11	1:C:156:ARG:CG	2.13	0.59
1:C:102:TRP:CG	1:C:107:MET:HE1	2.39	0.58
1:A:32:SER:O	3:A:604:5GP:N2	2.31	0.58
1:A:156:ARG:HG3	1:B:9:ARG:HD3	1.85	0.58
1:B:20:ILE:CD1	1:B:78:LEU:HG	2.32	0.58
1:C:29:ILE:HD12	1:C:77:ILE:HG21	1.85	0.58
1:B:34:ILE:CG2	1:B:39:MET:HG2	2.28	0.58
1:A:9:ARG:NH2	2:A:600:PO4:O2	2.34	0.58
1:A:51:PRO:HB2	1:A:53:HIS:CD2	2.39	0.58
1:A:102:TRP:CE3	1:A:107:MET:HE1	2.39	0.57
1:A:156:ARG:HD3	1:B:9:ARG:NH2	2.13	0.57
1:A:169:TRP:CZ2	1:A:178:LYS:HD2	2.39	0.57
1:B:34:ILE:HG12	1:B:43:ILE:HD12	1.86	0.57
1:A:160:ARG:NH2	6:A:829:HOH:O	2.37	0.57
1:A:136:VAL:O	1:C:140:ILE:HD13	2.05	0.56
1:B:120:ALA:O	1:B:124:ARG:HG3	2.05	0.56
1:B:34:ILE:HG23	1:B:39:MET:CG	2.30	0.56
1:C:22:ASP:OD1	1:C:23:ALA:N	2.39	0.56
1:B:156:ARG:HH11	1:C:9:ARG:HD3	1.69	0.56
1:C:70:ASP:HB3	6:C:963:HOH:O	2.05	0.56
1:A:74:ASP:HB3	6:A:878:HOH:O	2.06	0.56
1:B:44:GLN:HA	1:B:47:LYS:HE3	1.87	0.56
1:A:138:MET:HB3	1:B:138:MET:HB2	1.88	0.55
1:C:104:TYR:CD2	1:C:149:HIS:HB3	2.41	0.55
1:B:98:ASP:OD2	1:B:100:GLU:N	2.30	0.55
1:A:9:ARG:HH11	1:C:156:ARG:CZ	2.19	0.55
1:A:29:ILE:HD12	1:A:77:ILE:CG2	2.37	0.55
1:A:102:TRP:HB3	1:A:107:MET:SD	2.46	0.55
1:A:93:LEU:HD12	1:A:107:MET:HG2	1.90	0.54
1:B:27:LEU:HD12	1:B:55:ARG:O	2.06	0.54
1:C:105:ALA:O	1:C:109:ARG:HG3	2.07	0.54
1:B:170:LEU:O	1:B:177:VAL:N	2.37	0.54
1:A:87:THR:O	1:A:91:PHE:HD1	1.91	0.54
1:A:108:GLU:HG3	1:A:153:ILE:CG1	2.18	0.54
1:C:143:GLU:O	1:C:148:ARG:NH1	2.41	0.54
1:A:68:THR:CB	1:A:71:LEU:HD22	2.38	0.54
1:C:50:ARG:HD3	1:C:54:TRP:HE3	1.73	0.54
1:A:117:ILE:HG22	1:A:118:LEU:N	2.22	0.54
1:B:63:LEU:O	1:B:67:ILE:N	2.30	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:ASP:O	1:A:127:ALA:HB1	2.08	0.53
1:C:68:THR:HG22	1:C:71:LEU:CD2	2.39	0.53
1:A:14:ARG:HG3	1:A:15:HIS:N	2.21	0.53
1:A:102:TRP:CG	1:A:107:MET:HE1	2.44	0.53
1:B:53:HIS:HE1	6:B:881:HOH:O	1.92	0.53
1:C:120:ALA:O	1:C:123:GLN:HB2	2.08	0.53
1:C:144:ASN:HD22	1:C:147:ALA:HB2	1.72	0.53
1:B:34:ILE:CG2	1:B:40:ALA:HB2	2.26	0.53
1:B:159:GLN:HE22	1:C:9:ARG:HH11	1.57	0.52
1:A:28:TYR:HE1	1:A:30:ALA:HB2	1.75	0.52
1:C:28:TYR:CD1	1:C:30:ALA:HB2	2.44	0.52
1:A:9:ARG:CZ	1:C:159:GLN:HE22	2.22	0.52
1:A:27:LEU:HD11	1:A:57:ALA:CB	2.39	0.52
1:A:32:SER:HG	3:A:604:5GP:H1'	1.73	0.52
1:B:73:PRO:HA	1:B:126:PRO:O	2.10	0.52
1:B:47:LYS:O	1:B:50:ARG:N	2.33	0.51
1:A:71:LEU:O	1:A:126:PRO:HG2	2.11	0.51
1:A:117:ILE:HA	6:A:825:HOH:O	2.11	0.51
1:B:98:ASP:OD2	1:B:99:PRO:N	2.44	0.51
1:A:34:ILE:HD13	6:A:924:HOH:O	2.11	0.51
1:B:85:MET:O	1:B:89:LEU:HD12	2.11	0.51
1:C:121:ALA:N	1:C:124:ARG:HH11	2.07	0.51
1:C:167:GLU:HG2	1:C:178:LYS:CE	2.40	0.51
1:A:89:LEU:O	1:A:93:LEU:HG	2.10	0.51
1:B:9:ARG:HA	2:B:601:PO4:O4	2.11	0.51
1:C:90:LEU:O	1:C:94:GLY:N	2.43	0.51
1:B:79:LEU:HD23	1:B:79:LEU:C	2.36	0.51
1:B:62:HIS:O	1:B:65:THR:HB	2.11	0.50
1:B:102:TRP:CG	1:B:107:MET:HE1	2.46	0.50
1:A:72:ALA:O	1:A:75:ASP:HB2	2.11	0.50
1:B:44:GLN:O	1:B:48:ASP:OD2	2.30	0.50
1:A:9:ARG:HA	1:A:9:ARG:NE	2.11	0.50
1:B:56:THR:HG22	1:B:57:ALA:N	2.26	0.50
1:B:141:VAL:HA	1:B:151:ARG:NH2	2.27	0.50
1:C:75:ASP:HA	6:C:758:HOH:O	2.12	0.50
1:A:82:ILE:HD13	1:A:85:MET:HE3	1.91	0.49
1:C:79:LEU:C	1:C:79:LEU:HD23	2.38	0.49
1:C:98:ASP:O	1:C:101:GLN:HG2	2.12	0.49
1:B:28:TYR:HE1	1:B:30:ALA:HB2	1.77	0.49
1:A:101:GLN:HG3	1:A:101:GLN:O	2.11	0.49
1:B:83:THR:HG22	1:B:150:PHE:HZ	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:146:LEU:HD23	1:B:146:LEU:C	2.37	0.49
1:C:22:ASP:O	1:C:23:ALA:O	2.31	0.49
1:C:145:ARG:HG3	1:C:149:HIS:HE1	1.73	0.49
1:C:102:TRP:CZ3	1:C:146:LEU:HD12	2.48	0.49
1:A:26:VAL:HG21	1:A:54:TRP:HE1	1.77	0.48
1:A:53:HIS:H	1:A:53:HIS:HD2	1.57	0.48
1:A:67:ILE:HG22	1:A:125:CYS:HB2	1.95	0.48
1:A:60:TRP:CZ3	1:A:61:ARG:NH1	2.81	0.48
1:A:67:ILE:HG12	1:A:77:ILE:HD13	1.95	0.48
1:B:34:ILE:HG12	1:B:43:ILE:CD1	2.42	0.48
1:A:80:GLU:HA	1:A:81:CYS:HA	1.64	0.48
1:B:48:ASP:OD2	1:B:48:ASP:N	2.45	0.48
1:C:32:SER:OG	3:C:606:5GP:H1'	2.14	0.48
1:C:112:ASP:O	1:C:116:GLN:HG2	2.13	0.48
1:C:106:ALA:HA	1:C:109:ARG:CZ	2.44	0.48
1:A:108:GLU:OE2	1:A:156:ARG:NE	2.45	0.47
1:B:39:MET:HE2	1:B:39:MET:HB2	1.69	0.47
1:B:140:ILE:CD1	1:C:151:ARG:HB3	2.44	0.47
1:C:156:ARG:HG2	1:C:156:ARG:NH1	2.11	0.47
1:B:43:ILE:HG22	1:B:47:LYS:HZ1	1.79	0.47
1:A:60:TRP:HZ3	1:A:61:ARG:NH1	2.11	0.47
1:B:1:MET:C	1:B:2:ILE:HG13	2.38	0.47
1:B:43:ILE:C	1:B:47:LYS:HE2	2.39	0.47
1:A:160:ARG:HD2	6:A:1056:HOH:O	2.14	0.47
1:B:124:ARG:NH2	6:B:1065:HOH:O	2.31	0.47
1:A:115:ILE:O	1:A:119:ILE:HG13	2.14	0.47
1:B:138:MET:HB3	1:C:138:MET:HB2	1.95	0.47
1:A:108:GLU:HA	1:A:111:ILE:HG13	1.96	0.47
1:C:119:ILE:HD13	1:C:164:ALA:CB	2.44	0.47
1:A:26:VAL:HB	1:A:54:TRP:CD1	2.50	0.46
1:B:53:HIS:H	1:B:53:HIS:CD2	2.32	0.46
1:B:144:ASN:ND2	1:B:145:ARG:N	2.64	0.46
1:C:71:LEU:HD12	1:C:71:LEU:HA	1.68	0.46
1:C:107:MET:O	1:C:111:ILE:HD12	2.15	0.46
1:C:121:ALA:C	1:C:124:ARG:HD3	2.39	0.46
1:C:145:ARG:CG	1:C:149:HIS:HD1	2.28	0.46
1:C:93:LEU:HD12	1:C:93:LEU:HA	1.69	0.46
1:C:145:ARG:HG2	1:C:146:LEU:N	2.30	0.46
1:A:73:PRO:HA	1:A:126:PRO:O	2.16	0.46
1:B:98:ASP:H	1:B:101:GLN:CD	2.23	0.46
1:B:80:GLU:HG2	3:B:605:5GP:H3'	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:26:VAL:HB	1:C:54:TRP:CD1	2.51	0.46
1:A:65:THR:CG2	1:A:66:LEU:HG	2.25	0.46
1:A:68:THR:H	1:A:71:LEU:HD23	1.80	0.46
1:C:68:THR:HG23	1:C:70:ASP:H	1.80	0.46
1:C:110:ALA:O	1:C:113:ASP:HB2	2.16	0.46
1:A:53:HIS:CD2	1:A:53:HIS:N	2.80	0.45
1:B:13:SER:OG	2:B:601:PO4:O2	2.33	0.45
1:C:29:ILE:CD1	1:C:77:ILE:HG21	2.45	0.45
1:C:22:ASP:CG	1:C:23:ALA:H	2.23	0.45
1:A:27:LEU:HD23	1:A:29:ILE:HD11	1.98	0.45
1:B:61:ARG:NH1	1:B:89:LEU:HD23	2.30	0.45
1:B:63:LEU:CD2	1:B:66:LEU:HD12	2.46	0.45
1:B:99:PRO:HA	1:B:102:TRP:CG	2.51	0.45
1:B:179:ILE:CD1	1:C:175:ILE:HB	2.38	0.45
1:C:10:SER:OG	1:C:12:LYS:HG2	2.16	0.45
1:C:23:ALA:HA	1:C:24:PRO:HD3	1.60	0.45
1:A:27:LEU:CD2	1:A:29:ILE:HD11	2.46	0.45
1:C:167:GLU:HA	1:C:180:LYS:OXT	2.17	0.45
1:A:103:ASP:OD1	1:A:106:ALA:HB2	2.16	0.45
1:B:9:ARG:HA	1:B:9:ARG:NE	2.30	0.45
1:A:156:ARG:CG	1:B:9:ARG:NH2	2.80	0.45
1:B:141:VAL:CG2	1:B:151:ARG:HH22	2.25	0.45
1:B:56:THR:CG2	1:B:57:ALA:N	2.80	0.44
1:B:61:ARG:HH11	1:B:89:LEU:HD23	1.81	0.44
1:C:6:GLY:O	1:C:134:ASN:HA	2.18	0.44
1:C:121:ALA:O	1:C:124:ARG:HD3	2.18	0.44
1:A:120:ALA:O	1:A:123:GLN:HB2	2.17	0.44
1:B:83:THR:CG2	1:B:150:PHE:HZ	2.31	0.44
1:A:156:ARG:NH1	1:B:9:ARG:CD	2.80	0.44
1:B:94:GLY:HA3	1:B:102:TRP:CE2	2.52	0.44
1:A:156:ARG:CG	1:B:9:ARG:HH21	2.30	0.44
1:A:168:VAL:HB	1:A:180:LYS:HB3	1.99	0.44
1:A:98:ASP:O	1:A:101:GLN:HG2	2.17	0.44
1:B:31:THR:O	1:B:88:ASN:ND2	2.51	0.44
1:B:71:LEU:HD23	1:B:71:LEU:HA	1.84	0.44
1:A:41:ALA:HB3	1:A:42:ARG:NH2	2.32	0.44
1:B:79:LEU:HD23	1:B:80:GLU:N	2.33	0.44
1:B:91:PHE:CD2	1:B:102:TRP:HH2	2.36	0.44
1:B:98:ASP:CB	1:B:101:GLN:NE2	2.80	0.44
1:B:59:CYS:SG	1:B:60:TRP:N	2.90	0.43
1:C:167:GLU:HG2	1:C:178:LYS:HE3	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:THR:HG21	1:A:70:ASP:OD1	2.18	0.43
1:A:93:LEU:HD23	1:A:93:LEU:H	1.81	0.43
1:C:54:TRP:CD1	1:C:54:TRP:N	2.86	0.43
1:C:98:ASP:OD2	1:C:99:PRO:N	2.51	0.43
1:C:144:ASN:ND2	6:C:928:HOH:O	2.51	0.43
1:A:99:PRO:CA	1:A:102:TRP:CE2	3.01	0.43
1:A:146:LEU:O	1:A:149:HIS:N	2.52	0.43
1:C:29:ILE:O	1:C:80:GLU:N	2.36	0.43
1:C:145:ARG:C	1:C:149:HIS:HD1	2.26	0.43
1:B:34:ILE:CG2	1:B:40:ALA:N	2.82	0.43
1:B:85:MET:HG2	1:B:89:LEU:HD11	1.99	0.43
1:C:12:LYS:HE3	1:C:133:THR:O	2.18	0.43
1:B:90:LEU:CD2	1:B:111:ILE:HD11	2.43	0.43
1:C:24:PRO:HD2	1:C:74:ASP:O	2.19	0.43
1:C:111:ILE:O	1:C:115:ILE:HG12	2.17	0.42
1:A:9:ARG:CD	1:C:156:ARG:HD2	2.39	0.42
1:A:68:THR:HG22	1:A:70:ASP:H	1.84	0.42
1:C:80:GLU:HA	1:C:81:CYS:HA	1.75	0.42
1:C:83:THR:HG23	6:C:712:HOH:O	2.19	0.42
1:B:140:ILE:HG13	6:B:733:HOH:O	2.19	0.42
1:A:68:THR:HB	1:A:71:LEU:CD2	2.46	0.42
1:A:108:GLU:CG	1:A:153:ILE:HG12	2.18	0.42
1:C:17:GLU:O	1:C:20:ILE:HB	2.20	0.42
1:A:26:VAL:CG2	1:A:54:TRP:NE1	2.80	0.42
1:A:65:THR:O	1:A:65:THR:HG23	2.20	0.42
1:C:128:LYS:HE2	1:C:128:LYS:HB2	1.93	0.42
1:A:161:LEU:O	1:A:165:ALA:N	2.40	0.42
1:C:120:ALA:C	1:C:124:ARG:HH11	2.27	0.42
1:B:90:LEU:HD23	1:B:111:ILE:CD1	2.40	0.42
1:A:28:TYR:CE1	1:A:30:ALA:HB2	2.55	0.41
1:A:82:ILE:HA	1:A:85:MET:HE3	2.02	0.41
1:A:147:ALA:O	1:A:151:ARG:HB3	2.20	0.41
1:C:16:ALA:O	1:C:19:LEU:HB2	2.20	0.41
1:C:161:LEU:O	1:C:165:ALA:N	2.39	0.41
1:A:165:ALA:O	1:A:180:LYS:NZ	2.32	0.41
1:B:9:ARG:NE	2:B:601:PO4:O4	2.45	0.41
1:B:168:VAL:HG23	1:B:180:LYS:HD2	2.02	0.41
6:A:1095:HOH:O	1:B:14:ARG:CB	2.68	0.41
1:C:140:ILE:HD12	1:C:140:ILE:HA	1.71	0.41
1:A:111:ILE:O	1:A:115:ILE:HG12	2.20	0.41
1:B:145:ARG:HA	1:B:148:ARG:NH2	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:50:ARG:NH2	3:C:606:5GP:N7	2.67	0.41
1:A:33:GLN:NE2	6:A:877:HOH:O	2.53	0.41
1:A:1:MET:HG3	1:A:165:ALA:HA	2.03	0.41
1:B:62:HIS:C	1:B:64:ASP:H	2.29	0.41
1:B:146:LEU:O	1:B:150:PHE:N	2.49	0.41
1:A:102:TRP:CE3	1:A:107:MET:CE	3.03	0.40
1:A:156:ARG:HH11	1:B:9:ARG:HD2	1.83	0.40
1:B:44:GLN:HA	6:B:812:HOH:O	2.21	0.40
1:B:99:PRO:HA	1:B:102:TRP:CE2	2.55	0.40
1:A:98:ASP:C	1:A:100:GLU:H	2.29	0.40
1:A:169:TRP:CZ3	1:A:178:LYS:HB2	2.56	0.40
1:A:72:ALA:HA	1:A:73:PRO:HD3	1.92	0.40
1:A:180:LYS:HB2	1:B:175:ILE:HD13	2.04	0.40
1:C:29:ILE:O	1:C:79:LEU:HA	2.22	0.40
1:B:125:CYS:HA	1:B:126:PRO:HD3	1.84	0.40
1:C:82:ILE:O	1:C:86:VAL:HG23	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	164/180 (91%)	152 (93%)	11 (7%)	1 (1%)	21	23
1	B	178/180 (99%)	169 (95%)	9 (5%)	0	100	100
1	C	157/180 (87%)	149 (95%)	5 (3%)	3 (2%)	6	4
All	All	499/540 (92%)	470 (94%)	25 (5%)	4 (1%)	16	16

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	99	PRO

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Mol	Chain	Res	Type
1	C	22	ASP
1	C	23	ALA
1	C	24	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	130/143 (91%)	108 (83%)	22 (17%)	2	2
1	B	129/143 (90%)	100 (78%)	29 (22%)	1	1
1	C	120/143 (84%)	91 (76%)	29 (24%)	1	0
All	All	379/429 (88%)	299 (79%)	80 (21%)	1	1

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	9	ARG
1	A	14	ARG
1	A	27	LEU
1	A	47	LYS
1	A	61	ARG
1	A	65	THR
1	A	71	LEU
1	A	74	ASP
1	A	78	LEU
1	A	83	THR
1	A	90	LEU
1	A	93	LEU
1	A	111	ILE
1	A	117	ILE
1	A	123	GLN
1	A	124	ARG
1	A	138	MET
1	A	140	ILE

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Mol	Chain	Res	Type
1	A	151	ARG
1	A	172	VAL
1	A	179	ILE
1	B	1	MET
1	B	13	SER
1	B	19	LEU
1	B	32	SER
1	B	39	MET
1	B	42	ARG
1	B	43	ILE
1	B	61	ARG
1	B	68	THR
1	B	78	LEU
1	B	80	GLU
1	B	83	THR
1	B	84	THR
1	B	90	LEU
1	B	93	LEU
1	B	101	GLN
1	B	107	MET
1	B	108	GLU
1	B	112	ASP
1	B	116	GLN
1	B	123	GLN
1	B	124	ARG
1	B	131	LEU
1	B	138	MET
1	B	140	ILE
1	B	151	ARG
1	B	160	ARG
1	B	175	ILE
1	B	177	VAL
1	C	1	MET
1	C	20	ILE
1	C	22	ASP
1	C	32	SER
1	C	47	LYS
1	C	50	ARG
1	C	63	LEU
1	C	70	ASP
1	C	71	LEU
1	C	78	LEU

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Mol	Chain	Res	Type
1	C	83	THR
1	C	84	THR
1	C	85	MET
1	C	93	LEU
1	C	107	MET
1	C	111	ILE
1	C	123	GLN
1	C	124	ARG
1	C	128	LYS
1	C	138	MET
1	C	140	ILE
1	C	145	ARG
1	C	148	ARG
1	C	156	ARG
1	C	160	ARG
1	C	170	LEU
1	C	172	VAL
1	C	178	LYS
1	C	179	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	HIS
1	A	116	GLN
1	A	144	ASN
1	B	53	HIS
1	B	97	ASN
1	B	101	GLN
1	B	134	ASN
1	B	144	ASN
1	B	159	GLN
1	C	53	HIS
1	C	116	GLN
1	C	134	ASN
1	C	144	ASN
1	C	159	GLN

5.3.3 RNA ⓘ

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 7 ligands modelled in this entry, 1 is 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$
3	5GP	B	605	-	26,26,26	0.97	1 (3%)	39,40,40	2.82	4 (10%)
5	POP	C	608	4	6,8,8	1.23	0	12,13,13	3.89	3 (25%)
2	PO4	B	601	-	4,4,4	1.64	2 (50%)	6,6,6	0.67	0
3	5GP	A	604	-	26,26,26	0.88	1 (3%)	39,40,40	1.26	4 (10%)
2	PO4	A	600	-	4,4,4	1.76	1 (25%)	6,6,6	0.47	0
3	5GP	C	606	-	22,22,26	0.48	0	33,33,40	2.81	3 (9%)

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
3	5GP	B	605	-	-	3/10/26/26	0/3/3/3
5	POP	C	608	4	-	0/6/6/6	-
3	5GP	A	604	-	-	0/10/26/26	0/3/3/3
3	5GP	C	606	-	-	0/6/22/26	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	605	5GP	P-O5'	-3.67	1.48	1.60
3	A	604	5GP	P-O5'	-3.18	1.50	1.60
2	A	600	PO4	P-O3	2.32	1.61	1.54
2	B	601	PO4	P-O4	2.28	1.61	1.54
2	B	601	PO4	P-O3	2.04	1.60	1.54

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	605	5GP	OP2-P-OP1	-13.39	58.67	110.83
3	C	606	5GP	C1'-N9-C8	11.10	158.26	126.73
3	C	606	5GP	C1'-N9-C4	-10.83	94.49	126.49
3	B	605	5GP	OP2-P-OP3	10.04	145.45	107.80
5	C	608	POP	O6-P2-O4	9.37	147.35	110.83
5	C	608	POP	O6-P2-O5	-8.86	74.59	107.80
3	A	604	5GP	OP2-P-OP3	-4.45	91.11	107.80
3	A	604	5GP	OP2-P-OP1	3.84	125.78	110.83
5	C	608	POP	O2-P1-O1	2.96	122.35	110.83
3	A	604	5GP	C1'-N9-C8	2.74	134.52	126.73
3	B	605	5GP	O4'-C4'-C3'	2.62	110.36	105.15
3	A	604	5GP	C1'-N9-C4	-2.59	118.83	126.49
3	B	605	5GP	C3'-C2'-C1'	2.17	105.56	101.46
3	C	606	5GP	C2'-C1'-N9	-2.04	107.58	113.25

There are no chirality outliers.

All (3) torsion outliers are listed below:

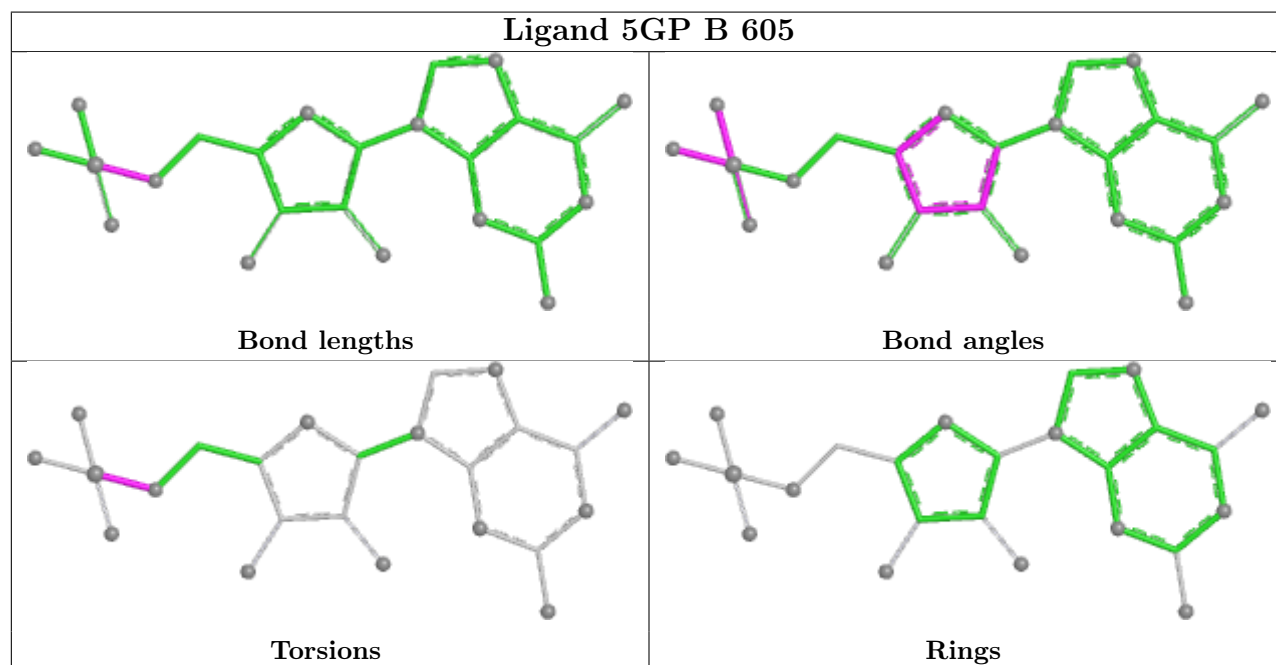
Mol	Chain	Res	Type	Atoms
3	B	605	5GP	C5'-O5'-P-OP3
3	B	605	5GP	C5'-O5'-P-OP1
3	B	605	5GP	C5'-O5'-P-OP2

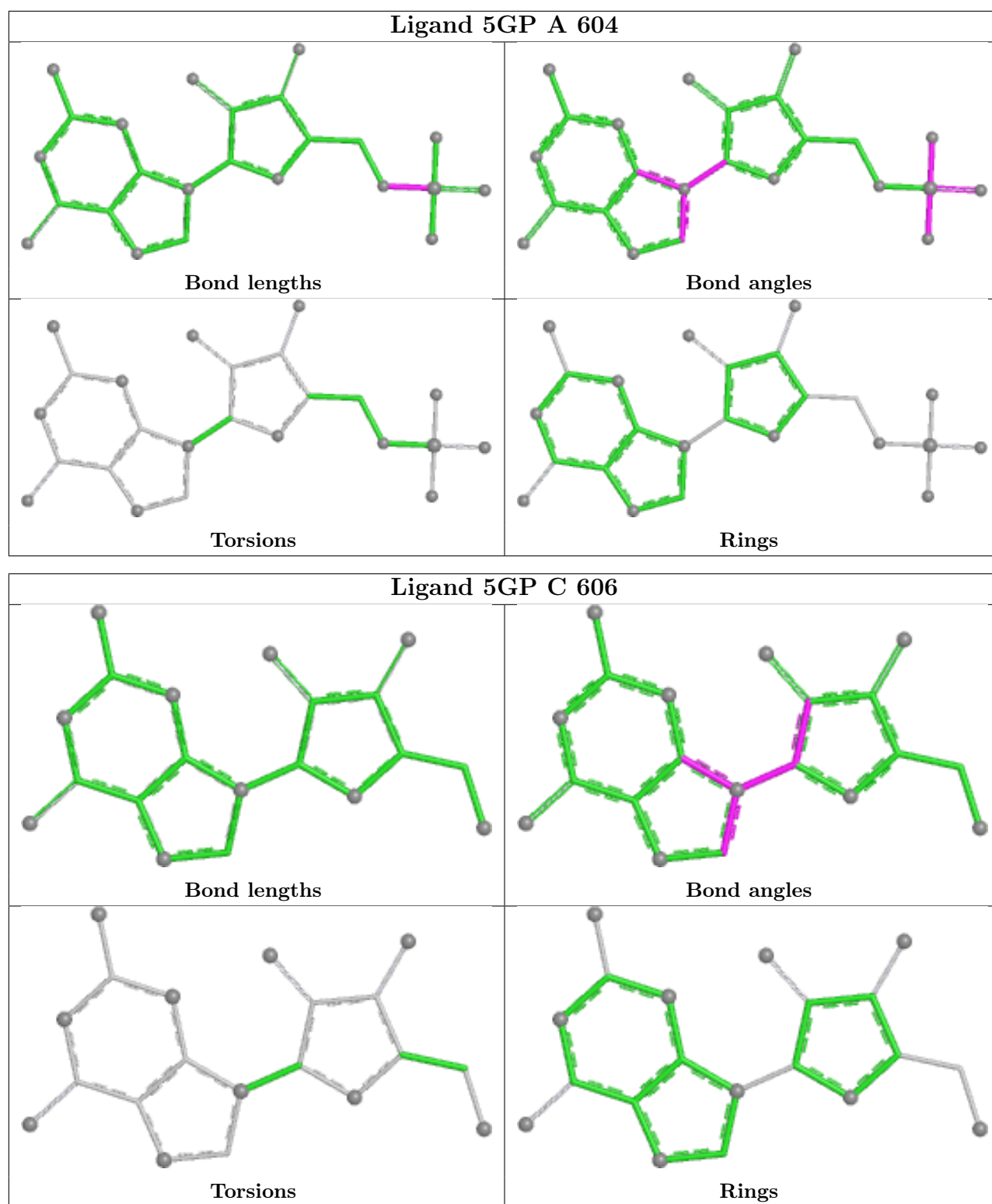
There are no ring outliers.

5 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	605	5GP	1	0
2	B	601	PO4	3	0
3	A	604	5GP	4	0
2	A	600	PO4	2	0
3	C	606	5GP	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 ⓘ

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, 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	170/180 (94%)	0.00	2 (1%) 76 74	16, 36, 66, 97	0
1	B	180/180 (100%)	-0.03	3 (1%) 69 66	13, 36, 78, 98	0
1	C	165/180 (91%)	0.03	7 (4%) 40 37	14, 38, 80, 99	0
All	All	515/540 (95%)	0.00	12 (2%) 61 58	13, 37, 78, 99	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	101	GLN	3.0
1	B	41	ALA	2.6
1	C	35	LEU	2.6
1	C	62	HIS	2.5
1	C	45	HIS	2.4
1	B	106	ALA	2.3
1	A	41	ALA	2.2
1	C	48	ASP	2.1
1	A	48	ASP	2.0
1	C	94	GLY	2.0
1	B	35	LEU	2.0
1	C	34	ILE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

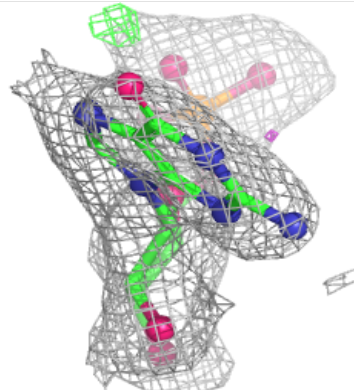
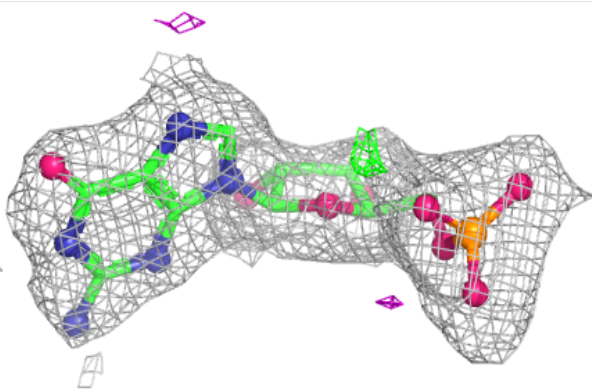
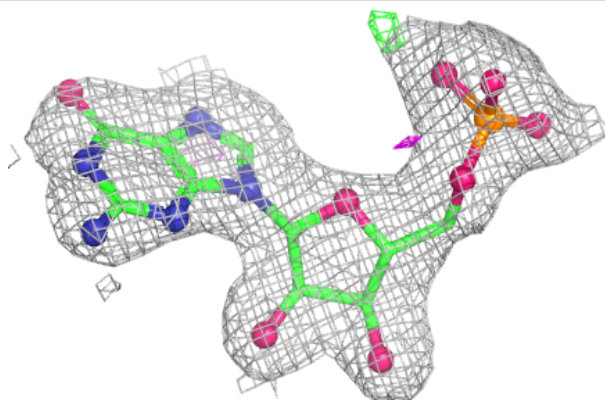
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	5GP	A	604	24/24	0.92	0.10	23,56,98,100	0
3	5GP	C	606	20/24	0.93	0.11	19,51,100,100	0
4	MG	C	607	1/1	0.93	0.16	41,41,41,41	1
2	PO4	B	601	5/5	0.94	0.14	29,31,63,94	0
2	PO4	A	600	5/5	0.94	0.14	43,44,57,98	0
3	5GP	B	605	24/24	0.95	0.08	17,30,97,100	0
5	POP	C	608	9/9	0.96	0.10	6,13,51,71	9

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.

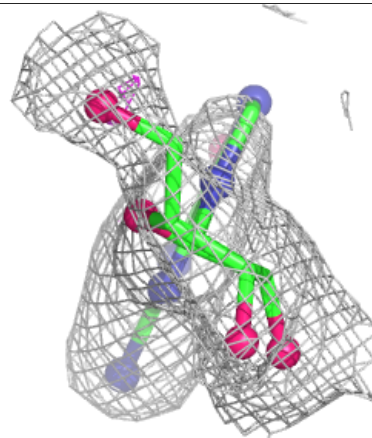
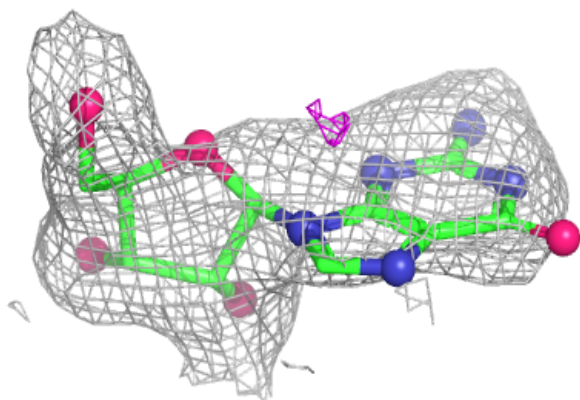
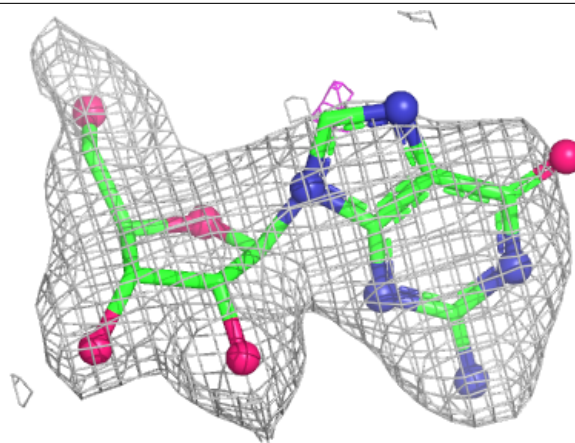
Electron density around 5GP A 604:

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

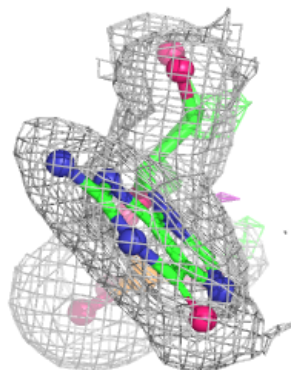
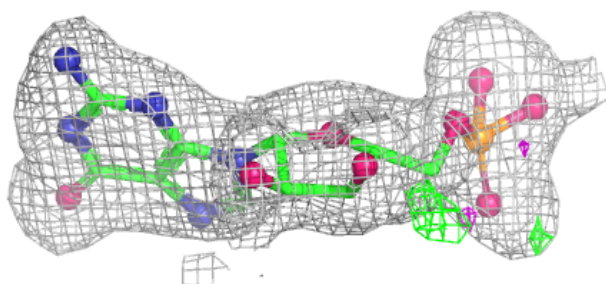


Electron density around 5GP C 606:

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

**Electron density around 5GP B 605:**

$2mF_o-DF_c$ (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.