



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 08:31 PM UTC

PDB ID : 9J0G / pdb_00009j0g
Title : Crystal structure of RhoA-TP1001 complex
Authors : Zhu, L.; Li, H.; Chang, L.; Hu, X.
Deposited on : 2024-08-02
Resolution : 3.10 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

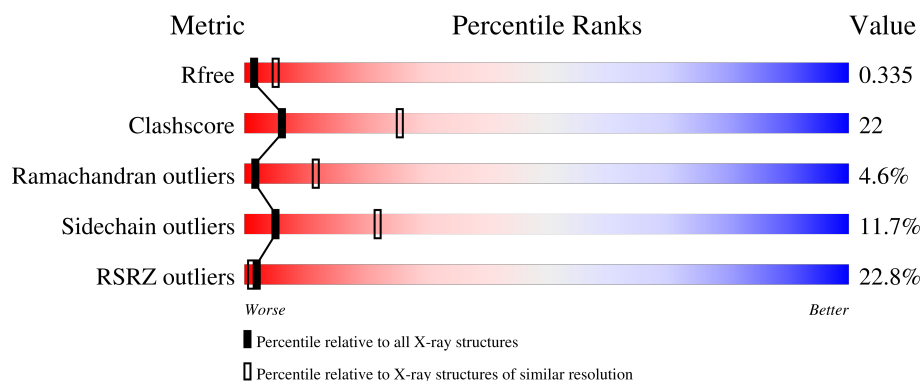
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	193	23% 44% 39% 6% 8%
1	B	193	22% 40% 45% 6% 8%
1	C	193	19% 49% 36% 10%
1	D	193	20% 39% 42% 10% 8%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5742 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.

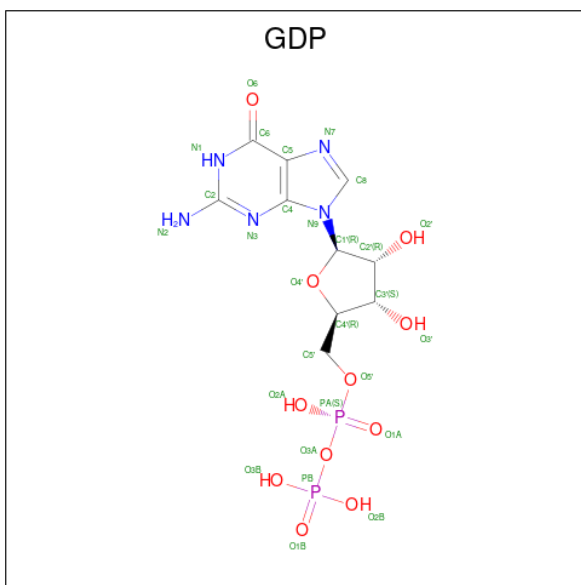
- Molecule 1 is a protein called Transforming protein RhoA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	178	Total	C	N	O	S	0	0	0
			1402	885	239	268	10			
1	B	178	Total	C	N	O	S	0	0	0
			1407	888	240	269	10			
1	C	174	Total	C	N	O	S	0	0	0
			1379	872	235	262	10			
1	D	177	Total	C	N	O	S	0	0	0
			1405	886	239	270	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	25	ASN	PHE	conflict	UNP P61586
B	25	ASN	PHE	conflict	UNP P61586
C	25	ASN	PHE	conflict	UNP P61586
D	25	ASN	PHE	conflict	UNP P61586

- Molecule 2 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).

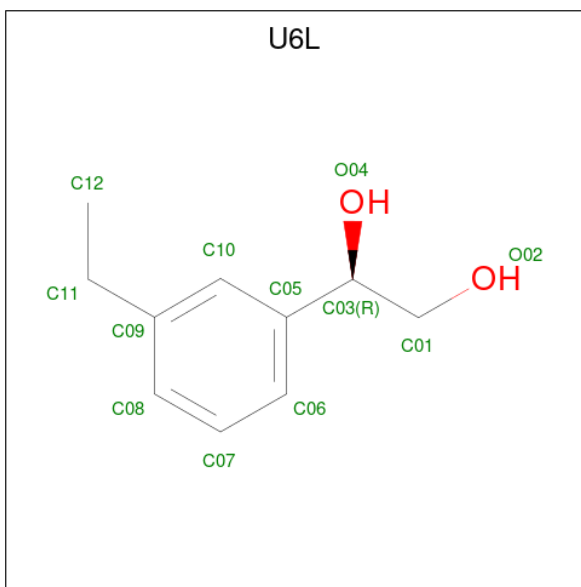


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 28	C 10	N 5	O 11	P 2	0	0
2	B	1	Total 28	C 10	N 5	O 11	P 2	0	0
2	C	1	Total 28	C 10	N 5	O 11	P 2	0	0
2	D	1	Total 28	C 10	N 5	O 11	P 2	0	0

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

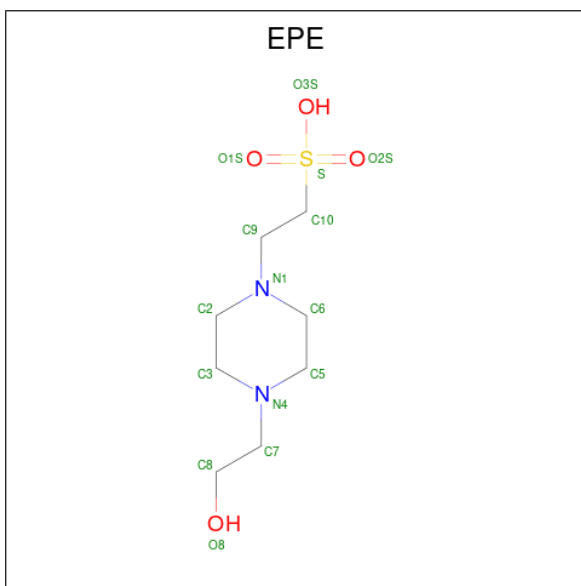
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mg 1 1	0	0
3	B	2	Total Mg 2 2	0	0
3	C	1	Total Mg 1 1	0	0
3	D	2	Total Mg 2 2	0	0

- Molecule 4 is (1 {R})-1-(3-ethylphenyl)ethane-1,2-diol (CCD ID: U6L) (formula: C₁₀H₁₄O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	C	O	0	0
			12	10	2		

- Molecule 5 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	D	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

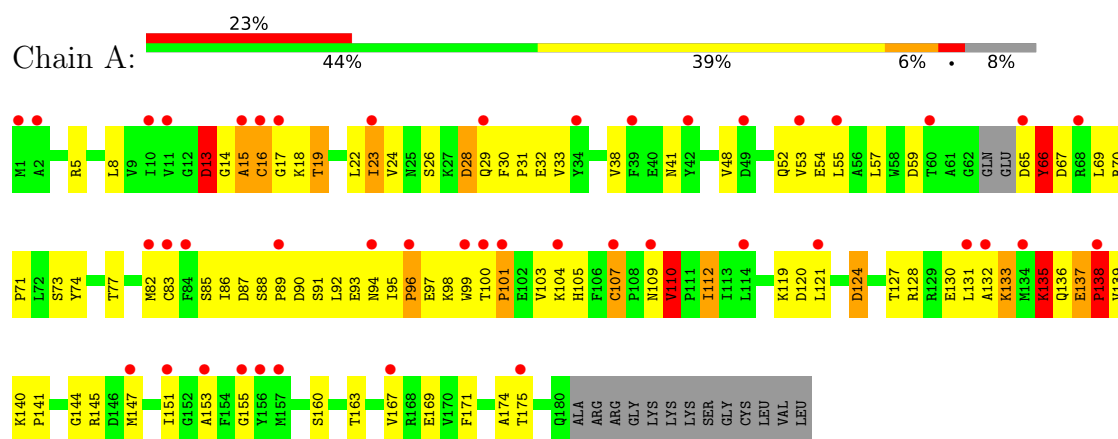
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	3	Total 3	O 3	0	0
6	B	1	Total 1	O 1	0	0

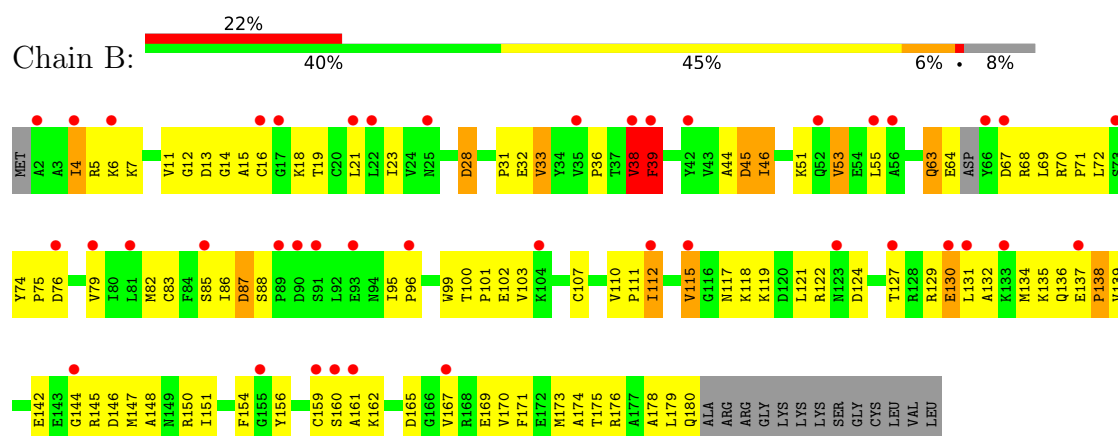
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

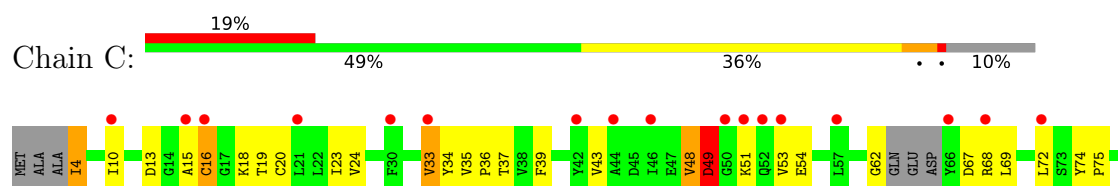
• Molecule 1: Transforming protein RhoA

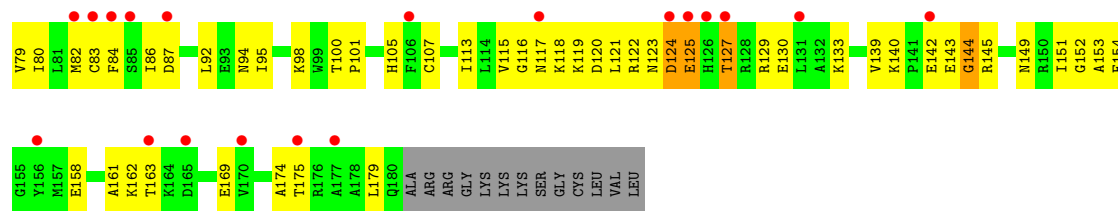


• Molecule 1: Transforming protein RhoA

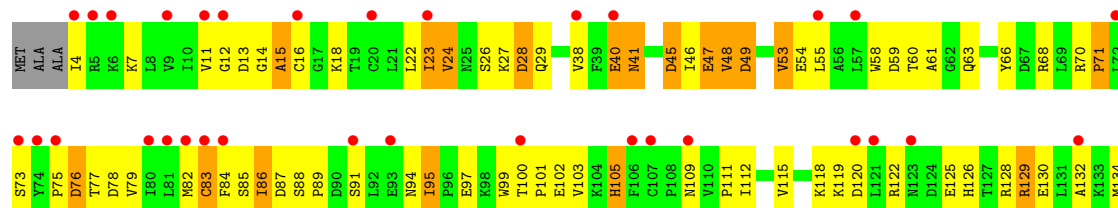


• Molecule 1: Transforming protein RhoA





● Molecule 1: Transforming protein RhoA



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	81.24Å 124.80Å 173.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.08 – 3.10 44.08 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (44.08-3.10) 99.9 (44.08-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.52 (at 3.12Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.311 , 0.354 0.314 , 0.335	Depositor DCC
R_{free} test set	872 reflections (5.32%)	wwPDB-VP
Wilson B-factor (Å ²)	39.5	Xtriage
Anisotropy	0.023	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 75.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.80	EDS
Total number of atoms	5742	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 21.54 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.8581e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: EPE, U6L, MG, GDP

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	0.54	0/1428	1.27	11/1931 (0.6%)
1	B	0.59	0/1433	1.19	4/1937 (0.2%)
1	C	0.54	0/1405	1.14	2/1899 (0.1%)
1	D	0.57	0/1432	1.24	6/1937 (0.3%)
All	All	0.56	0/5698	1.21	23/7704 (0.3%)

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	41	ASN	N-CA-CB	13.02	131.85	110.37
1	D	41	ASN	N-CA-C	-9.14	92.57	108.20
1	A	16	CYS	N-CA-C	-8.75	103.11	113.88
1	A	83	CYS	N-CA-C	8.69	123.58	109.59
1	A	31	PRO	N-CA-CB	7.15	106.92	102.92
1	D	105	HIS	CB-CA-C	5.92	118.62	109.03
1	A	124	ASP	CA-CB-CG	5.81	118.41	112.60
1	A	13	ASP	CA-CB-CG	5.75	118.35	112.60
1	D	40	GLU	N-CA-CB	-5.72	100.83	110.49
1	D	28	ASP	CA-CB-CG	5.62	118.22	112.60
1	A	96	PRO	N-CA-C	5.61	120.91	113.40
1	D	40	GLU	N-CA-C	5.56	122.64	110.80
1	B	45	ASP	CA-CB-CG	5.46	118.06	112.60
1	B	28	ASP	CA-CB-CG	5.45	118.05	112.60
1	B	67	ASP	CA-CB-CG	5.43	118.03	112.60
1	A	28	ASP	CB-CA-C	5.42	119.00	111.63
1	A	59	ASP	CA-CB-CG	5.42	118.02	112.60
1	B	39	PHE	CA-CB-CG	5.37	119.17	113.80
1	A	67	ASP	CA-CB-CG	5.36	117.96	112.60
1	C	175	THR	CA-CB-OG1	-5.27	101.69	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	37	THR	CA-CB-OG1	-5.25	101.72	109.60
1	A	66	TYR	N-CA-CB	5.13	119.16	110.49
1	A	32	GLU	CB-CA-C	5.01	118.01	109.24

There are no chirality outliers.

There are no planarity outliers.

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	1402	0	1397	60	0
1	B	1407	0	1401	68	0
1	C	1379	0	1377	54	0
1	D	1405	0	1397	72	0
2	A	28	0	12	2	0
2	B	28	0	12	0	0
2	C	28	0	12	3	0
2	D	28	0	12	4	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
3	C	1	0	0	0	0
3	D	2	0	0	0	0
4	D	12	0	0	0	0
5	D	15	0	18	3	0
6	A	3	0	0	0	0
6	B	1	0	0	0	0
All	All	5742	0	5638	255	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 22.

All (255) 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:D:38:VAL:HG12	1:D:40:GLU:HG3	1.28	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:38:VAL:HG12	1:D:40:GLU:CG	1.90	1.01
1:C:20:CYS:O	1:C:24:VAL:HG23	1.69	0.91
1:B:11:VAL:O	1:B:18:LYS:HD3	1.72	0.89
1:D:38:VAL:CG1	1:D:40:GLU:CG	2.54	0.85
1:A:16:CYS:HA	1:A:85:SER:HB3	1.57	0.84
1:D:38:VAL:CG1	1:D:40:GLU:HG2	2.10	0.82
1:A:128:ARG:O	1:A:132:ALA:HB3	1.81	0.80
1:D:13:ASP:N	1:D:16:CYS:SG	2.55	0.80
1:D:83:CYS:HB3	1:D:115:VAL:HB	1.61	0.79
1:D:84:PHE:HB2	1:D:91:SER:OG	1.86	0.76
1:A:132:ALA:HA	1:A:135:LYS:HA	1.67	0.75
1:C:169:GLU:OE1	1:C:169:GLU:N	2.20	0.75
1:B:21:LEU:HD12	1:B:167:VAL:HG13	1.72	0.71
1:D:137:GLU:HG2	1:D:138:PRO:HD2	1.73	0.70
1:B:13:ASP:OD1	1:B:99:TRP:NE1	2.24	0.70
1:A:109:ASN:O	1:A:110:VAL:HG12	1.93	0.69
1:A:70:ARG:N	1:A:71:PRO:HD2	2.07	0.69
1:C:122:ARG:NH1	1:C:139:VAL:O	2.26	0.69
1:D:119:LYS:HD3	1:D:158:GLU:HB3	1.76	0.68
1:B:7:LYS:HD3	1:B:76:ASP:HB3	1.76	0.67
1:C:79:VAL:HG21	1:C:174:ALA:HB1	1.76	0.67
1:D:100:THR:N	1:D:101:PRO:HD2	2.10	0.67
1:B:21:LEU:HD22	1:B:117:ASN:HD21	1.60	0.66
1:B:23:ILE:HG21	1:B:31:PRO:HB3	1.79	0.65
1:A:13:ASP:HB3	1:A:99:TRP:CH2	2.31	0.65
1:A:140:LYS:HG3	1:A:141:PRO:HD2	1.78	0.64
1:B:79:VAL:HB	1:B:178:ALA:HB2	1.79	0.64
1:B:103:VAL:HG21	1:B:112:ILE:HD11	1.80	0.64
1:A:24:VAL:HG22	1:A:30:PHE:HA	1.80	0.63
1:B:70:ARG:N	1:B:71:PRO:HD2	2.13	0.62
1:B:13:ASP:O	1:B:16:CYS:SG	2.51	0.62
1:B:148:ALA:O	1:B:151:ILE:HG13	1.99	0.62
1:D:97:GLU:O	1:D:101:PRO:HG3	2.00	0.61
1:D:58:TRP:CH2	5:D:205:EPE:H52	2.36	0.61
1:C:162:LYS:HD3	1:D:120:ASP:OD2	2.01	0.61
1:B:83:CYS:SG	1:B:115:VAL:HG12	2.41	0.60
1:B:129:ARG:O	1:B:131:LEU:N	2.34	0.60
1:A:103:VAL:O	1:A:107:CYS:N	2.34	0.60
1:B:118:LYS:HB3	1:B:121:LEU:HD13	1.81	0.60
1:D:11:VAL:HG22	1:D:82:MET:HA	1.82	0.59
1:B:4:ILE:O	1:B:5:ARG:HB2	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:147:MET:O	1:B:148:ALA:C	2.46	0.59
1:D:140:LYS:O	1:D:143:GLU:HB2	2.03	0.59
1:C:94:ASN:O	1:C:98:LYS:HB3	2.03	0.58
1:D:141:PRO:O	1:D:144:GLY:N	2.33	0.58
1:C:74:TYR:N	1:C:75:PRO:CD	2.66	0.58
1:D:125:GLU:O	1:D:126:HIS:C	2.45	0.58
1:B:160:SER:C	1:B:162:LYS:H	2.12	0.58
1:D:23:ILE:O	1:D:26:SER:N	2.36	0.58
1:A:65:ASP:O	1:A:66:TYR:CG	2.56	0.57
1:A:95:ILE:HG22	1:A:147:MET:HE2	1.85	0.57
1:A:55:LEU:HD23	1:A:57:LEU:HD21	1.87	0.57
1:A:140:LYS:HG3	1:A:141:PRO:CD	2.35	0.57
1:A:141:PRO:HB3	1:A:145:ARG:HH11	1.70	0.57
1:B:100:THR:HB	1:B:101:PRO:HD3	1.86	0.57
1:D:45:ASP:O	1:D:47:GLU:HG2	2.05	0.57
1:A:94:ASN:O	1:A:97:GLU:O	2.23	0.56
1:C:125:GLU:O	1:C:129:ARG:HD2	2.04	0.56
1:A:107:CYS:SG	1:A:110:VAL:HG11	2.45	0.56
1:B:71:PRO:HA	1:B:74:TYR:CD2	2.40	0.56
1:D:22:LEU:O	1:D:23:ILE:C	2.47	0.56
1:B:44:ALA:HB3	1:B:55:LEU:HB2	1.85	0.56
1:B:96:PRO:HD3	1:B:147:MET:SD	2.46	0.56
1:A:130:GLU:HG3	1:A:133:LYS:HD2	1.88	0.55
1:C:125:GLU:HA	1:C:129:ARG:NH1	2.22	0.55
1:D:160:SER:HB3	1:D:165:ASP:HB2	1.89	0.55
1:C:120:ASP:OD1	1:C:121:LEU:HG	2.06	0.55
1:D:59:ASP:OD1	1:D:60:THR:N	2.39	0.54
1:B:36:PRO:C	1:B:38:VAL:H	2.15	0.54
1:D:38:VAL:HG11	1:D:40:GLU:HG2	1.89	0.54
1:D:111:PRO:CB	1:D:177:ALA:HB1	2.38	0.54
1:B:117:ASN:HA	1:B:159:CYS:O	2.07	0.54
1:C:142:GLU:C	1:C:144:GLY:H	2.16	0.54
1:C:151:ILE:HD11	1:C:153:ALA:HB2	1.88	0.54
1:A:18:LYS:HB2	2:A:201:GDP:O1B	2.07	0.54
1:A:24:VAL:O	1:A:28:ASP:N	2.34	0.54
1:C:84:PHE:CE2	1:C:115:VAL:O	2.61	0.54
1:A:144:GLY:O	1:A:147:MET:N	2.39	0.54
1:A:100:THR:HB	1:A:101:PRO:HD3	1.89	0.53
1:A:160:SER:HB3	1:A:163:THR:OG1	2.08	0.53
1:B:72:LEU:O	1:B:75:PRO:HD2	2.08	0.53
1:C:151:ILE:HG13	1:C:152:GLY:N	2.22	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:14:GLY:HA2	2:D:202:GDP:O2B	2.09	0.53
1:A:86:ILE:HG23	1:A:139:VAL:HB	1.90	0.53
1:D:176:ARG:O	1:D:180:GLN:HG3	2.08	0.53
1:A:23:ILE:O	1:A:24:VAL:C	2.48	0.53
1:B:129:ARG:C	1:B:131:LEU:H	2.17	0.53
1:C:74:TYR:N	1:C:75:PRO:HD3	2.24	0.53
1:A:103:VAL:C	1:A:105:HIS:H	2.16	0.53
1:B:169:GLU:O	1:B:173:MET:HG2	2.08	0.53
1:D:53:VAL:HG11	1:D:179:LEU:HD21	1.90	0.53
1:A:92:LEU:O	1:A:96:PRO:HD2	2.08	0.53
1:B:145:ARG:O	1:B:146:ASP:C	2.48	0.52
1:D:88:SER:O	1:D:91:SER:HB3	2.09	0.52
1:C:23:ILE:CD1	1:C:36:PRO:HD2	2.40	0.52
1:C:24:VAL:HG21	1:C:161:ALA:HB1	1.90	0.52
1:D:12:GLY:O	1:D:99:TRP:CZ2	2.63	0.52
1:D:85:SER:O	1:D:88:SER:N	2.43	0.52
1:D:86:ILE:HG21	1:D:122:ARG:HB2	1.92	0.52
1:D:157:MET:HE3	1:D:173:MET:SD	2.49	0.52
1:B:179:LEU:O	1:B:180:GLN:C	2.52	0.52
1:C:16:CYS:SG	1:C:117:ASN:ND2	2.83	0.52
2:C:201:GDP:O2A	2:C:201:GDP:O3'	2.23	0.52
1:D:27:LYS:C	1:D:29:GLN:H	2.18	0.51
1:D:78:ASP:HB3	1:D:178:ALA:HB1	1.92	0.51
1:A:71:PRO:HA	1:A:74:TYR:HD2	1.75	0.51
1:A:121:LEU:HB3	1:A:127:THR:HG21	1.92	0.51
1:C:123:ASN:O	1:C:124:ASP:C	2.53	0.51
1:A:53:VAL:HG21	1:A:175:THR:HG23	1.92	0.51
1:A:128:ARG:O	1:A:132:ALA:CB	2.55	0.51
1:A:141:PRO:HB3	1:A:145:ARG:NH1	2.25	0.51
1:C:48:VAL:O	1:C:49:ASP:C	2.53	0.51
1:C:62:GLY:HA2	1:C:67:ASP:OD1	2.10	0.51
1:D:27:LYS:C	1:D:29:GLN:N	2.68	0.51
1:A:70:ARG:HG2	1:A:74:TYR:OH	2.11	0.51
1:A:85:SER:HB2	1:A:87:ASP:OD1	2.10	0.51
1:A:14:GLY:O	1:A:15:ALA:HB3	2.11	0.51
1:A:71:PRO:HA	1:A:74:TYR:CD2	2.46	0.51
1:B:170:VAL:HG23	1:B:171:PHE:CD1	2.46	0.51
1:C:19:THR:OG1	2:C:201:GDP:O1B	2.29	0.51
1:D:140:LYS:N	1:D:143:GLU:OE1	2.26	0.50
1:D:141:PRO:O	1:D:142:GLU:C	2.53	0.50
1:D:162:LYS:N	2:D:202:GDP:O6	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:24:VAL:HG11	1:D:167:VAL:HG11	1.94	0.50
1:B:55:LEU:HD21	1:B:175:THR:HG23	1.93	0.50
1:C:100:THR:N	1:C:101:PRO:HD2	2.27	0.50
1:A:5:ARG:NH2	1:A:54:GLU:OE1	2.46	0.49
1:B:4:ILE:HD12	1:B:53:VAL:HG22	1.94	0.49
1:D:27:LYS:O	1:D:29:GLN:N	2.45	0.49
1:D:128:ARG:HG2	1:D:138:PRO:HG3	1.93	0.49
1:A:169:GLU:CD	1:A:169:GLU:N	2.70	0.49
1:D:75:PRO:O	1:D:76:ASP:HB2	2.13	0.49
1:D:122:ARG:NE	1:D:139:VAL:O	2.22	0.49
1:B:19:THR:HG21	1:B:36:PRO:O	2.13	0.48
1:D:18:LYS:HG3	2:D:202:GDP:O3B	2.13	0.48
1:B:95:ILE:HG22	1:B:151:ILE:HG21	1.95	0.48
1:C:20:CYS:SG	1:C:35:VAL:HG11	2.54	0.48
1:D:66:TYR:C	1:D:68:ARG:H	2.20	0.48
1:B:11:VAL:HG23	1:B:12:GLY:N	2.28	0.48
1:C:20:CYS:C	1:C:24:VAL:HG23	2.37	0.48
1:B:129:ARG:C	1:B:131:LEU:N	2.72	0.48
1:C:82:MET:O	1:C:115:VAL:HG23	2.14	0.48
1:B:45:ASP:C	1:B:46:ILE:HG12	2.39	0.47
1:D:100:THR:N	1:D:101:PRO:CD	2.77	0.47
1:C:87:ASP:OD1	1:C:127:THR:HG23	2.14	0.47
1:A:144:GLY:O	1:A:145:ARG:C	2.56	0.47
1:A:103:VAL:O	1:A:105:HIS:N	2.48	0.47
1:D:91:SER:O	1:D:94:ASN:N	2.48	0.47
1:A:89:PRO:HB3	1:A:139:VAL:HG22	1.97	0.47
1:B:159:CYS:HB2	1:B:165:ASP:O	2.14	0.47
1:D:101:PRO:O	1:D:102:GLU:C	2.57	0.47
1:A:153:ALA:C	1:A:155:GLY:N	2.71	0.47
1:C:10:ILE:HD13	1:C:18:LYS:HB3	1.96	0.47
1:A:88:SER:O	1:A:91:SER:OG	2.31	0.46
1:A:94:ASN:O	1:A:98:LYS:O	2.32	0.46
1:A:8:LEU:HD23	1:A:8:LEU:C	2.40	0.46
1:D:125:GLU:O	1:D:129:ARG:N	2.37	0.46
1:B:14:GLY:O	1:B:15:ALA:C	2.58	0.46
1:B:95:ILE:HB	1:B:96:PRO:CD	2.45	0.46
1:C:82:MET:CG	1:C:95:ILE:HD12	2.45	0.46
1:C:153:ALA:O	1:C:154:PHE:C	2.59	0.46
1:A:147:MET:HE3	1:A:151:ILE:HG23	1.98	0.46
1:A:124:ASP:OD2	1:A:127:THR:HG23	2.16	0.46
1:C:15:ALA:O	1:C:16:CYS:C	2.58	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:160:SER:C	1:B:162:LYS:N	2.74	0.46
1:C:101:PRO:O	1:C:105:HIS:N	2.49	0.46
1:D:60:THR:HG22	1:D:70:ARG:HG3	1.97	0.46
1:A:70:ARG:N	1:A:71:PRO:CD	2.77	0.46
1:B:173:MET:O	1:B:174:ALA:C	2.59	0.45
1:D:166:GLY:HA2	1:D:169:GLU:OE1	2.15	0.45
1:D:171:PHE:HA	1:D:174:ALA:HB3	1.97	0.45
1:B:32:GLU:O	1:B:33:VAL:HB	2.17	0.45
1:B:79:VAL:HG21	1:B:174:ALA:O	2.16	0.45
1:B:103:VAL:O	1:B:107:CYS:N	2.46	0.45
1:B:129:ARG:O	1:B:132:ALA:N	2.48	0.45
1:D:85:SER:O	1:D:87:ASP:N	2.50	0.45
1:A:98:LYS:C	1:A:101:PRO:HD2	2.41	0.45
1:B:144:GLY:HA3	1:B:156:TYR:CZ	2.51	0.45
1:D:46:ILE:HD13	1:D:55:LEU:HD23	1.99	0.45
1:A:136:GLN:O	1:A:137:GLU:C	2.60	0.44
1:B:83:CYS:SG	1:B:115:VAL:CG1	3.04	0.44
1:B:121:LEU:HA	1:B:124:ASP:HB2	1.99	0.44
1:C:95:ILE:HD13	1:C:95:ILE:HA	1.87	0.44
1:D:7:LYS:HG2	1:D:77:THR:HA	1.99	0.44
1:C:124:ASP:HB3	1:C:127:THR:HB	1.98	0.44
1:D:23:ILE:O	1:D:24:VAL:C	2.60	0.44
1:D:118:LYS:HE2	2:D:202:GDP:C4	2.53	0.44
1:A:19:THR:OG1	2:A:201:GDP:O1B	2.35	0.44
1:A:167:VAL:O	1:A:171:PHE:HD1	2.00	0.44
1:B:13:ASP:O	1:B:18:LYS:NZ	2.50	0.44
1:A:16:CYS:SG	1:A:18:LYS:HG3	2.58	0.44
1:B:119:LYS:C	1:B:121:LEU:H	2.26	0.44
1:C:82:MET:HG2	1:C:95:ILE:HD12	2.00	0.44
1:B:4:ILE:HB	1:B:53:VAL:HG13	1.98	0.43
1:D:132:ALA:C	1:D:134:MET:H	2.26	0.43
1:B:117:ASN:OD1	1:B:159:CYS:SG	2.76	0.43
1:C:79:VAL:HG22	1:C:80:ILE:N	2.33	0.43
1:A:70:ARG:HG2	1:A:74:TYR:CZ	2.54	0.43
1:B:122:ARG:NH1	1:B:139:VAL:O	2.52	0.43
1:C:19:THR:CB	2:C:201:GDP:O1B	2.67	0.43
1:B:11:VAL:HG22	1:B:82:MET:HA	1.99	0.43
1:B:21:LEU:HD11	1:B:170:VAL:HG21	1.99	0.43
1:C:145:ARG:O	1:C:149:ASN:ND2	2.46	0.43
1:C:4:ILE:HG22	1:C:51:LYS:HE3	2.01	0.43
1:C:53:VAL:HG21	1:C:179:LEU:HD11	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:140:LYS:C	1:C:142:GLU:H	2.25	0.43
1:C:142:GLU:C	1:C:144:GLY:N	2.77	0.43
1:A:90:ASP:O	1:A:93:GLU:HB3	2.19	0.42
1:A:103:VAL:C	1:A:105:HIS:N	2.77	0.42
1:B:7:LYS:HB3	1:B:76:ASP:O	2.19	0.42
1:C:4:ILE:HA	1:C:53:VAL:HA	2.00	0.42
1:A:82:MET:HE3	1:A:112:ILE:HG21	2.01	0.42
1:C:92:LEU:HD22	1:C:139:VAL:HG11	2.01	0.42
1:C:16:CYS:SG	1:C:18:LYS:HG2	2.59	0.42
1:C:82:MET:SD	1:C:95:ILE:HG23	2.58	0.42
1:B:85:SER:HB2	1:B:88:SER:OG	2.20	0.42
1:D:41:ASN:HA	5:D:205:EPE:O2S	2.20	0.42
1:D:48:VAL:O	1:D:49:ASP:C	2.61	0.42
1:B:6:LYS:HE2	1:B:179:LEU:HD23	2.02	0.42
1:C:33:VAL:O	1:C:34:TYR:C	2.62	0.42
1:C:145:ARG:HE	1:C:145:ARG:HB2	1.65	0.41
1:D:61:ALA:HB3	1:D:66:TYR:CG	2.54	0.41
1:D:89:PRO:HB2	1:D:136:GLN:OE1	2.20	0.41
1:B:87:ASP:O	1:B:131:LEU:HD11	2.20	0.41
1:B:111:PRO:HG3	1:B:180:GLN:OE1	2.20	0.41
1:D:129:ARG:HG2	1:D:130:GLU:N	2.36	0.41
1:D:7:LYS:N	1:D:78:ASP:OD2	2.39	0.41
1:D:53:VAL:HG11	1:D:179:LEU:CD2	2.50	0.41
5:D:205:EPE:H101	5:D:205:EPE:H61	1.80	0.41
1:C:130:GLU:O	1:C:133:LYS:HB2	2.20	0.41
1:C:23:ILE:HG12	1:C:39:PHE:CD1	2.55	0.41
1:B:6:LYS:HD2	1:B:178:ALA:HB1	2.03	0.41
1:B:71:PRO:HA	1:B:74:TYR:HD2	1.82	0.41
1:B:110:VAL:O	1:B:111:PRO:C	2.63	0.41
1:C:86:ILE:O	1:C:86:ILE:HG22	2.20	0.41
1:D:78:ASP:O	1:D:111:PRO:HD2	2.19	0.41
1:A:17:GLY:O	1:A:18:LYS:C	2.63	0.41
1:A:174:ALA:O	1:A:175:THR:C	2.64	0.41
1:B:64:GLU:OE1	1:B:68:ARG:HB2	2.20	0.41
1:C:16:CYS:SG	1:C:83:CYS:HB3	2.61	0.41
1:D:129:ARG:O	1:D:132:ALA:HB3	2.21	0.41
1:B:173:MET:SD	1:B:176:ARG:NH2	2.94	0.41
1:C:116:GLY:HA3	1:C:158:GLU:HG2	2.01	0.41
1:D:86:ILE:HG23	1:D:139:VAL:O	2.20	0.40
1:A:23:ILE:O	1:A:26:SER:N	2.54	0.40
1:D:95:ILE:HD13	1:D:95:ILE:HA	1.98	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:LEU:HD22	1:A:138:PRO:HG3	2.03	0.40
1:B:148:ALA:HB2	1:B:156:TYR:HB2	2.03	0.40
1:D:4:ILE:O	1:D:54:GLU:N	2.54	0.40
1:B:38:VAL:O	1:B:39:PHE:CB	2.70	0.40
1:D:61:ALA:HB3	1:D:66:TYR:CD2	2.56	0.40
1:B:70:ARG:N	1:B:71:PRO:CD	2.83	0.40
1:C:43:VAL:HG11	1:C:54:GLU:HB3	2.04	0.40
1:D:14:GLY:O	1:D:15:ALA:HB2	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	174/193 (90%)	127 (73%)	37 (21%)	10 (6%)	1	8
1	B	174/193 (90%)	131 (75%)	34 (20%)	9 (5%)	1	10
1	C	170/193 (88%)	140 (82%)	24 (14%)	6 (4%)	3	16
1	D	175/193 (91%)	130 (74%)	38 (22%)	7 (4%)	2	13
All	All	693/772 (90%)	528 (76%)	133 (19%)	32 (5%)	2	12

All (32) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	66	TYR
1	A	135	LYS
1	B	63	GLN
1	B	38	VAL
1	B	39	PHE
1	B	130	GLU
1	C	49	ASP

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Mol	Chain	Res	Type
1	C	124	ASP
1	C	125	GLU
1	C	143	GLU
1	D	15	ALA
1	D	28	ASP
1	A	15	ALA
1	A	48	VAL
1	B	135	LYS
1	B	161	ALA
1	D	23	ILE
1	D	63	GLN
1	D	71	PRO
1	A	33	VAL
1	A	104	LYS
1	A	110	VAL
1	B	102	GLU
1	D	76	ASP
1	A	138	PRO
1	D	86	ILE
1	A	137	GLU
1	A	23	ILE
1	B	33	VAL
1	C	33	VAL
1	C	144	GLY
1	B	138	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	154/167 (92%)	135 (88%)	19 (12%)	4	20
1	B	155/167 (93%)	134 (86%)	21 (14%)	4	16
1	C	153/167 (92%)	139 (91%)	14 (9%)	8	31
1	D	156/167 (93%)	138 (88%)	18 (12%)	5	23
All	All	618/668 (92%)	546 (88%)	72 (12%)	5	22

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ASP
1	A	19	THR
1	A	22	LEU
1	A	29	GLN
1	A	38	VAL
1	A	41	ASN
1	A	52	GLN
1	A	69	LEU
1	A	73	SER
1	A	77	THR
1	A	101	PRO
1	A	107	CYS
1	A	110	VAL
1	A	112	ILE
1	A	119	LYS
1	A	120	ASP
1	A	133	LYS
1	A	135	LYS
1	A	138	PRO
1	B	4	ILE
1	B	28	ASP
1	B	38	VAL
1	B	46	ILE
1	B	51	LYS
1	B	53	VAL
1	B	63	GLN
1	B	69	LEU
1	B	86	ILE
1	B	87	ASP
1	B	112	ILE
1	B	115	VAL
1	B	127	THR
1	B	130	GLU
1	B	134	MET
1	B	136	GLN
1	B	137	GLU
1	B	138	PRO
1	B	142	GLU
1	B	150	ARG
1	B	154	PHE
1	C	4	ILE
1	C	13	ASP

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Mol	Chain	Res	Type
1	C	16	CYS
1	C	48	VAL
1	C	49	ASP
1	C	68	ARG
1	C	69	LEU
1	C	72	LEU
1	C	107	CYS
1	C	113	ILE
1	C	118	LYS
1	C	119	LYS
1	C	127	THR
1	C	163	THR
1	D	24	VAL
1	D	45	ASP
1	D	47	GLU
1	D	48	VAL
1	D	49	ASP
1	D	53	VAL
1	D	71	PRO
1	D	73	SER
1	D	79	VAL
1	D	83	CYS
1	D	95	ILE
1	D	103	VAL
1	D	105	HIS
1	D	109	ASN
1	D	112	ILE
1	D	129	ARG
1	D	173	MET
1	D	175	THR

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

Mol	Chain	Res	Type
1	A	25	ASN
1	B	25	ASN
1	B	41	ASN
1	C	41	ASN
1	C	105	HIS
1	D	117	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 12 ligands modelled in this entry, 6 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$
2	GDP	C	201	3	29,30,30	0.63	0	45,47,47	0.72	0
5	EPE	D	205	-	15,15,15	0.78	1 (6%)	19,20,20	1.20	1 (5%)
2	GDP	D	202	3	29,30,30	0.50	0	45,47,47	0.66	0
4	U6L	D	201	3	11,12,12	0.44	0	12,15,15	0.27	0
2	GDP	B	301	3	29,30,30	0.48	0	45,47,47	0.64	0
2	GDP	A	201	3	29,30,30	0.54	0	45,47,47	0.77	1 (2%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GDP	C	201	3	-	6/16/32/32	0/3/3/3
5	EPE	D	205	-	-	4/9/19/19	0/1/1/1
2	GDP	D	202	3	-	7/16/32/32	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	U6L	D	201	3	-	0/8/8/8	0/1/1/1
2	GDP	B	301	3	-	2/16/32/32	0/3/3/3
2	GDP	A	201	3	-	4/16/32/32	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	205	EPE	O3S-S	2.28	1.55	1.47

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	201	GDP	C2'-C3'-C4'	-2.32	98.12	102.61
5	D	205	EPE	C9-N1-C6	2.04	116.68	111.24

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	201	GDP	C5'-O5'-PA-O3A
2	A	201	GDP	O4'-C4'-C5'-O5'
2	C	201	GDP	O4'-C4'-C5'-O5'
2	D	202	GDP	C5'-O5'-PA-O3A
2	D	202	GDP	C5'-O5'-PA-O1A
2	D	202	GDP	O4'-C4'-C5'-O5'
5	D	205	EPE	C10-C9-N1-C6
5	D	205	EPE	S-C10-C9-N1
2	A	201	GDP	C3'-C4'-C5'-O5'
2	D	202	GDP	C3'-C4'-C5'-O5'
2	C	201	GDP	C3'-C4'-C5'-O5'
2	C	201	GDP	C4'-C5'-O5'-PA
5	D	205	EPE	C10-C9-N1-C2
2	C	201	GDP	PA-O3A-PB-O1B
2	A	201	GDP	C5'-O5'-PA-O1A
2	D	202	GDP	C5'-O5'-PA-O2A
2	B	301	GDP	PB-O3A-PA-O2A
2	D	202	GDP	PB-O3A-PA-O2A
2	D	202	GDP	C4'-C5'-O5'-PA
2	C	201	GDP	PA-O3A-PB-O2B
2	C	201	GDP	PA-O3A-PB-O3B
2	B	301	GDP	PB-O3A-PA-O1A

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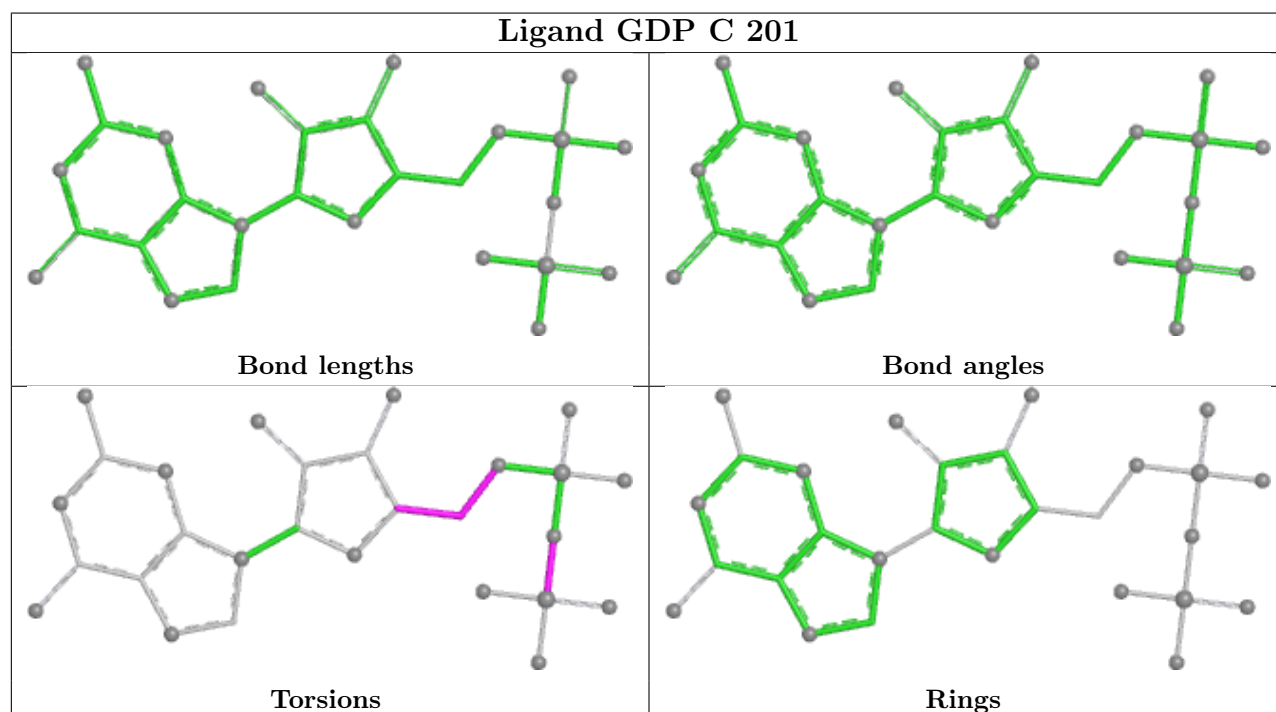
Mol	Chain	Res	Type	Atoms
5	D	205	EPE	N4-C7-C8-O8

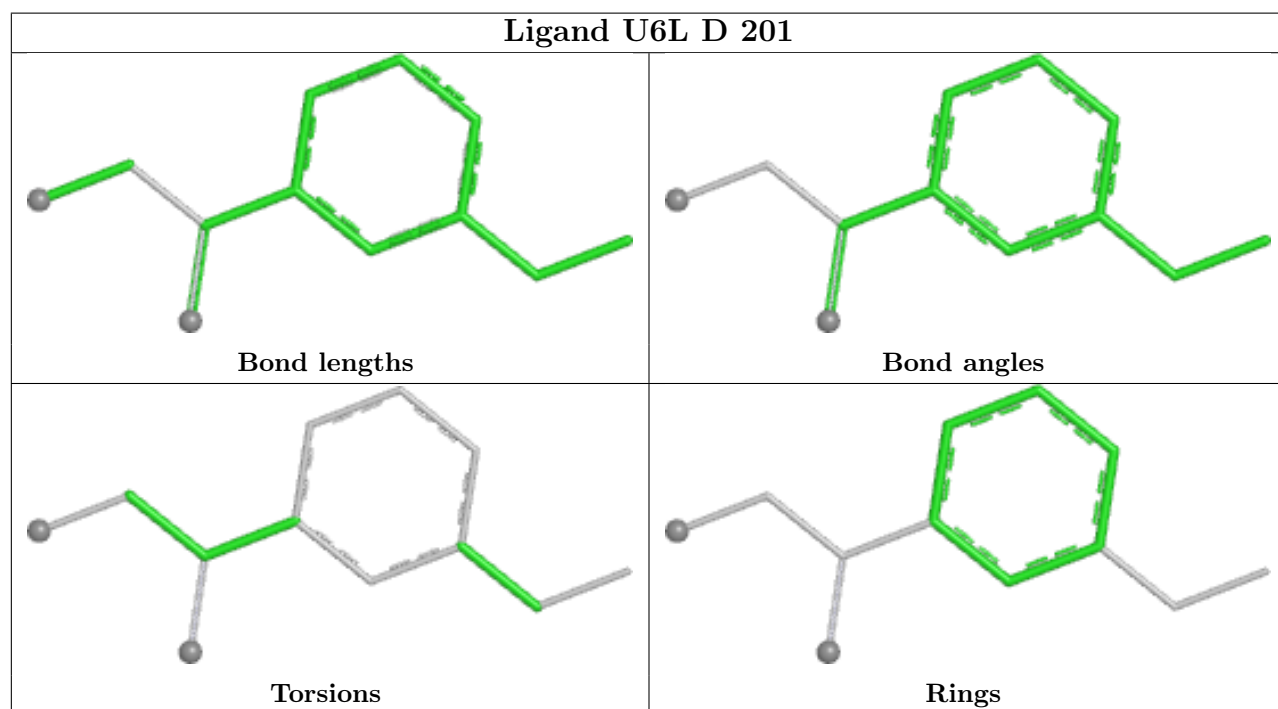
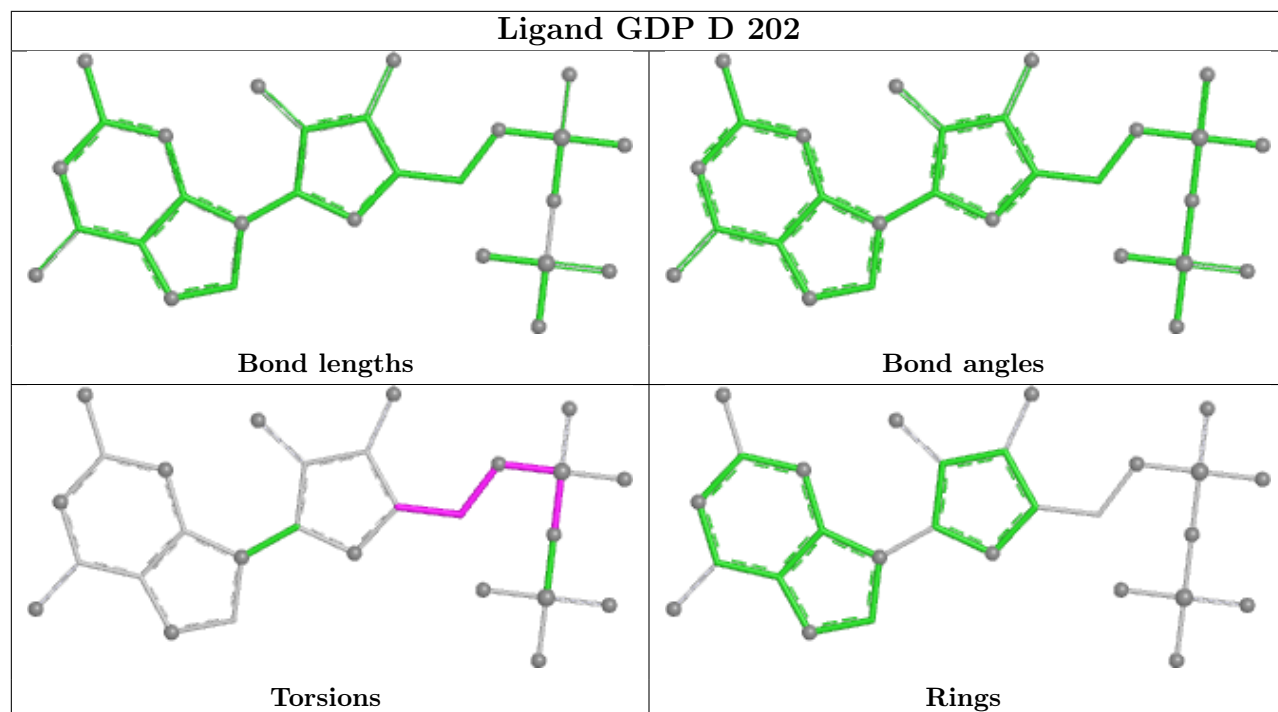
There are no ring outliers.

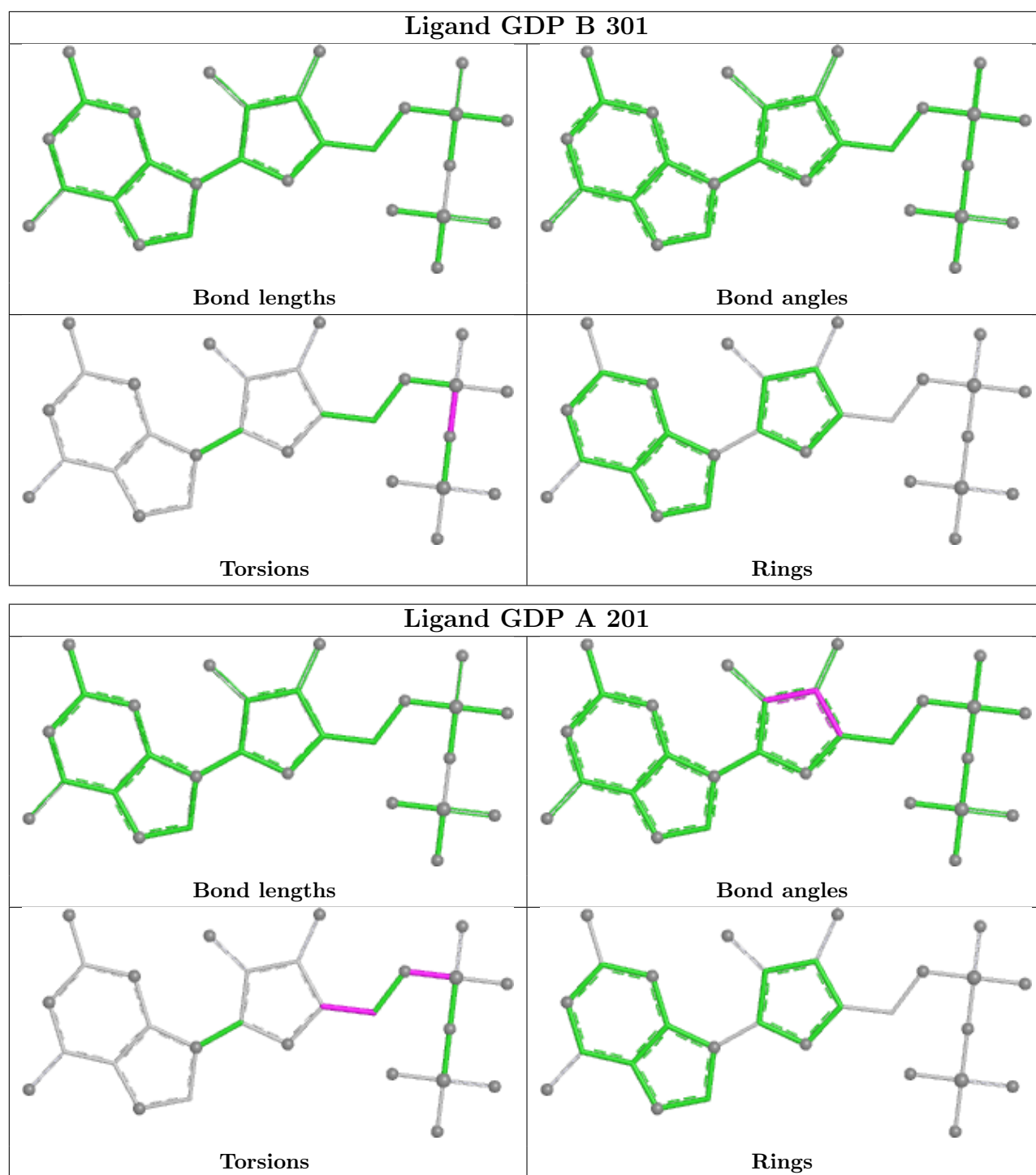
4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	201	GDP	3	0
5	D	205	EPE	3	0
2	D	202	GDP	4	0
2	A	201	GDP	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	178/193 (92%)	1.55	44 (24%) 2 1	18, 32, 55, 70	0
1	B	178/193 (92%)	1.42	42 (23%) 2 1	14, 31, 50, 59	0
1	C	174/193 (90%)	1.40	36 (20%) 2 1	11, 30, 54, 76	0
1	D	177/193 (91%)	1.33	39 (22%) 2 1	16, 33, 49, 62	0
All	All	707/772 (91%)	1.43	161 (22%) 2 1	11, 32, 53, 76	0

All (161) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	161	ALA	5.6
1	A	83	CYS	4.7
1	A	156	TYR	4.6
1	A	10	ILE	4.3
1	C	15	ALA	4.3
1	C	16	CYS	4.2
1	A	16	CYS	4.1
1	C	131	LEU	4.0
1	D	84	PHE	4.0
1	B	76	ASP	4.0
1	D	121	LEU	3.9
1	C	68	ARG	3.8
1	C	72	LEU	3.8
1	A	23	ILE	3.6
1	C	42	TYR	3.5
1	A	114	LEU	3.4
1	C	44	ALA	3.4
1	A	96	PRO	3.4
1	B	17	GLY	3.4
1	A	11	VAL	3.4
1	D	165	ASP	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	30	PHE	3.3
1	B	137	GLU	3.3
1	A	53	VAL	3.3
1	C	21	LEU	3.3
1	D	81	LEU	3.3
1	D	160	SER	3.2
1	C	126	HIS	3.2
1	A	153	ALA	3.2
1	D	174	ALA	3.2
1	A	39	PHE	3.1
1	A	15	ALA	3.1
1	A	42	TYR	3.1
1	D	57	LEU	3.1
1	A	107	CYS	3.1
1	A	89	PRO	3.1
1	A	1	MET	3.0
1	D	11	VAL	3.0
1	C	83	CYS	3.0
1	C	170	VAL	3.0
1	D	16	CYS	3.0
1	D	175	THR	2.9
1	B	66	TYR	2.9
1	A	100	THR	2.9
1	D	106	PHE	2.9
1	B	22	LEU	2.9
1	B	89	PRO	2.8
1	B	67	ASP	2.8
1	B	38	VAL	2.8
1	C	51	LYS	2.8
1	D	55	LEU	2.8
1	C	57	LEU	2.8
1	A	17	GLY	2.7
1	C	124	ASP	2.7
1	A	131	LEU	2.7
1	A	167	VAL	2.7
1	B	167	VAL	2.7
1	D	23	ILE	2.7
1	A	101	PRO	2.6
1	B	73	SER	2.6
1	A	94	ASN	2.6
1	A	2	ALA	2.6
1	C	175	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	133	LYS	2.6
1	A	155	GLY	2.6
1	B	127	THR	2.6
1	D	12	GLY	2.6
1	D	73	SER	2.5
1	D	9	VAL	2.5
1	A	132	ALA	2.5
1	A	60	THR	2.5
1	B	42	TYR	2.5
1	A	34	TYR	2.5
1	D	93	GLU	2.5
1	D	6	LYS	2.5
1	D	38	VAL	2.5
1	B	96	PRO	2.5
1	C	142	GLU	2.4
1	B	115	VAL	2.4
1	B	93	GLU	2.4
1	B	2	ALA	2.4
1	B	155	GLY	2.4
1	A	82	MET	2.4
1	B	39	PHE	2.4
1	B	55	LEU	2.3
1	D	75	PRO	2.3
1	B	16	CYS	2.3
1	D	83	CYS	2.3
1	B	123	ASN	2.3
1	D	40	GLU	2.3
1	C	46	ILE	2.3
1	C	50	GLY	2.3
1	D	80	ILE	2.3
1	A	134	MET	2.3
1	D	82	MET	2.3
1	D	109	ASN	2.3
1	C	125	GLU	2.3
1	B	104	LYS	2.3
1	C	106	PHE	2.3
1	A	99	TRP	2.3
1	C	163	THR	2.3
1	A	151	ILE	2.2
1	C	177	ALA	2.2
1	B	91	SER	2.2
1	D	20	CYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	90	ASP	2.2
1	B	112	ILE	2.2
1	D	72	LEU	2.2
1	B	144	GLY	2.2
1	A	104	LYS	2.2
1	D	4	ILE	2.2
1	A	55	LEU	2.2
1	B	159	CYS	2.2
1	D	132	ALA	2.2
1	D	74	TYR	2.2
1	A	147	MET	2.2
1	B	6	LYS	2.2
1	C	117	ASN	2.2
1	D	120	ASP	2.2
1	A	175	THR	2.1
1	C	33	VAL	2.1
1	C	66	TYR	2.1
1	B	4	ILE	2.1
1	A	29	GLN	2.1
1	B	52	GLN	2.1
1	B	56	ALA	2.1
1	B	130	GLU	2.1
1	A	65	ASP	2.1
1	C	53	VAL	2.1
1	D	173	MET	2.1
1	D	153	ALA	2.1
1	A	157	MET	2.1
1	C	52	GLN	2.1
1	B	85	SER	2.1
1	A	109	ASN	2.1
1	A	49	ASP	2.1
1	B	81	LEU	2.1
1	D	107	CYS	2.1
1	D	161	ALA	2.1
1	A	138	PRO	2.1
1	C	85	SER	2.1
1	D	91	SER	2.1
1	B	35	VAL	2.1
1	B	25	ASN	2.1
1	A	84	PHE	2.1
1	A	121	LEU	2.0
1	D	123	ASN	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	87	ASP	2.0
1	C	127	THR	2.0
1	C	156	TYR	2.0
1	C	165	ASP	2.0
1	D	100	THR	2.0
1	C	84	PHE	2.0
1	A	68	ARG	2.0
1	B	21	LEU	2.0
1	B	131	LEU	2.0
1	B	160	SER	2.0
1	C	10	ILE	2.0
1	C	82	MET	2.0
1	B	79	VAL	2.0
1	D	5	ARG	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($\text{\AA}^2$)	Q<0.9
4	U6L	D	201	12/12	0.77	0.20	32,35,36,37	0
5	EPE	D	205	15/15	0.77	0.20	43,50,51,51	0
2	GDP	C	201	28/28	0.86	0.15	20,26,30,30	0
2	GDP	B	301	28/28	0.89	0.14	22,28,34,36	0
3	MG	B	302	1/1	0.89	0.24	2,2,2,2	0
3	MG	A	202	1/1	0.91	0.08	15,15,15,15	0
2	GDP	D	202	28/28	0.91	0.12	22,26,33,34	0
2	GDP	A	201	28/28	0.92	0.11	18,23,29,31	0
3	MG	B	303	1/1	0.92	0.18	20,20,20,20	0

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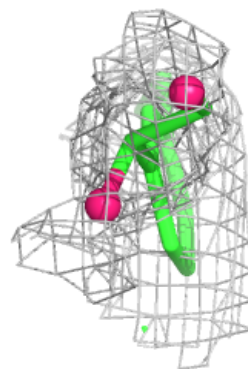
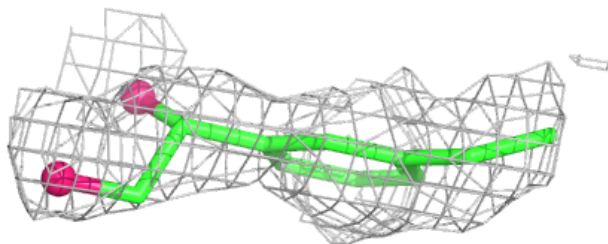
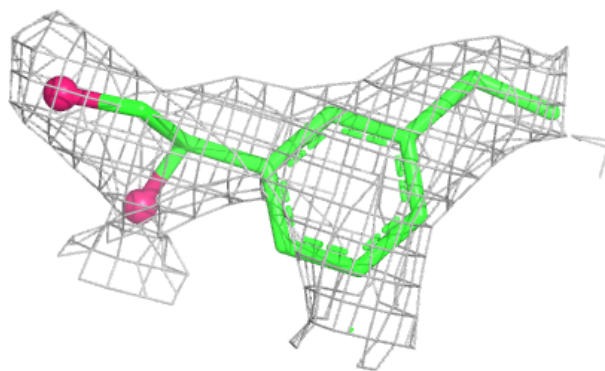
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MG	D	204	1/1	0.93	0.13	18,18,18,18	0
3	MG	D	203	1/1	0.96	0.13	10,10,10,10	0
3	MG	C	202	1/1	0.99	0.19	3,3,3,3	0

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.

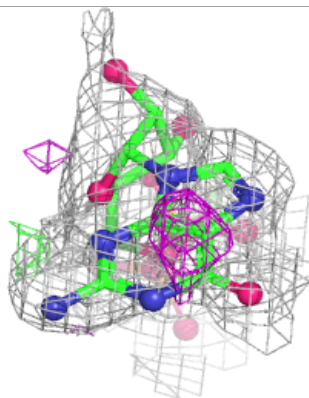
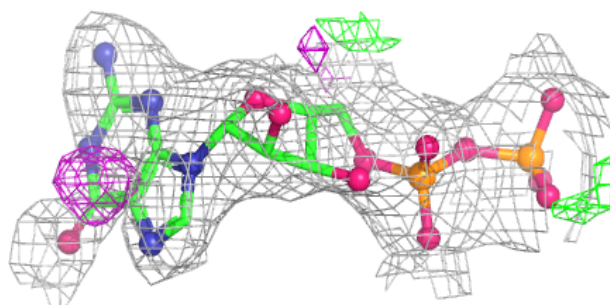
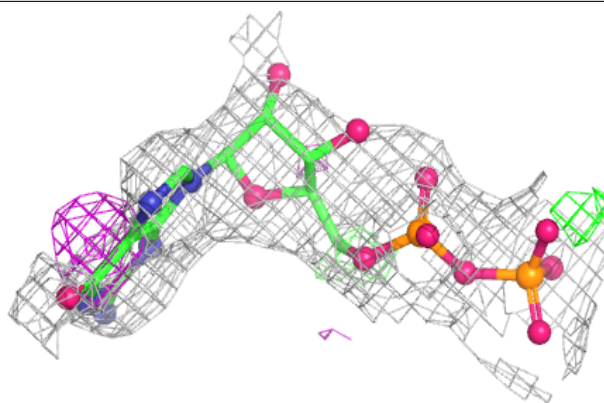
Electron density around U6L D 201:

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

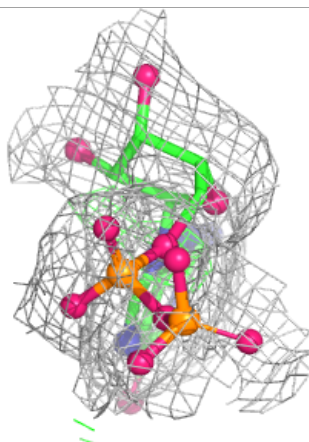
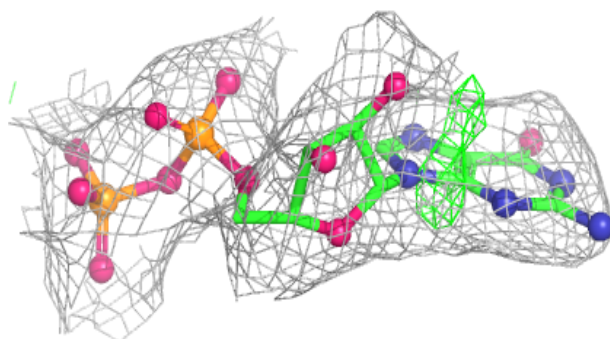
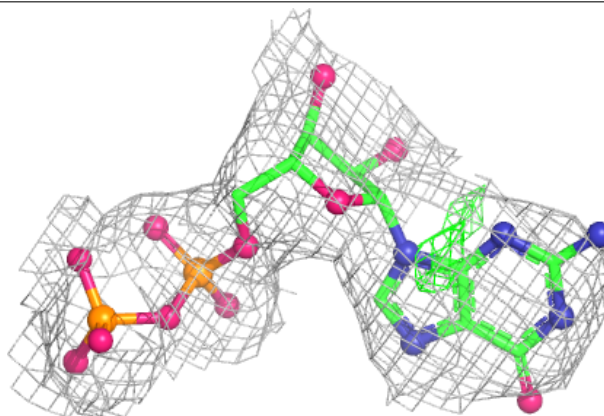


Electron density around GDP C 201:

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

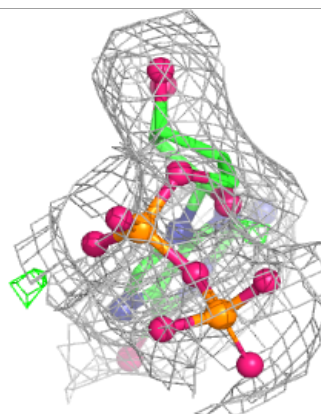
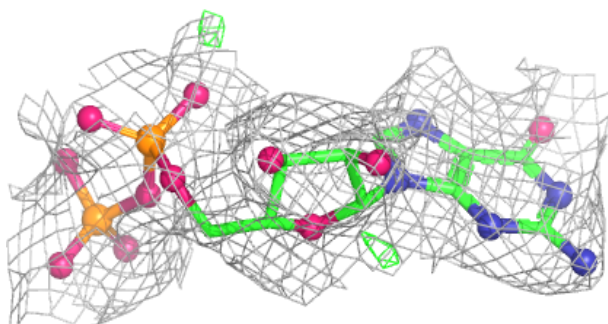
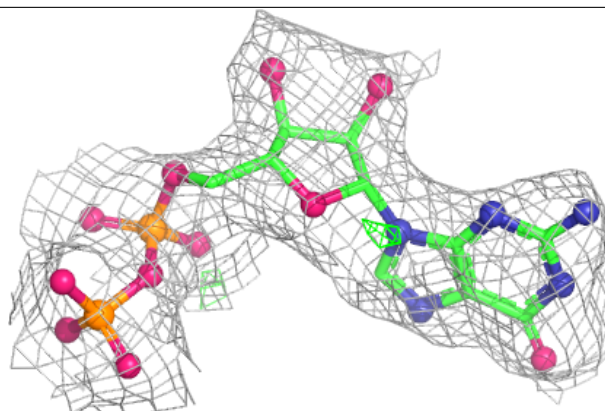
**Electron density around GDP B 301:**

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

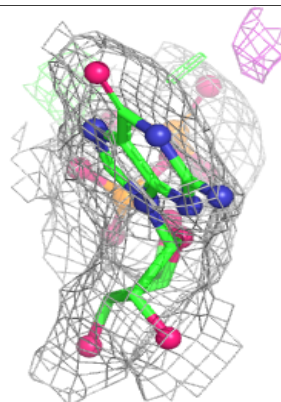
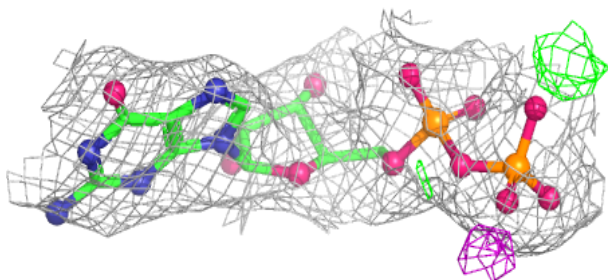
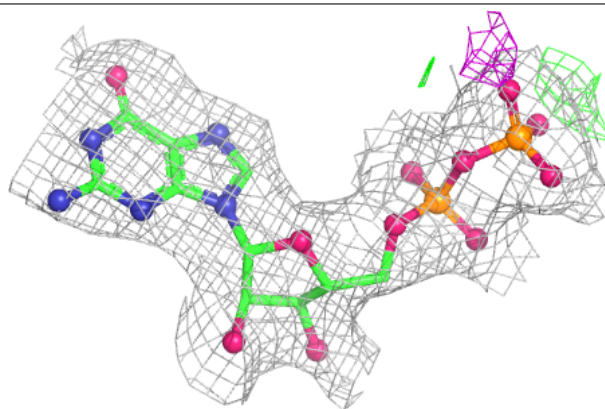


Electron density around GDP D 202:

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

**Electron density around GDP A 201:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.