



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2026 – 01:42 AM UTC

PDB ID : 5U7R / pdb_00005u7r
Title : Identification of A New Class of Potent Cdc7 Inhibitors Designed by Putative Pharmacophore Model: Synthesis and Biological Evaluation of 2,3-Dihydrothieno[3,2-d]pyrimidin-4(1H)-ones
Authors : Hoffman, I.D.; Skene, R.J.
Deposited on : 2016-12-12
Resolution : 3.33 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

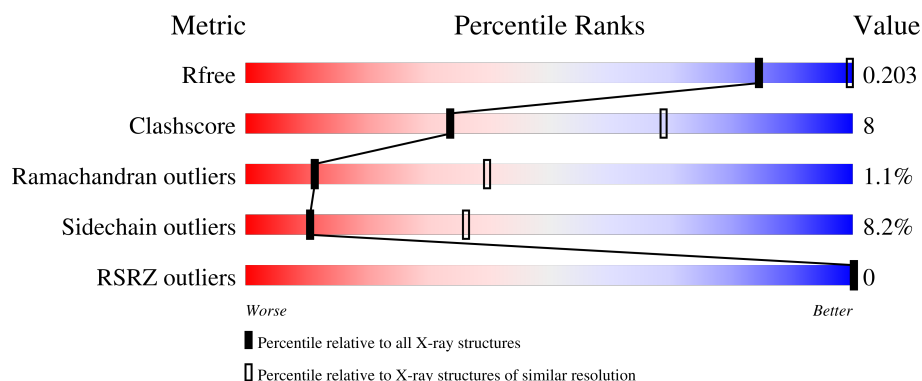
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.33 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	1434 (3.38-3.30)
Clashscore	190562	1479 (3.38-3.30)
Ramachandran outliers	187476	1456 (3.38-3.30)
Sidechain outliers	187428	1455 (3.38-3.30)
RSRZ outliers	180081	1434 (3.38-3.30)

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	395	
1	B	395	
1	C	395	
1	D	395	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12680 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.

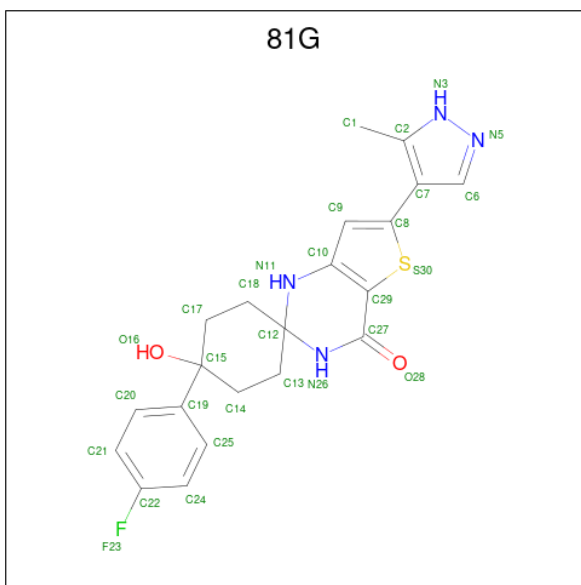
- Molecule 1 is a protein called Rho-associated protein kinase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	387	Total	C	N	O	S	0	0	0
			3140	2013	527	580	20			
1	B	387	Total	C	N	O	S	0	0	0
			3140	2013	527	580	20			
1	C	387	Total	C	N	O	S	0	0	0
			3140	2013	527	580	20			
1	D	382	Total	C	N	O	S	0	0	0
			3104	1991	522	571	20			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	270	TYR	PHE	conflict	UNP O75116
B	270	TYR	PHE	conflict	UNP O75116
C	270	TYR	PHE	conflict	UNP O75116
D	270	TYR	PHE	conflict	UNP O75116

- Molecule 2 is (1s,4s)-4-(4-fluorophenyl)-4-hydroxy-6'-(5-methyl-1H-pyrazol-4-yl)-1'H-spiro[cyclohexane-1,2'-thieno[3,2-d]pyrimidin]-4'(3'H)-one (CCD ID: 81G) (formula: C₂₁H₂₁FN₄O₂S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	S	0	0
			29	21	1	4	2	1		
2	B	1	Total	C	F	N	O	S	0	0
			29	21	1	4	2	1		
2	C	1	Total	C	F	N	O	S	0	0
			29	21	1	4	2	1		
2	D	1	Total	C	F	N	O	S	0	0
			29	21	1	4	2	1		

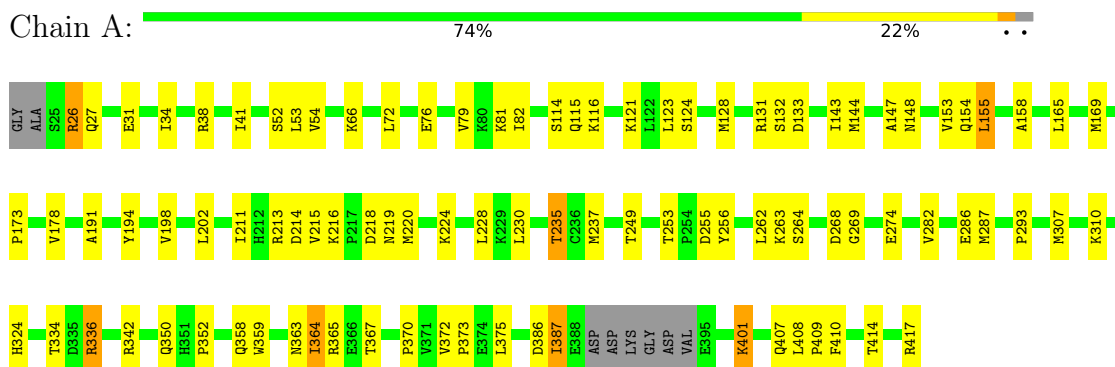
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	23	Total	O	0	0
			23	23		
3	B	4	Total	O	0	0
			4	4		
3	C	4	Total	O	0	0
			4	4		
3	D	9	Total	O	0	0
			9	9		

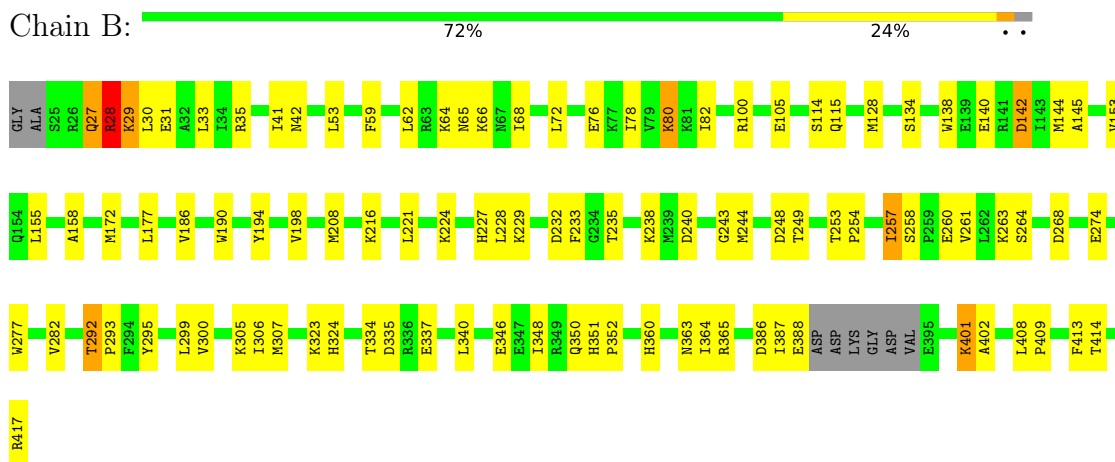
3 Residue-property plots

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.

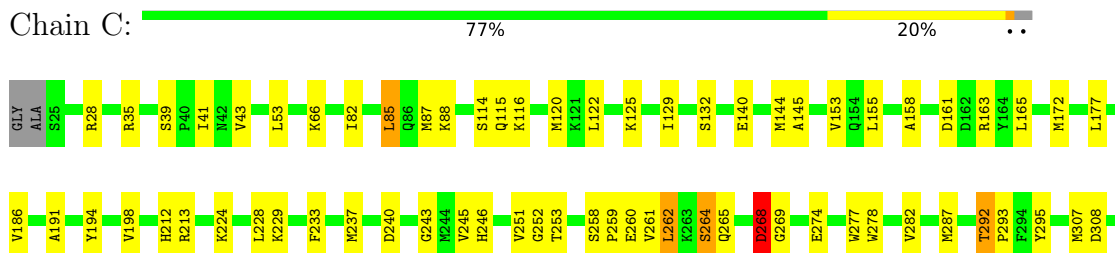
• Molecule 1: Rho-associated protein kinase 2

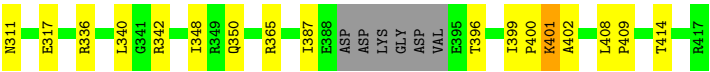


• Molecule 1: Rho-associated protein kinase 2

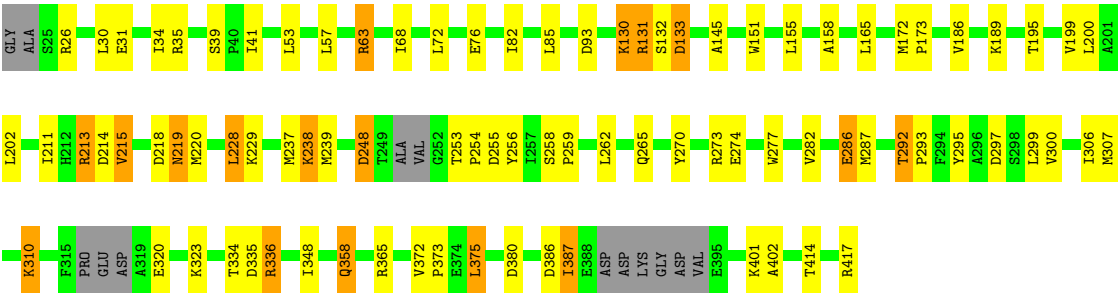


• Molecule 1: Rho-associated protein kinase 2





● Molecule 1: Rho-associated protein kinase 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	90.20Å 146.14Å 110.90Å 90.00° 96.03° 90.00°	Depositor
Resolution (Å)	49.40 – 3.33 49.40 – 3.33	Depositor EDS
% Data completeness (in resolution range)	99.5 (49.40-3.33) 99.6 (49.40-3.33)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 3.33Å)	Xtriage
Refinement program	REFMAC 5.7.0025	Depositor
R, R_{free}	0.200 , 0.241 0.199 , 0.203	Depositor DCC
R_{free} test set	2060 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	105.4	Xtriage
Anisotropy	0.056	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 121.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12680	wwPDB-VP
Average B, all atoms (Å ²)	129.0	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 33.63 % 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 7.7567e-04. 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 81G

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.39	0/3218	0.81	0/4349
1	B	0.40	0/3218	0.82	3/4349 (0.1%)
1	C	0.40	0/3218	0.84	2/4349 (0.0%)
1	D	0.39	0/3179	0.82	2/4291 (0.0%)
All	All	0.40	0/12833	0.82	7/17338 (0.0%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	364	ILE	N-CA-C	5.82	116.00	110.42
1	D	253	THR	CA-C-N	5.78	125.31	119.19
1	D	253	THR	C-N-CA	5.78	125.31	119.19
1	C	39	SER	CA-C-N	5.25	124.96	119.82
1	C	39	SER	C-N-CA	5.25	124.96	119.82
1	B	351	HIS	CA-C-N	5.00	124.72	119.82
1	B	351	HIS	C-N-CA	5.00	124.72	119.82

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3140	0	3065	60	0
1	B	3140	0	3065	51	0
1	C	3140	0	3065	42	0
1	D	3104	0	3032	50	0
2	A	29	0	0	1	0
2	B	29	0	0	0	0
2	C	29	0	0	0	0
2	D	29	0	0	0	0
3	A	23	0	0	1	0
3	B	4	0	0	0	0
3	C	4	0	0	0	0
3	D	9	0	0	0	0
All	All	12680	0	12227	192	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 8.

All (192) 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:26:ARG:HH11	1:A:26:ARG:HG3	1.29	0.95
1:A:41:ILE:HD13	1:D:82:ILE:HD11	1.50	0.92
1:D:63:ARG:HG2	1:D:63:ARG:HH11	1.36	0.88
1:B:41:ILE:HD13	1:C:82:ILE:HD11	1.65	0.78
1:D:72:LEU:O	1:D:76:GLU:HB2	1.83	0.78
1:D:292:THR:HG23	1:D:295:TYR:HB2	1.72	0.71
1:C:307:MET:HA	1:C:307:MET:HE2	1.73	0.71
1:A:256:TYR:OH	1:A:286:GLU:OE1	2.07	0.70
1:A:31:GLU:HG2	1:B:35:ARG:HD2	1.74	0.69
1:B:401:LYS:HG2	1:B:402:ALA:H	1.57	0.69
1:A:82:ILE:HD11	1:D:41:ILE:HD13	1.75	0.68
1:B:82:ILE:HD11	1:C:41:ILE:HD13	1.75	0.68
1:A:213:ARG:NH1	1:A:237:MET:HB2	2.10	0.67
1:B:53:LEU:HD13	1:C:53:LEU:HD13	1.77	0.67
1:B:240:ASP:HB2	1:B:244:MET:H	1.61	0.66
1:D:256:TYR:OH	1:D:286:GLU:OE1	2.13	0.65
1:A:26:ARG:HG3	1:A:26:ARG:NH1	2.03	0.64
1:A:216:LYS:NZ	1:A:253:THR:HG21	2.12	0.64
1:C:144:MET:HB2	1:C:153:VAL:O	1.98	0.64
1:A:143:ILE:O	1:A:147:ALA:HB2	1.97	0.63
1:A:401:LYS:HD3	1:A:401:LYS:N	2.14	0.63
1:B:144:MET:HG3	1:B:155:LEU:HB2	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:259:PRO:HA	1:C:307:MET:HE1	1.81	0.62
1:D:213:ARG:HH22	1:D:237:MET:CG	2.12	0.62
1:D:173:PRO:HB2	1:D:375:LEU:HD11	1.82	0.61
1:C:292:THR:HG23	1:C:295:TYR:HB2	1.82	0.61
1:C:336:ARG:O	1:C:342:ARG:HD2	2.01	0.61
1:D:131:ARG:HG2	1:D:133:ASP:HB2	1.82	0.60
1:B:76:GLU:O	1:B:80:LYS:HG2	2.02	0.60
1:A:263:LYS:HG2	1:A:307:MET:HE1	1.83	0.60
1:D:239:MET:HE1	1:D:273:ARG:HH21	1.68	0.59
1:B:27:GLN:HG2	1:B:30:LEU:HB3	1.85	0.59
1:D:63:ARG:HH11	1:D:63:ARG:CG	2.09	0.59
1:B:216:LYS:NZ	1:B:253:THR:HG21	2.16	0.59
1:A:158:ALA:O	1:A:414:THR:HA	2.03	0.59
1:A:155:LEU:HD21	1:A:158:ALA:HB2	1.84	0.59
1:D:145:ALA:HB2	1:D:155:LEU:HD23	1.84	0.58
1:D:211:ILE:HG22	1:D:273:ARG:HA	1.86	0.58
1:A:363:ASN:O	1:A:365:ARG:N	2.36	0.58
1:B:114:SER:O	1:B:115:GLN:HB2	2.04	0.58
1:B:138:TRP:O	1:B:142:ASP:HB2	2.03	0.57
1:A:128:MET:HE3	1:A:410:PHE:CE1	2.40	0.57
1:B:292:THR:HG23	1:B:295:TYR:HB2	1.87	0.57
1:B:307:MET:HA	1:B:307:MET:HE2	1.88	0.56
1:C:114:SER:O	1:C:115:GLN:HB2	2.04	0.56
1:A:191:ALA:HA	1:A:287:MET:HE3	1.86	0.56
1:B:340:LEU:HG	1:B:348:ILE:HD13	1.87	0.55
1:A:214:ASP:HA	3:A:609:HOH:O	2.07	0.55
1:C:144:MET:HG3	1:C:155:LEU:HB2	1.87	0.55
1:A:307:MET:HA	1:A:307:MET:HE2	1.89	0.55
1:D:130:LYS:O	1:D:131:ARG:HB2	2.06	0.55
1:A:72:LEU:O	1:A:76:GLU:HB2	2.07	0.55
1:D:259:PRO:HA	1:D:307:MET:HE1	1.90	0.54
1:A:216:LYS:HZ2	1:A:253:THR:HG21	1.73	0.53
1:A:53:LEU:HD13	1:D:53:LEU:HD13	1.89	0.53
1:A:214:ASP:O	1:A:219:ASN:ND2	2.42	0.53
1:C:262:LEU:CD1	1:C:307:MET:HE3	2.39	0.53
1:B:28:ARG:O	1:B:31:GLU:HG2	2.08	0.53
1:C:212:HIS:O	1:C:213:ARG:HB2	2.09	0.52
1:D:131:ARG:C	1:D:133:ASP:H	2.18	0.52
1:B:254:PRO:HG2	1:B:299:LEU:HD22	1.91	0.52
1:C:145:ALA:HB2	1:C:155:LEU:HD23	1.92	0.52
1:C:213:ARG:HH12	1:C:251:VAL:HB	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:401:LYS:H	1:B:401:LYS:HD2	1.75	0.51
1:C:35:ARG:HH12	1:D:35:ARG:HG3	1.75	0.51
1:C:246:HIS:CE1	1:C:268:ASP:HB3	2.46	0.51
1:D:63:ARG:HE	1:D:72:LEU:HD13	1.75	0.51
1:A:211:ILE:HG13	1:A:213:ARG:HD2	1.92	0.51
1:D:151:TRP:CH2	1:D:365:ARG:HG3	2.45	0.51
1:C:262:LEU:HD11	1:C:307:MET:HE3	1.91	0.51
1:D:238:LYS:HD2	1:D:238:LYS:H	1.75	0.51
1:A:214:ASP:HB2	1:A:235:THR:OG1	2.10	0.51
1:B:194:TYR:O	1:B:198:VAL:HG23	2.11	0.50
1:B:72:LEU:O	1:B:76:GLU:HB2	2.12	0.50
1:A:144:MET:HE1	1:A:169:MET:SD	2.52	0.50
1:B:140:GLU:HA	1:B:233:PHE:HB2	1.94	0.50
1:B:144:MET:HB2	1:B:153:VAL:O	2.11	0.50
1:B:417:ARG:HH21	1:D:417:ARG:HG2	1.76	0.50
1:B:401:LYS:HG2	1:B:402:ALA:N	2.23	0.49
1:A:194:TYR:CZ	1:A:370:PRO:HG2	2.47	0.49
1:B:186:VAL:HG13	1:B:190:TRP:HB2	1.93	0.49
1:C:120:MET:HE3	1:C:122:LEU:HD11	1.94	0.49
1:D:220:MET:HB3	1:D:228:LEU:HD21	1.95	0.49
1:D:172:MET:SD	1:D:229:LYS:HD2	2.52	0.49
1:A:114:SER:O	1:A:115:GLN:HB2	2.13	0.49
1:A:202:LEU:HD21	1:A:215:VAL:HG21	1.93	0.49
1:B:360:HIS:HB2	1:B:363:ASN:HB3	1.94	0.49
1:C:240:ASP:HB2	1:C:243:GLY:H	1.78	0.49
1:C:282:VAL:HG13	1:C:293:PRO:HD2	1.94	0.49
1:A:264:SER:HB2	1:A:269:GLY:HA3	1.94	0.48
1:B:27:GLN:O	1:B:29:LYS:N	2.46	0.48
1:B:228:LEU:O	1:B:365:ARG:NH1	2.45	0.48
1:C:194:TYR:O	1:C:198:VAL:HG23	2.13	0.48
1:A:124:SER:O	1:A:128:MET:HB2	2.14	0.48
1:D:211:ILE:CG2	1:D:273:ARG:HA	2.43	0.48
1:A:173:PRO:HG3	1:A:224:LYS:HG3	1.95	0.48
1:B:28:ARG:CZ	1:B:28:ARG:HB2	2.44	0.48
1:A:224:LYS:HB3	1:C:224:LYS:HE3	1.95	0.48
1:B:346:GLU:O	1:B:350:GLN:HG2	2.14	0.48
1:D:63:ARG:HG2	1:D:63:ARG:NH1	2.16	0.48
1:A:128:MET:HE3	1:A:410:PHE:HE1	1.76	0.48
1:A:153:VAL:HG23	1:A:230:LEU:O	2.14	0.47
1:D:213:ARG:HH22	1:D:237:MET:HG2	1.78	0.47
1:C:125:LYS:NZ	1:C:161:ASP:O	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:202:LEU:HD21	1:D:215:VAL:HG21	1.96	0.47
1:A:143:ILE:O	1:A:147:ALA:CB	2.62	0.47
1:C:158:ALA:O	1:C:414:THR:HA	2.15	0.47
1:D:375:LEU:HD13	1:D:380:ASP:HB3	1.95	0.47
1:A:131:ARG:C	1:A:133:ASP:H	2.22	0.47
1:C:129:ILE:HD11	1:C:409:PRO:HG2	1.97	0.47
1:D:274:GLU:HA	1:D:277:TRP:HD1	1.80	0.47
1:A:121:LYS:HE2	2:A:500:81G:O28	2.14	0.47
1:B:145:ALA:HB2	1:B:155:LEU:HD23	1.97	0.47
1:D:228:LEU:HD22	1:D:229:LYS:N	2.30	0.47
1:A:213:ARG:NH1	1:A:235:THR:O	2.49	0.46
1:C:191:ALA:HA	1:C:287:MET:HE3	1.97	0.46
1:C:264:SER:HB2	1:C:269:GLY:HA3	1.97	0.46
1:B:324:HIS:ND1	1:B:352:PRO:HB2	2.31	0.46
1:C:186:VAL:HG11	1:C:287:MET:HG2	1.97	0.46
1:A:154:GLN:HG3	1:A:417:ARG:HH22	1.80	0.46
1:B:227:HIS:HB3	1:B:365:ARG:NH1	2.30	0.46
1:B:258:SER:OG	1:B:261:VAL:HG23	2.15	0.46
1:D:358:GLN:H	1:D:358:GLN:CD	2.23	0.46
1:A:363:ASN:O	1:A:364:ILE:C	2.59	0.46
1:A:274:GLU:HG3	1:A:336:ARG:HB3	1.98	0.46
1:A:401:LYS:N	1:A:401:LYS:CD	2.79	0.46
1:B:33:LEU:HB3	1:B:41:ILE:HD11	1.97	0.45
1:B:282:VAL:HG13	1:B:293:PRO:HD2	1.99	0.45
1:B:128:MET:HB3	1:B:134:SER:HB3	1.98	0.45
1:C:260:GLU:HG2	1:C:261:VAL:N	2.32	0.45
1:A:173:PRO:HG3	1:A:224:LYS:HE3	1.99	0.45
1:A:359:TRP:HB3	1:A:367:THR:HG21	1.99	0.45
1:D:195:THR:O	1:D:199:VAL:HG23	2.17	0.45
1:A:26:ARG:HH11	1:A:27:GLN:H	1.65	0.45
1:A:324:HIS:ND1	1:A:352:PRO:HB2	2.31	0.44
1:B:158:ALA:HB1	1:B:413:PHE:CE2	2.52	0.44
1:C:259:PRO:HA	1:C:307:MET:CE	2.45	0.44
1:A:386:ASP:C	1:A:387:ILE:HG13	2.38	0.44
1:C:140:GLU:HA	1:C:233:PHE:HB2	1.98	0.44
1:B:172:MET:HE2	1:B:229:LYS:HB2	2.00	0.44
1:C:399:ILE:HA	1:C:400:PRO:HD3	1.77	0.44
1:B:274:GLU:HA	1:B:277:TRP:CD1	2.53	0.44
1:A:282:VAL:HG13	1:A:293:PRO:CD	2.48	0.44
1:D:306:ILE:HG22	1:D:307:MET:HE2	2.00	0.44
1:C:213:ARG:NH1	1:C:251:VAL:HB	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:401:LYS:HD3	1:A:401:LYS:H	1.79	0.44
1:A:52:SER:HB3	1:A:407:GLN:HB2	1.99	0.43
1:A:54:VAL:HG21	1:A:79:VAL:HG21	2.00	0.43
1:D:310:LYS:HD2	1:D:310:LYS:H	1.82	0.43
1:A:144:MET:HE3	1:A:153:VAL:HB	1.99	0.43
1:C:308:ASP:OD2	1:C:311:ASN:HB2	2.18	0.43
1:D:213:ARG:HH22	1:D:237:MET:HG3	1.80	0.43
1:D:214:ASP:O	1:D:219:ASN:ND2	2.51	0.43
1:D:386:ASP:O	1:D:387:ILE:HG23	2.18	0.43
1:B:221:LEU:O	1:B:228:LEU:HA	2.18	0.43
1:B:261:VAL:O	1:B:264:SER:HB3	2.18	0.43
1:B:408:LEU:N	1:B:409:PRO:CD	2.82	0.43
1:D:336:ARG:H	1:D:336:ARG:HG3	1.59	0.43
1:B:30:LEU:HB2	1:C:85:LEU:HD12	2.00	0.43
1:C:116:LYS:H	1:C:116:LYS:HG2	1.66	0.43
1:D:186:VAL:HG11	1:D:287:MET:HG2	2.01	0.43
1:D:401:LYS:HG3	1:D:402:ALA:H	1.83	0.43
1:D:228:LEU:HD22	1:D:229:LYS:H	1.85	0.42
1:C:172:MET:HE2	1:C:229:LYS:HB2	2.01	0.42
1:A:372:VAL:HA	1:A:373:PRO:HD3	1.90	0.42
1:D:57:LEU:HB3	1:D:68:ILE:HG23	2.02	0.42
1:A:194:TYR:O	1:A:198:VAL:HG23	2.20	0.42
1:B:100:ARG:HA	1:B:105:GLU:HA	2.01	0.42
1:C:340:LEU:HG	1:C:348:ILE:HD13	2.01	0.42
1:C:401:LYS:HG3	1:C:402:ALA:H	1.85	0.42
1:B:274:GLU:H	1:B:274:GLU:CD	2.28	0.42
1:A:408:LEU:N	1:A:409:PRO:CD	2.83	0.41
1:D:158:ALA:O	1:D:414:THR:HA	2.20	0.41
1:D:372:VAL:HA	1:D:373:PRO:HD3	1.93	0.41
1:B:158:ALA:O	1:B:414:THR:HA	2.20	0.41
1:D:254:PRO:O	1:D:255:ASP:C	2.63	0.41
1:A:123:LEU:HB2	1:A:165:LEU:HB3	2.02	0.41
1:B:65:ASN:HB3	1:B:68:ILE:HB	2.03	0.41
1:D:131:ARG:C	1:D:133:ASP:N	2.79	0.41
1:D:282:VAL:HG13	1:D:293:PRO:HD2	2.03	0.41
1:A:401:LYS:CD	1:A:401:LYS:H	2.33	0.41
1:C:408:LEU:N	1:C:409:PRO:CD	2.84	0.41
1:B:59:PHE:HB2	1:B:62:LEU:HD12	2.02	0.41
1:C:259:PRO:HD3	1:C:278:TRP:CE2	2.56	0.41
1:C:274:GLU:HA	1:C:277:TRP:CD1	2.56	0.41
1:D:401:LYS:HG3	1:D:402:ALA:N	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:257:ILE:HG22	1:B:306:ILE:HD13	2.02	0.40
1:A:202:LEU:HD11	1:A:215:VAL:HG21	2.03	0.40
1:A:220:MET:HB3	1:A:228:LEU:HD11	2.04	0.40
1:A:262:LEU:HD23	1:A:307:MET:HE3	2.03	0.40
1:D:248:ASP:OD1	1:D:248:ASP:N	2.54	0.40
1:B:253:THR:HA	1:B:254:PRO:HD3	1.98	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	383/395 (97%)	365 (95%)	16 (4%)	2 (0%)	24	55
1	B	383/395 (97%)	356 (93%)	21 (6%)	6 (2%)	7	32
1	C	383/395 (97%)	352 (92%)	26 (7%)	5 (1%)	9	35
1	D	374/395 (95%)	349 (93%)	21 (6%)	4 (1%)	11	40
All	All	1523/1580 (96%)	1422 (93%)	84 (6%)	17 (1%)	11	40

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	364	ILE
1	B	28	ARG
1	B	224	LYS
1	A	132	SER
1	B	235	THR
1	B	268	ASP
1	C	87	MET
1	C	132	SER
1	C	268	ASP

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Mol	Chain	Res	Type
1	C	28	ARG
1	D	26	ARG
1	B	232	ASP
1	D	132	SER
1	D	320	GLU
1	D	131	ARG
1	B	243	GLY
1	C	252	GLY

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	341/346 (99%)	318 (93%)	23 (7%)	15	43
1	B	341/346 (99%)	313 (92%)	28 (8%)	10	36
1	C	341/346 (99%)	318 (93%)	23 (7%)	15	43
1	D	337/346 (97%)	300 (89%)	37 (11%)	6	23
All	All	1360/1384 (98%)	1249 (92%)	111 (8%)	10	36

All (111) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	ARG
1	A	34	ILE
1	A	38	ARG
1	A	66	LYS
1	A	81	LYS
1	A	116	LYS
1	A	148	ASN
1	A	155	LEU
1	A	178	VAL
1	A	218	ASP
1	A	235	THR
1	A	249	THR
1	A	255	ASP

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Mol	Chain	Res	Type
1	A	268	ASP
1	A	310	LYS
1	A	334	THR
1	A	336	ARG
1	A	342	ARG
1	A	350	GLN
1	A	358	GLN
1	A	375	LEU
1	A	387	ILE
1	A	401	LYS
1	B	27	GLN
1	B	28	ARG
1	B	29	LYS
1	B	42	ASN
1	B	64	LYS
1	B	66	LYS
1	B	78	ILE
1	B	80	LYS
1	B	142	ASP
1	B	177	LEU
1	B	208	MET
1	B	238	LYS
1	B	248	ASP
1	B	249	THR
1	B	257	ILE
1	B	260	GLU
1	B	263	LYS
1	B	292	THR
1	B	300	VAL
1	B	305	LYS
1	B	323	LYS
1	B	334	THR
1	B	335	ASP
1	B	337	GLU
1	B	386	ASP
1	B	387	ILE
1	B	388	GLU
1	B	401	LYS
1	C	43	VAL
1	C	66	LYS
1	C	85	LEU
1	C	88	LYS

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Mol	Chain	Res	Type
1	C	163	ARG
1	C	165	LEU
1	C	177	LEU
1	C	228	LEU
1	C	237	MET
1	C	245	VAL
1	C	253	THR
1	C	258	SER
1	C	262	LEU
1	C	264	SER
1	C	265	GLN
1	C	268	ASP
1	C	292	THR
1	C	317	GLU
1	C	350	GLN
1	C	365	ARG
1	C	387	ILE
1	C	396	THR
1	C	401	LYS
1	D	30	LEU
1	D	31	GLU
1	D	34	ILE
1	D	39	SER
1	D	63	ARG
1	D	85	LEU
1	D	93	ASP
1	D	130	LYS
1	D	133	ASP
1	D	165	LEU
1	D	189	LYS
1	D	200	LEU
1	D	213	ARG
1	D	215	VAL
1	D	218	ASP
1	D	219	ASN
1	D	228	LEU
1	D	238	LYS
1	D	248	ASP
1	D	258	SER
1	D	262	LEU
1	D	265	GLN
1	D	270	TYR

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Mol	Chain	Res	Type
1	D	286	GLU
1	D	292	THR
1	D	297	ASP
1	D	299	LEU
1	D	300	VAL
1	D	310	LYS
1	D	323	LYS
1	D	334	THR
1	D	335	ASP
1	D	336	ARG
1	D	348	ILE
1	D	358	GLN
1	D	375	LEU
1	D	387	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	ASN
1	A	51	ASN
1	A	107	GLN
1	A	183	ASN
1	A	343	ASN
1	A	350	GLN
1	A	360	HIS
1	A	363	ASN
1	B	65	ASN
1	B	107	GLN
1	B	115	GLN
1	B	183	ASN
1	B	225	HIS
1	B	309	HIS
1	B	327	ASN
1	C	27	GLN
1	C	86	GLN
1	C	107	GLN
1	C	219	ASN
1	C	246	HIS
1	C	327	ASN
1	C	351	HIS
1	C	358	GLN
1	C	406	ASN

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Mol	Chain	Res	Type
1	D	27	GLN
1	D	51	ASN
1	D	154	GLN
1	D	183	ASN
1	D	219	ASN
1	D	246	HIS
1	D	265	GLN
1	D	356	ASN
1	D	358	GLN
1	D	360	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	81G	B	500	-	32,33,33	0.84	1 (3%)	42,51,51	1.67	9 (21%)
2	81G	C	500	-	32,33,33	0.82	1 (3%)	42,51,51	1.59	10 (23%)
2	81G	A	500	-	32,33,33	0.82	1 (3%)	42,51,51	1.72	11 (26%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	81G	D	500	-	32,33,33	0.81	1 (3%)	42,51,51	1.65	10 (23%)

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	81G	B	500	-	-	4/10/38/38	0/5/5/5
2	81G	C	500	-	-	2/10/38/38	0/5/5/5
2	81G	A	500	-	-	4/10/38/38	0/5/5/5
2	81G	D	500	-	-	4/10/38/38	0/5/5/5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	500	81G	C9-C10	-2.41	1.37	1.42
2	B	500	81G	C9-C10	-2.35	1.37	1.42
2	D	500	81G	C9-C10	-2.32	1.37	1.42
2	A	500	81G	C9-C10	-2.05	1.38	1.42

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	81G	C18-C12-N26	-5.22	100.90	109.61
2	A	500	81G	C18-C12-N26	-5.11	101.09	109.61
2	D	500	81G	C18-C12-N26	-4.75	101.70	109.61
2	C	500	81G	C18-C12-N26	-4.62	101.91	109.61
2	B	500	81G	C17-C15-C19	-3.86	103.37	111.05
2	A	500	81G	C17-C15-C19	-3.48	104.14	111.05
2	D	500	81G	C17-C15-C19	-3.43	104.24	111.05
2	B	500	81G	C6-C7-C2	3.05	107.48	104.90
2	C	500	81G	C17-C15-C19	-2.83	105.43	111.05
2	A	500	81G	C10-C9-C8	2.77	115.96	111.72
2	C	500	81G	C6-C7-C2	2.76	107.24	104.90
2	B	500	81G	C6-N5-N3	2.64	109.15	105.19
2	A	500	81G	C6-N5-N3	2.61	109.10	105.19
2	B	500	81G	C2-N3-N5	-2.59	108.25	112.26
2	C	500	81G	C6-N5-N3	2.59	109.08	105.19
2	D	500	81G	C10-C9-C8	2.57	115.66	111.72
2	A	500	81G	C2-N3-N5	-2.53	108.34	112.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	500	81G	C10-C9-C8	2.49	115.53	111.72
2	C	500	81G	C2-N3-N5	-2.48	108.43	112.26
2	A	500	81G	C29-C27-N26	2.46	118.77	113.96
2	D	500	81G	C14-C15-C19	2.44	115.89	111.05
2	B	500	81G	C7-C6-N5	-2.35	107.83	111.12
2	A	500	81G	C13-C14-C15	-2.32	110.43	112.25
2	D	500	81G	C29-C27-N26	2.30	118.46	113.96
2	D	500	81G	C2-N3-N5	-2.29	108.72	112.26
2	D	500	81G	C6-N5-N3	2.27	108.59	105.19
2	D	500	81G	C6-C7-C2	2.27	106.82	104.90
2	A	500	81G	C24-C22-C21	-2.23	119.88	122.80
2	A	500	81G	C6-C7-C2	2.23	106.78	104.90
2	C	500	81G	C7-C6-N5	-2.21	108.03	111.12
2	A	500	81G	C27-C29-S30	2.18	128.64	124.97
2	C	500	81G	C29-C27-N26	2.13	118.12	113.96
2	B	500	81G	C10-C9-C8	2.10	114.94	111.72
2	B	500	81G	C29-C27-N26	2.10	118.05	113.96
2	D	500	81G	C25-C19-C15	-2.09	117.90	121.11
2	B	500	81G	C25-C19-C15	-2.06	117.93	121.11
2	C	500	81G	C24-C22-C21	-2.05	120.11	122.80
2	A	500	81G	C7-C6-N5	-2.02	108.30	111.12
2	C	500	81G	C13-C12-N26	2.01	112.96	109.61
2	D	500	81G	C24-C22-C21	-2.00	120.17	122.80

There are no chirality outliers.

All (14) torsion outliers are listed below:

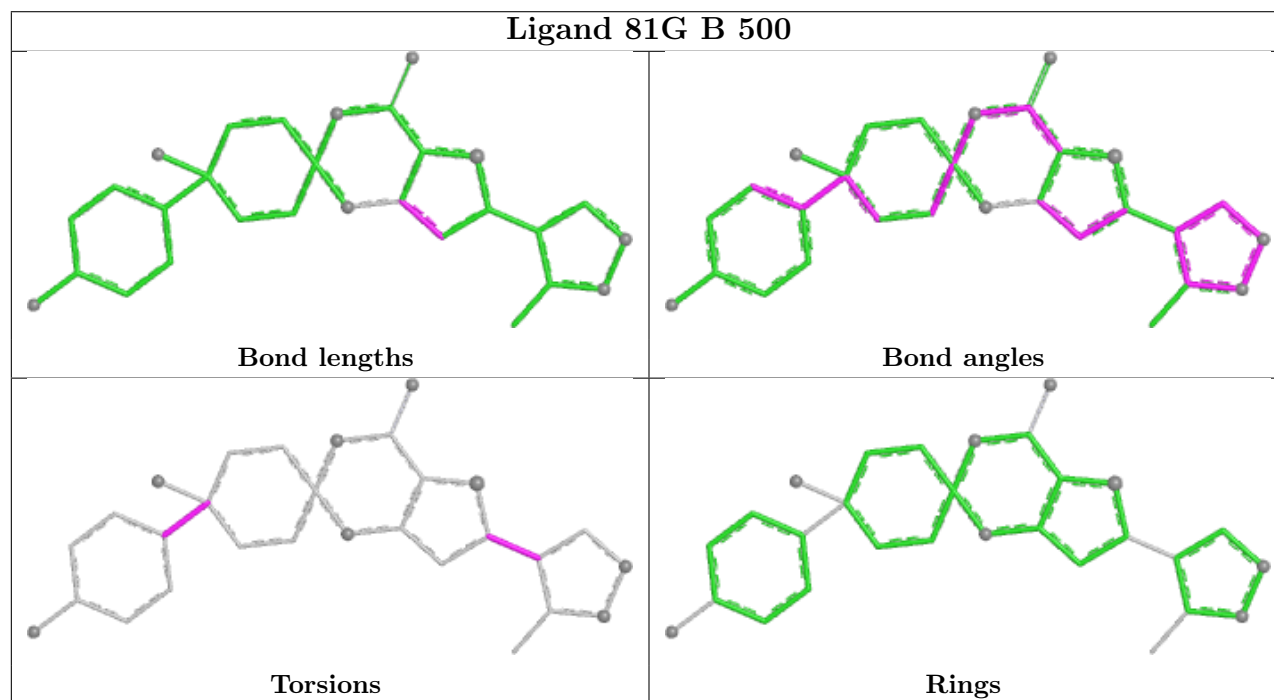
Mol	Chain	Res	Type	Atoms
2	A	500	81G	C6-C7-C8-C9
2	A	500	81G	C6-C7-C8-S30
2	B	500	81G	C6-C7-C8-C9
2	D	500	81G	C6-C7-C8-C9
2	D	500	81G	C6-C7-C8-S30
2	A	500	81G	C2-C7-C8-S30
2	D	500	81G	C2-C7-C8-C9
2	D	500	81G	C2-C7-C8-S30
2	B	500	81G	C6-C7-C8-S30
2	C	500	81G	C6-C7-C8-C9
2	C	500	81G	C6-C7-C8-S30
2	A	500	81G	C2-C7-C8-C9
2	B	500	81G	C2-C7-C8-C9
2	B	500	81G	C14-C15-C19-C20

There are no ring outliers.

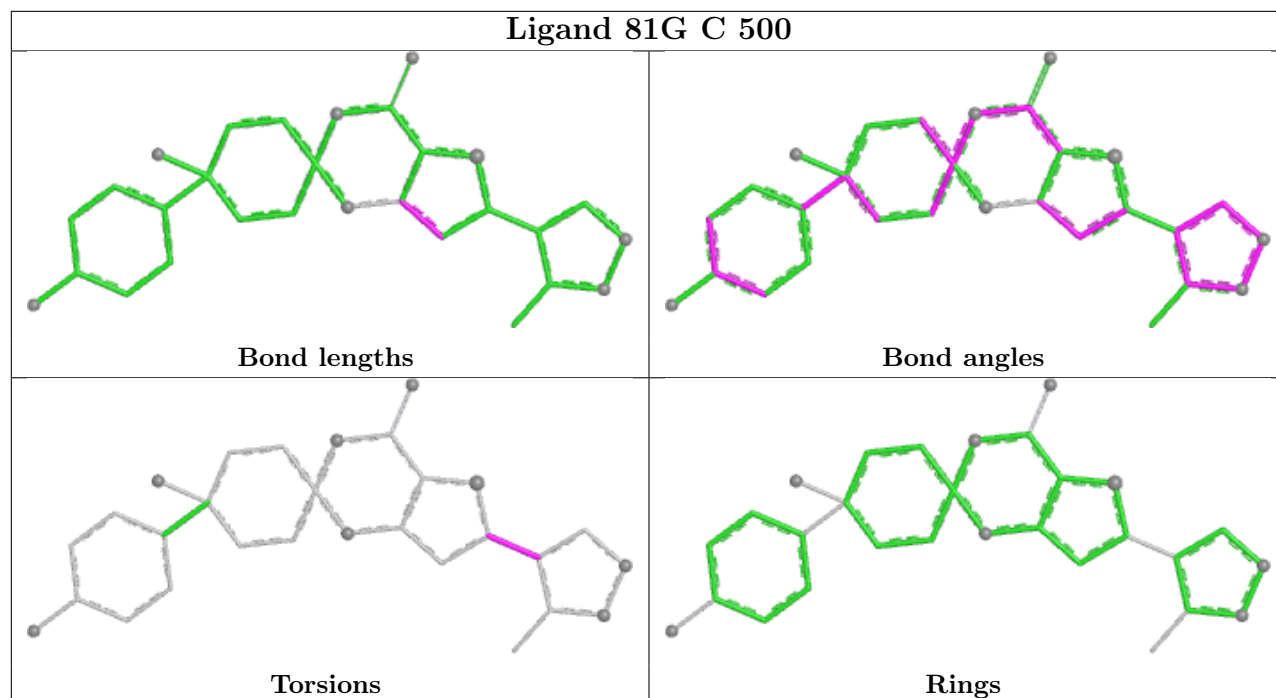
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	81G	1	0

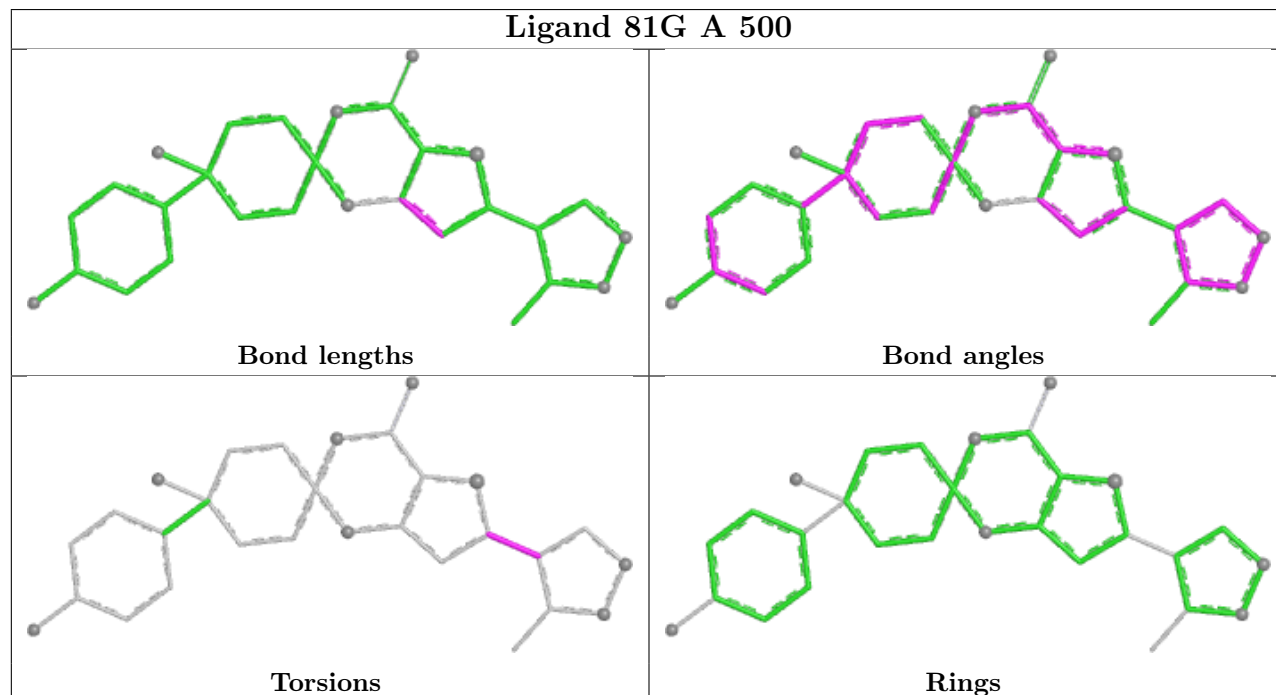
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

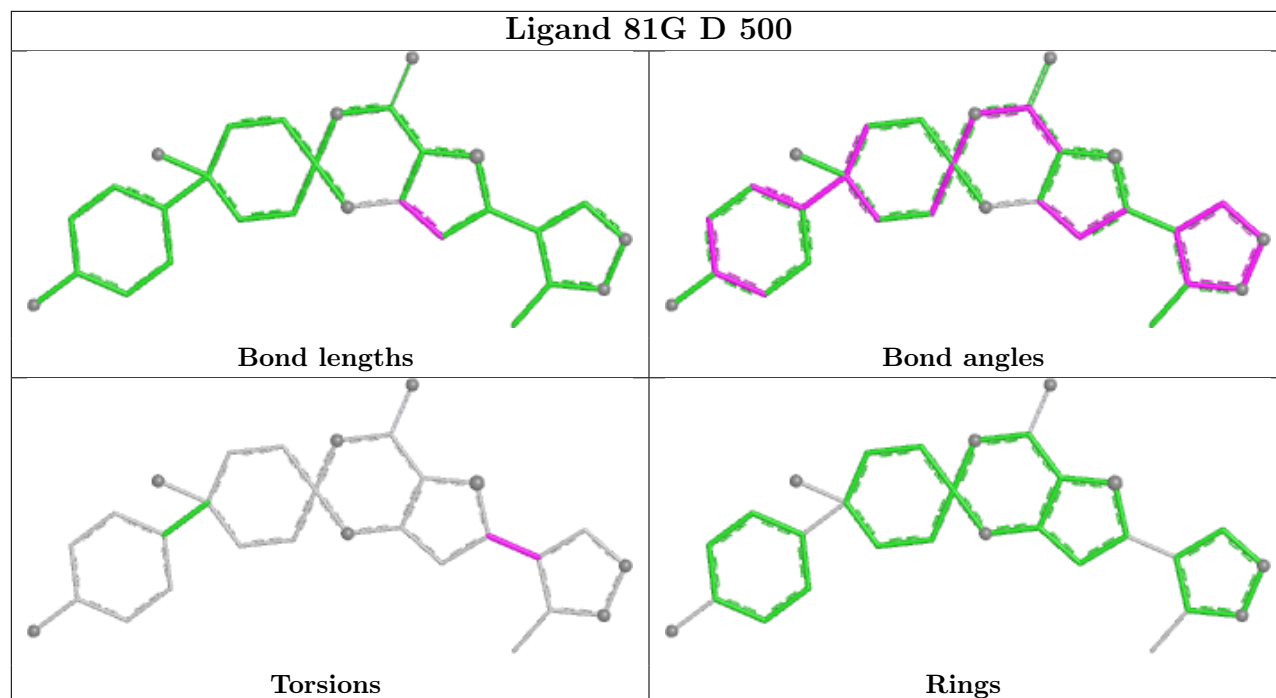


Ligand 81G C 500



Ligand 81G A 500





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 [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	387/395 (97%)	-0.75	0 100 100	85, 118, 157, 205	0
1	B	387/395 (97%)	-0.60	0 100 100	93, 132, 181, 236	0
1	C	387/395 (97%)	-0.55	0 100 100	98, 131, 180, 229	0
1	D	382/395 (96%)	-0.64	0 100 100	86, 122, 167, 207	0
All	All	1543/1580 (97%)	-0.64	0 100 100	85, 126, 174, 236	0

There are no RSRZ outliers to report.

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](#)

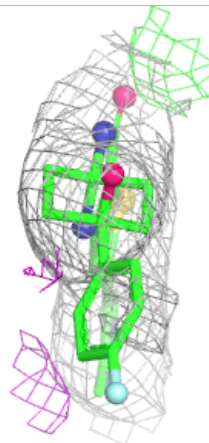
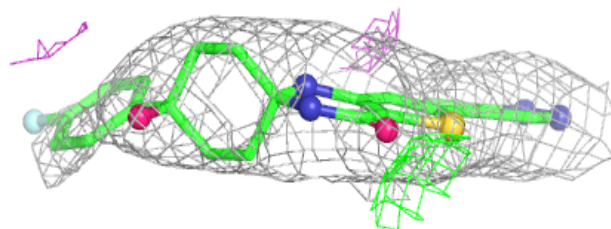
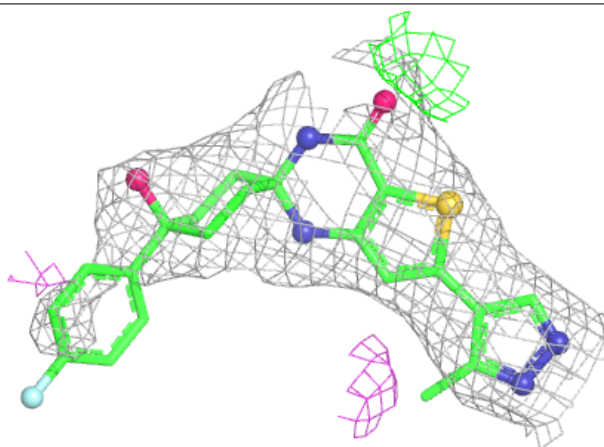
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	81G	C	500	29/29	0.94	0.11	105,128,176,186	0
2	81G	B	500	29/29	0.95	0.10	107,126,179,183	0
2	81G	D	500	29/29	0.95	0.09	100,118,151,156	0
2	81G	A	500	29/29	0.96	0.09	101,108,141,148	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.

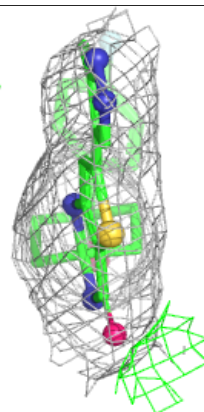
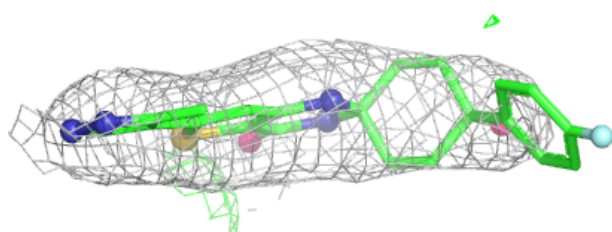
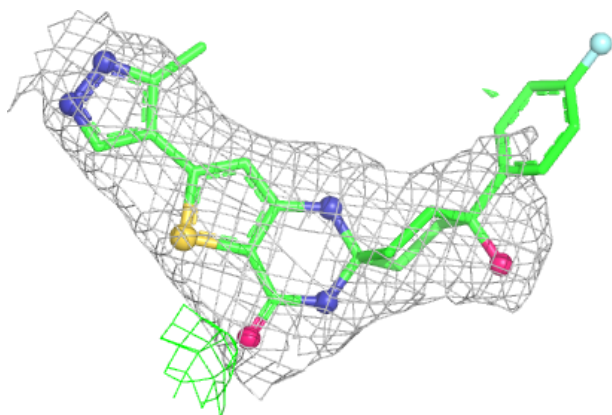
Electron density around 81G C 500:

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

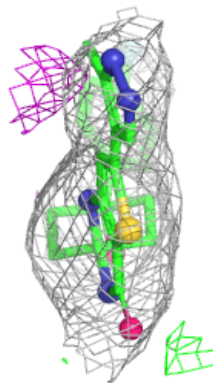
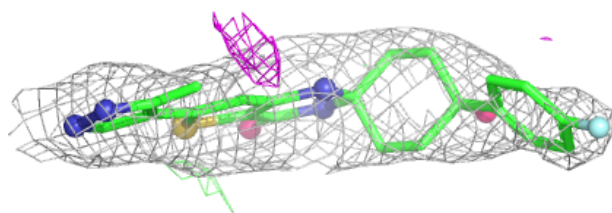
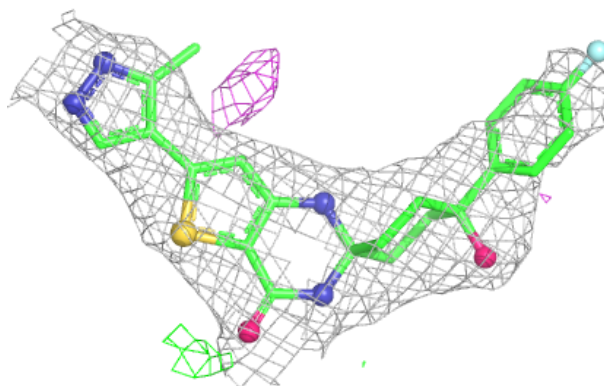


Electron density around 81G B 500:

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

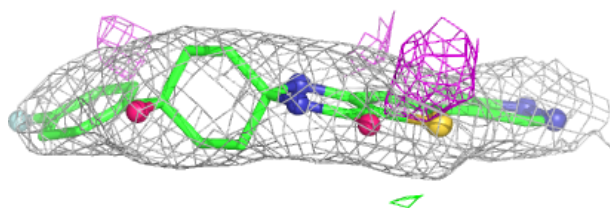
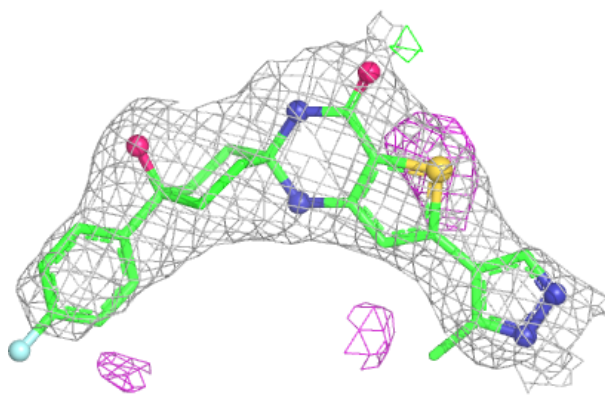
**Electron density around 81G D 500:**

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



Electron density around 81G A 500:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.