Full wwPDB X-ray Structure Validation Report (i)

Jun 26, 2024 – 05:15 AM EDT

PDB ID	:	7NXJ
Title	:	Crystal structure of human Cdk13/Cyclin K in complex with the inhibitor THZ531
Authors	:	Anand, K.; Greifenberg, A.K.; Kaltheuner, I.H.; Geyer, M.
Deposited on	:	2021-03-18
Resolution	:	2.36 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 (i) 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 (i)) were used in the production of this report:

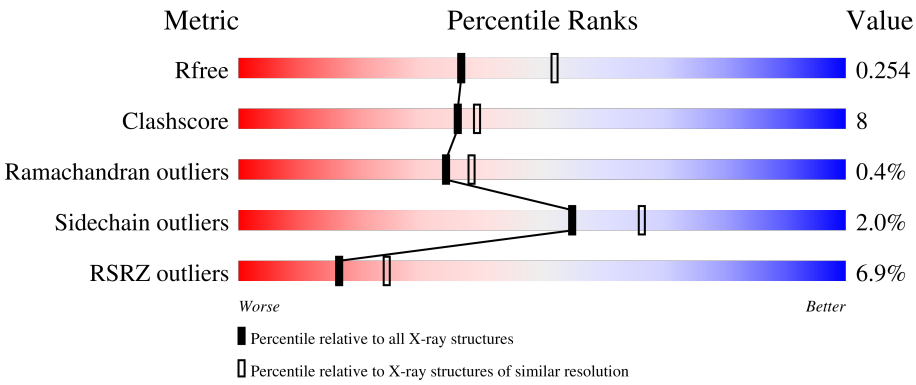
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.36 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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 <=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	347	11% 73% 20% • 6%
1	C	347	6% 75% 17% • 7%
2	B	268	4% 78% 13% 9%
2	D	268	3% 82% 10% 8%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 9402 atoms, of which 35 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 Cyclin-dependent kinase 13.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
1	A	325	Total C N O P S 2585 1665 429 471 1 19	0	0	0
1	C	323	Total C H N O P S 2623 1665 35 435 468 1 19	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	693	GLY	-	expression tag	UNP Q14004
C	693	GLY	-	expression tag	UNP Q14004

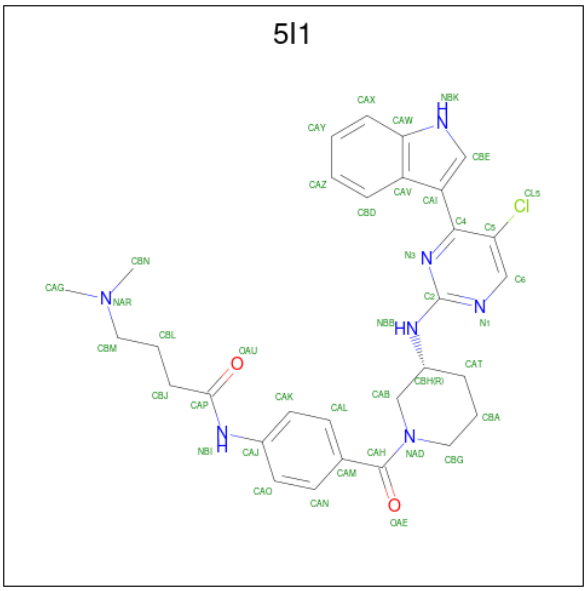
- Molecule 2 is a protein called Cyclin-K.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
2	B	244	Total C N O S 2003 1303 333 356 11	0	0	0
2	D	247	Total C N O S 2028 1320 338 357 13	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	GLY	-	expression tag	UNP O75909
D	0	GLY	-	expression tag	UNP O75909

- Molecule 3 is N-[4-[(3R)-3-[[5-chloranyl-4-(1H-indol-3-yl)pyrimidin-2-yl]amino]piperidin-1-yl]carbonylphenyl]-4-(dimethylamino)butanamide (three-letter code: 5I1) (formula: $C_{30}H_{34}ClN_7O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Cl	N	O		
3	A	1	40	30	1	7	2	0	0
			Total	C	Cl	N	O		
3	C	1	40	30	1	7	2	0	0

- Molecule 4 is water.

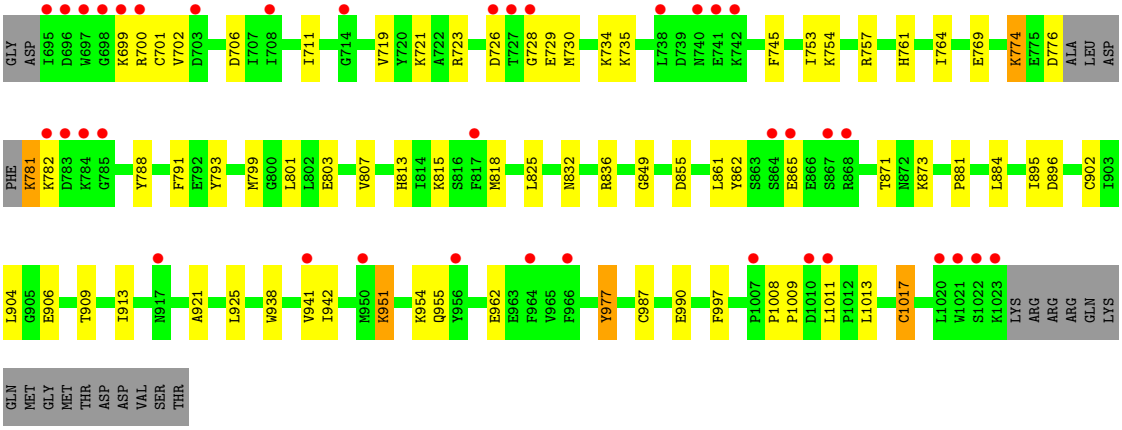
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	16	16	16	0	0
			Total	O		
4	B	17	17	17	0	0
			Total	O		
4	C	25	25	25	0	0
			Total	O		
4	D	25	25	25	0	0

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.

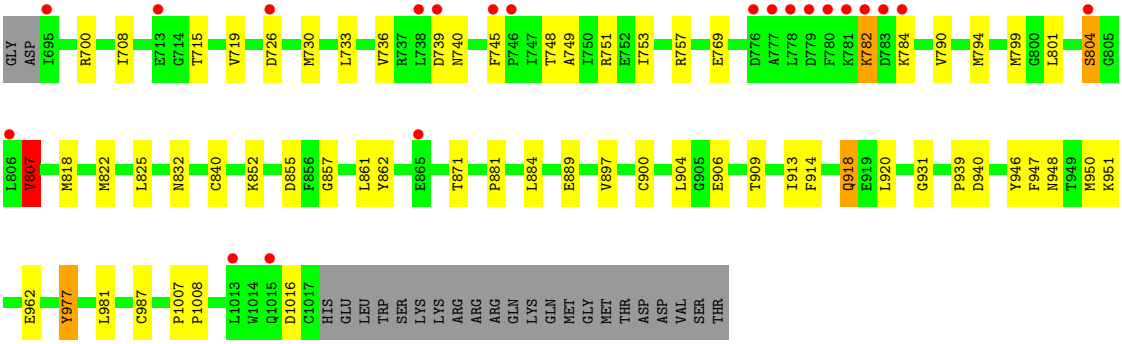
• Molecule 1: Cyclin-dependent kinase 13

Chain A:



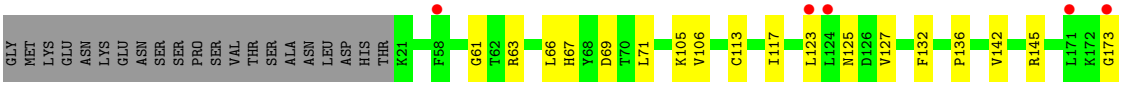
• Molecule 1: Cyclin-dependent kinase 13

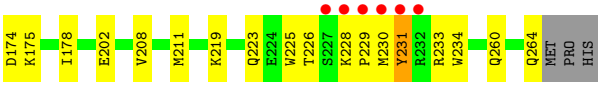
Chain C:



• Molecule 2: Cyclin-K

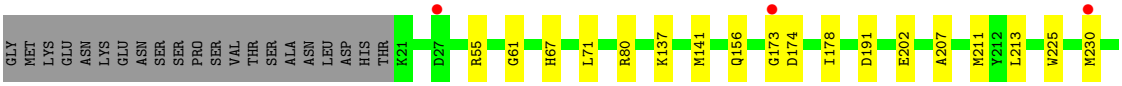
Chain B:

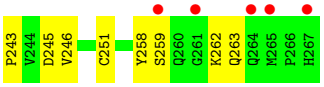




• Molecule 2: Cyclin-K

Chain D:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	49.88Å 149.45Å 92.01Å 90.00° 94.61° 90.00°	Depositor
Resolution (Å)	45.86 – 2.36 45.86 – 2.36	Depositor EDS
% Data completeness (in resolution range)	98.2 (45.86-2.36) 98.2 (45.86-2.36)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.64 (at 2.37Å)	Xtriage
Refinement program	PHENIX (1.18.2_3874: ???)	Depositor
R, R_{free}	0.212 , 0.252 0.218 , 0.254	Depositor DCC
R_{free} test set	1353 reflections (2.50%)	wwPDB-VP
Wilson B-factor (Å^2)	57.5	Xtriage
Anisotropy	0.422	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 43.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	9402	wwPDB-VP
Average B, all atoms (Å^2)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.40% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: TPO, 5I1

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	0.30	0/2635	0.49	0/3565
1	C	0.32	0/2637	0.50	0/3565
2	B	0.29	0/2058	0.46	0/2791
2	D	0.25	0/2085	0.40	0/2827
All	All	0.29	0/9415	0.47	0/12748

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2585	0	2525	54	0
1	C	2588	35	2560	51	2
2	B	2003	0	1954	25	2
2	D	2028	0	1982	17	0
3	A	40	0	0	8	0
3	C	40	0	0	2	0
4	A	16	0	0	0	0
4	B	17	0	0	0	0
4	C	25	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	25	0	0	1	0
All	All	9367	35	9021	146	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (146) 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:1017:CYS:SG	3:A:2000:5I1:CBL	2.01	1.49
2:B:173:GLY:HA3	2:B:225:TRP:CD1	1.53	1.42
1:C:745:PHE:CE1	1:C:749:ALA:HB3	1.84	1.11
1:C:736:VAL:HG11	1:C:745:PHE:CE2	1.89	1.07
1:A:1017:CYS:SG	3:A:2000:5I1:CBM	2.52	0.98
2:B:173:GLY:CA	2:B:225:TRP:CD1	2.46	0.98
1:C:745:PHE:CE1	1:C:749:ALA:CB	2.47	0.98
1:C:801:LEU:O	1:C:807:VAL:HG21	1.63	0.97
1:C:745:PHE:HE1	1:C:749:ALA:HB3	1.30	0.93
1:C:745:PHE:HE1	1:C:749:ALA:CB	1.83	0.90
2:B:174:ASP:O	2:B:178:ILE:HG12	1.74	0.88
1:C:708:ILE:HD12	1:C:730:MET:HE1	1.57	0.87
1:A:1017:CYS:SG	3:A:2000:5I1:CAP	2.63	0.87
2:D:173:GLY:HA3	2:D:225:TRP:CD1	2.10	0.85
2:B:63:ARG:HH11	2:B:123:LEU:HG	1.44	0.82
1:C:736:VAL:HG11	1:C:745:PHE:CZ	2.17	0.78
1:A:1017:CYS:SG	3:A:2000:5I1:CBJ	2.73	0.76
2:B:173:GLY:HA3	2:B:225:TRP:HD1	1.44	0.76
1:C:947:PHE:O	1:C:951:LYS:HG3	1.85	0.75
2:B:63:ARG:NH1	2:B:123:LEU:HG	2.05	0.70
1:C:700:ARG:NH1	1:C:769:GLU:OE1	2.24	0.70
1:A:700:ARG:NH2	1:A:769:GLU:OE2	2.25	0.69
1:A:801:LEU:HD21	1:A:1011:LEU:HD13	1.74	0.69
1:C:832:ASN:HA	1:C:862:TYR:CE1	2.30	0.67
1:C:948:ASN:O	1:C:951:LYS:NZ	2.28	0.67
1:A:745:PHE:O	2:B:105:LYS:NZ	2.24	0.65
1:A:754:LYS:NZ	2:B:106:VAL:O	2.30	0.65
1:C:736:VAL:CG1	1:C:745:PHE:CE2	2.74	0.65
1:C:708:ILE:CD1	1:C:730:MET:HE1	2.27	0.64
1:A:723:ARG:HH12	1:A:728:GLY:C	2.01	0.64
1:A:723:ARG:HD3	1:A:730:MET:SD	2.40	0.62
1:A:723:ARG:HH12	1:A:729:GLU:N	1.97	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:745:PHE:CE1	1:C:749:ALA:HB1	2.32	0.62
1:C:782:LYS:HD2	1:C:784:LYS:HB2	1.82	0.62
2:D:55:ARG:NH2	2:D:263:GLN:OE1	2.34	0.61
1:A:711:ILE:HD11	1:A:719:VAL:HG12	1.81	0.61
1:A:734:LYS:HE2	1:A:791:PHE:CE1	2.38	0.58
1:C:920:LEU:HD23	1:C:950:MET:HB2	1.85	0.58
1:A:776:ASP:C	1:A:781:LYS:O	2.42	0.58
1:A:942:ILE:HD12	2:D:67:HIS:CD2	2.39	0.57
1:A:776:ASP:CB	1:A:781:LYS:O	2.52	0.57
1:A:1017:CYS:SG	3:A:2000:5I1:OAU	2.61	0.57
1:A:818:MET:HG3	1:A:904:LEU:HD13	1.86	0.57
3:A:2000:5I1:CBE	3:A:2000:5I1:CL5	2.89	0.57
2:B:125:ASN:OD1	2:B:127:VAL:HG12	2.06	0.56
2:D:174:ASP:O	2:D:178:ILE:HG13	2.05	0.56
2:D:80:ARG:NH1	4:D:301:HOH:O	2.28	0.56
1:A:881:PRO:HD2	1:A:884:LEU:HD12	1.86	0.56
1:C:822:MET:HE3	1:C:900:CYS:HB3	1.89	0.55
1:A:721:LYS:HE2	1:A:793:TYR:CZ	2.42	0.55
2:B:142:VAL:O	2:B:145:ARG:HG2	2.07	0.55
2:B:223:GLN:O	2:B:229:PRO:HA	2.08	0.54
1:C:799:MET:HB2	1:C:840:CYS:HB3	1.90	0.54
1:A:921:ALA:O	1:A:925:LEU:HD13	2.07	0.54
1:C:804:SER:O	1:C:804:SER:OG	2.12	0.53
2:D:61:GLY:HA3	2:D:71:LEU:HD21	1.90	0.53
1:A:990:GLU:OE2	1:A:990:GLU:N	2.34	0.52
1:C:719:VAL:HG11	3:C:2000:5I1:CAZ	2.40	0.52
1:A:954:LYS:HG2	1:A:955:GLN:N	2.24	0.52
2:D:173:GLY:HA3	2:D:225:TRP:HD1	1.68	0.52
2:D:191:ASP:OD2	2:D:258:TYR:OH	2.25	0.52
2:B:208:VAL:HA	2:B:211:MET:HE2	1.91	0.52
1:C:794:MET:SD	1:C:852:LYS:HD2	2.49	0.52
2:B:208:VAL:HA	2:B:211:MET:CE	2.40	0.51
3:A:2000:5I1:CBG	3:A:2000:5I1:CAN	2.90	0.50
2:D:213:LEU:HB2	2:D:251:CYS:SG	2.51	0.50
1:A:849:GLY:HA3	1:A:1008:PRO:HB3	1.94	0.50
1:A:832:ASN:HB3	1:A:862:TYR:O	2.12	0.49
1:A:1013:LEU:HD23	1:A:1013:LEU:O	2.12	0.49
1:C:700:ARG:NH2	1:C:769:GLU:OE1	2.43	0.49
1:C:906:GLU:O	1:C:909:THR:O	2.31	0.49
1:C:962:GLU:OE1	1:C:962:GLU:N	2.45	0.49
2:B:67:HIS:CE1	2:B:69:ASP:HB2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:913:ILE:HD12	1:C:914:PHE:N	2.29	0.48
2:D:137:LYS:O	2:D:141:MET:HG2	2.15	0.47
1:C:977:TYR:CE1	1:C:987:CYS:HB3	2.49	0.47
1:C:832:ASN:O	1:C:861:LEU:HD12	2.15	0.47
2:D:61:GLY:HA3	2:D:71:LEU:CD2	2.45	0.47
1:A:702:VAL:HG21	1:A:788:TYR:CE1	2.50	0.46
1:A:753:ILE:O	1:A:757:ARG:HG3	2.15	0.46
2:B:173:GLY:HA3	2:B:225:TRP:NE1	2.18	0.46
1:C:881:PRO:HD2	1:C:884:LEU:HD12	1.96	0.46
2:B:61:GLY:HA3	2:B:71:LEU:HD21	1.97	0.46
1:A:951:LYS:HE2	1:A:951:LYS:HB2	1.49	0.46
1:A:902:CYS:HA	1:A:913:ILE:HD11	1.97	0.46
1:C:751:ARG:HD2	1:C:857:GLY:O	2.16	0.46
1:A:711:ILE:CG2	1:A:721:LYS:HB2	2.45	0.46
3:C:2000:5I1:OAU	3:C:2000:5I1:CAK	2.64	0.46
1:A:938:TRP:O	1:A:941:VAL:HG22	2.15	0.46
1:A:706:ASP:O	1:A:723:ARG:N	2.39	0.45
1:C:753:ILE:O	1:C:757:ARG:HG3	2.16	0.45
2:B:202:GLU:OE1	2:B:202:GLU:N	2.43	0.45
1:A:807:VAL:HG21	1:A:1011:LEU:CD2	2.46	0.45
1:A:938:TRP:CG	1:A:941:VAL:HG13	2.52	0.45
2:B:175:LYS:HE3	2:B:175:LYS:HB3	1.68	0.45
2:B:132:PHE:HB3	2:B:136:PRO:HA	1.98	0.45
1:A:873:LYS:HE2	1:A:873:LYS:HB3	1.83	0.45
1:A:1017:CYS:CB	3:A:2000:5I1:CBL	2.90	0.45
2:B:219:LYS:HE3	2:B:219:LYS:HB2	1.59	0.45
1:A:799:MET:O	1:A:803:GLU:HG2	2.17	0.44
1:C:913:ILE:HD13	1:C:914:PHE:CD2	2.53	0.44
1:C:931:GLY:O	1:C:981:LEU:HD11	2.17	0.44
1:A:774:LYS:HE3	1:A:774:LYS:HB2	1.52	0.44
1:A:938:TRP:CD2	1:A:941:VAL:HG13	2.52	0.44
1:C:939:PRO:HD2	4:C:2105:HOH:O	2.18	0.44
1:A:723:ARG:HH22	1:A:728:GLY:HA2	1.82	0.44
2:B:113:CYS:O	2:B:117:ILE:HD12	2.18	0.44
1:C:818:MET:HG3	1:C:904:LEU:HD13	1.99	0.44
1:C:769:GLU:HG3	1:C:790:VAL:HB	2.00	0.43
2:D:202:GLU:OE1	2:D:202:GLU:N	2.38	0.43
1:A:832:ASN:O	1:A:861:LEU:HG	2.18	0.43
1:C:719:VAL:HA	1:C:733:LEU:O	2.18	0.43
2:D:243:PRO:HG2	2:D:246:VAL:HG23	2.00	0.43
1:C:736:VAL:HG11	1:C:745:PHE:CD2	2.47	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:977:TYR:CD1	1:C:987:CYS:HB3	2.53	0.43
1:A:906:GLU:O	1:A:909:THR:O	2.35	0.43
2:B:231:TYR:HD1	2:B:231:TYR:HA	1.65	0.43
2:D:259:SER:OG	2:D:262:LYS:HE2	2.17	0.43
1:A:776:ASP:C	1:A:782:LYS:HB2	2.38	0.43
2:B:226:THR:HG21	2:B:234:TRP:HB2	2.00	0.43
1:A:701:CYS:SG	1:A:702:VAL:N	2.91	0.43
2:D:207:ALA:O	2:D:211:MET:HG3	2.19	0.43
1:C:715:THR:HG21	1:C:748:THR:HG21	2.00	0.42
1:A:942:ILE:HD12	2:D:67:HIS:NE2	2.34	0.42
2:D:230:MET:SD	2:D:230:MET:N	2.84	0.42
1:C:946:TYR:O	1:C:950:MET:HG2	2.20	0.42
1:C:918:GLN:OE1	1:C:920:LEU:HB3	2.20	0.42
1:A:813:HIS:CE1	1:A:1009:PRO:HB3	2.55	0.41
1:C:739:ASP:OD1	1:C:740:ASN:N	2.53	0.41
1:C:801:LEU:HD23	1:C:801:LEU:HA	1.94	0.41
1:A:699:LYS:HD2	1:A:699:LYS:HA	1.76	0.41
1:A:977:TYR:CE1	1:A:987:CYS:HB3	2.55	0.41
1:C:913:ILE:CD1	1:C:914:PHE:CD2	3.03	0.41
1:C:977:TYR:CZ	1:C:987:CYS:HB3	2.55	0.41
2:B:211:MET:HB2	2:B:211:MET:HE3	1.94	0.41
1:A:735:LYS:HD2	1:A:788:TYR:CE2	2.55	0.41
1:A:761:HIS:HB3	1:A:764:ILE:HG12	2.03	0.41
1:A:815:LYS:HE2	1:A:997:PHE:O	2.20	0.41
1:A:825:LEU:HD21	1:A:896:ASP:HB3	2.02	0.41
1:A:836:ARG:HG2	1:A:895:ILE:HD11	2.03	0.41
2:B:66:LEU:HD23	2:B:66:LEU:HA	1.92	0.41
1:C:825:LEU:HD22	1:C:900:CYS:SG	2.61	0.41
1:C:897:VAL:HG11	1:C:987:CYS:O	2.22	0.40
1:C:1016:ASP:OD1	1:C:1016:ASP:N	2.50	0.40
1:C:700:ARG:CZ	1:C:769:GLU:OE1	2.69	0.40
1:C:1007:PRO:HA	1:C:1008:PRO:HD3	1.90	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:230:MET:CB	1:C:889:GLU:OE1[1_655]	1.49	0.71
2:B:230:MET:CB	1:C:889:GLU:CD[1_655]	2.04	0.16

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	320/347 (92%)	306 (96%)	13 (4%)	1 (0%)	41	47
1	C	320/347 (92%)	308 (96%)	10 (3%)	2 (1%)	25	27
2	B	242/268 (90%)	238 (98%)	2 (1%)	2 (1%)	19	20
2	D	245/268 (91%)	243 (99%)	2 (1%)	0	100	100
All	All	1127/1230 (92%)	1095 (97%)	27 (2%)	5 (0%)	34	38

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	231	TYR
2	B	233	ARG
1	C	804	SER
1	A	865	GLU
1	C	807	VAL

5.3.2 Protein sidechains (i)

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	273/310 (88%)	265 (97%)	8 (3%)	42	52
1	C	276/310 (89%)	269 (98%)	7 (2%)	47	58
2	B	212/241 (88%)	209 (99%)	3 (1%)	67	78
2	D	215/241 (89%)	213 (99%)	2 (1%)	78	87
All	All	976/1102 (89%)	956 (98%)	20 (2%)	55	66

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	726	ASP
1	A	774	LYS
1	A	781	LYS
1	A	855	ASP
1	A	951	LYS
1	A	962	GLU
1	A	977	TYR
1	A	1017	CYS
2	B	228	LYS
2	B	260	GLN
2	B	264	GLN
1	C	726	ASP
1	C	782	LYS
1	C	807	VAL
1	C	855	ASP
1	C	918	GLN
1	C	940	ASP
1	C	977	TYR
2	D	156	GLN
2	D	245	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	67	HIS
2	B	260	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	$\#\|Z\| > 2$
1	TPO	A	871	1	8,10,11	1.60	1 (12%)	10,14,16	1.74	1 (10%)
1	TPO	C	871	1	8,10,11	1.10	0	10,14,16	1.70	1 (10%)

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
1	TPO	A	871	1	-	1/9/11/13	-
1	TPO	C	871	1	-	1/9/11/13	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	871	TPO	P-O1P	3.42	1.61	1.50

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	871	TPO	P-OG1-CB	-4.72	108.95	123.21
1	C	871	TPO	P-OG1-CB	-4.59	109.34	123.21

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	871	TPO	O-C-CA-CB
1	C	871	TPO	O-C-CA-CB

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

2 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	$\#\|Z\| > 2$
3	5I1	C	2000	-	41,44,44	2.27	9 (21%)	52,61,61	1.86	11 (21%)
3	5I1	A	2000	-	41,44,44	0.47	0	52,61,61	0.95	2 (3%)

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	5I1	C	2000	-	-	3/22/36/36	0/5/5/5
3	5I1	A	2000	-	-	3/22/36/36	0/5/5/5

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	2000	5I1	C2-NBB	8.81	1.46	1.34
3	C	2000	5I1	CAH-NAD	5.37	1.46	1.34
3	C	2000	5I1	CAP-NBI	5.10	1.46	1.35
3	C	2000	5I1	CAB-NAD	3.73	1.51	1.46
3	C	2000	5I1	CAJ-NBI	2.77	1.47	1.41
3	C	2000	5I1	CAM-CAH	2.71	1.54	1.50
3	C	2000	5I1	CBG-NAD	2.44	1.51	1.47
3	C	2000	5I1	C5-CL5	2.19	1.78	1.73
3	C	2000	5I1	OAU-CAP	-2.18	1.18	1.23

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	2000	5I1	N1-C2-N3	-6.07	120.80	126.55
3	C	2000	5I1	C6-C5-C4	-5.54	117.67	120.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	2000	5I1	C6-C5-C4	-4.46	118.20	120.39
3	C	2000	5I1	CBJ-CAP-NBI	3.78	121.25	114.59
3	C	2000	5I1	CAJ-NBI-CAP	-3.74	120.96	127.50
3	C	2000	5I1	C2-NBB-CBH	-3.46	118.47	124.31
3	C	2000	5I1	C6-N1-C2	3.10	120.55	115.88
3	C	2000	5I1	C5-C6-N1	-3.01	120.25	122.84
3	C	2000	5I1	CAM-CAH-NAD	2.85	122.33	118.72
3	C	2000	5I1	CAI-C4-N3	2.17	118.06	114.93
3	A	2000	5I1	N1-C2-N3	-2.13	124.53	126.55
3	C	2000	5I1	NBB-C2-N1	2.11	120.24	116.65
3	C	2000	5I1	OAU-CAP-NBI	-2.02	119.94	123.63

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	2000	5I1	CAT-CBH-NBB-C2
3	C	2000	5I1	CAB-CBH-NBB-C2
3	A	2000	5I1	CBL-CBM-NAR-CBN
3	A	2000	5I1	CAT-CBH-NBB-C2
3	C	2000	5I1	N1-C2-NBB-CBH
3	A	2000	5I1	CAB-CBH-NBB-C2

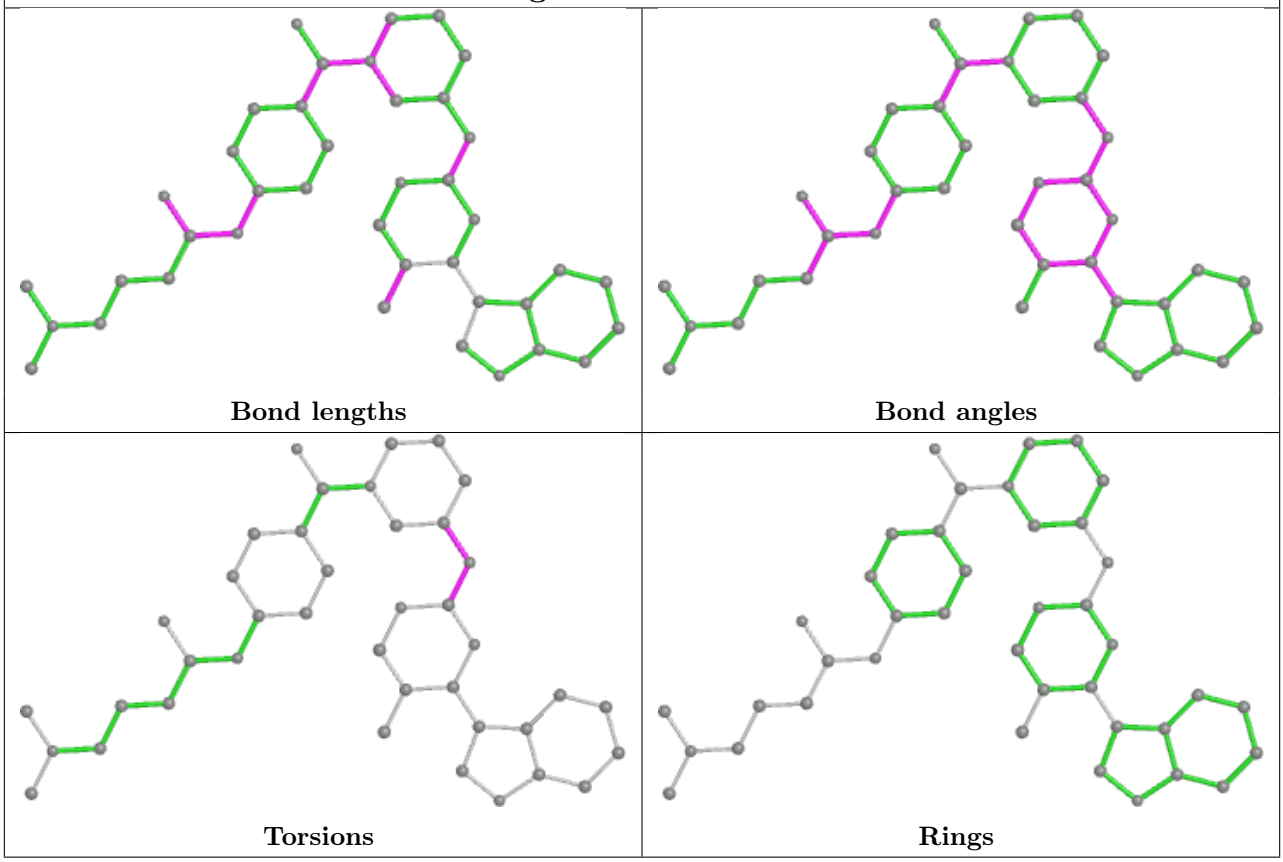
There are no ring outliers.

2 monomers are involved in 10 short contacts:

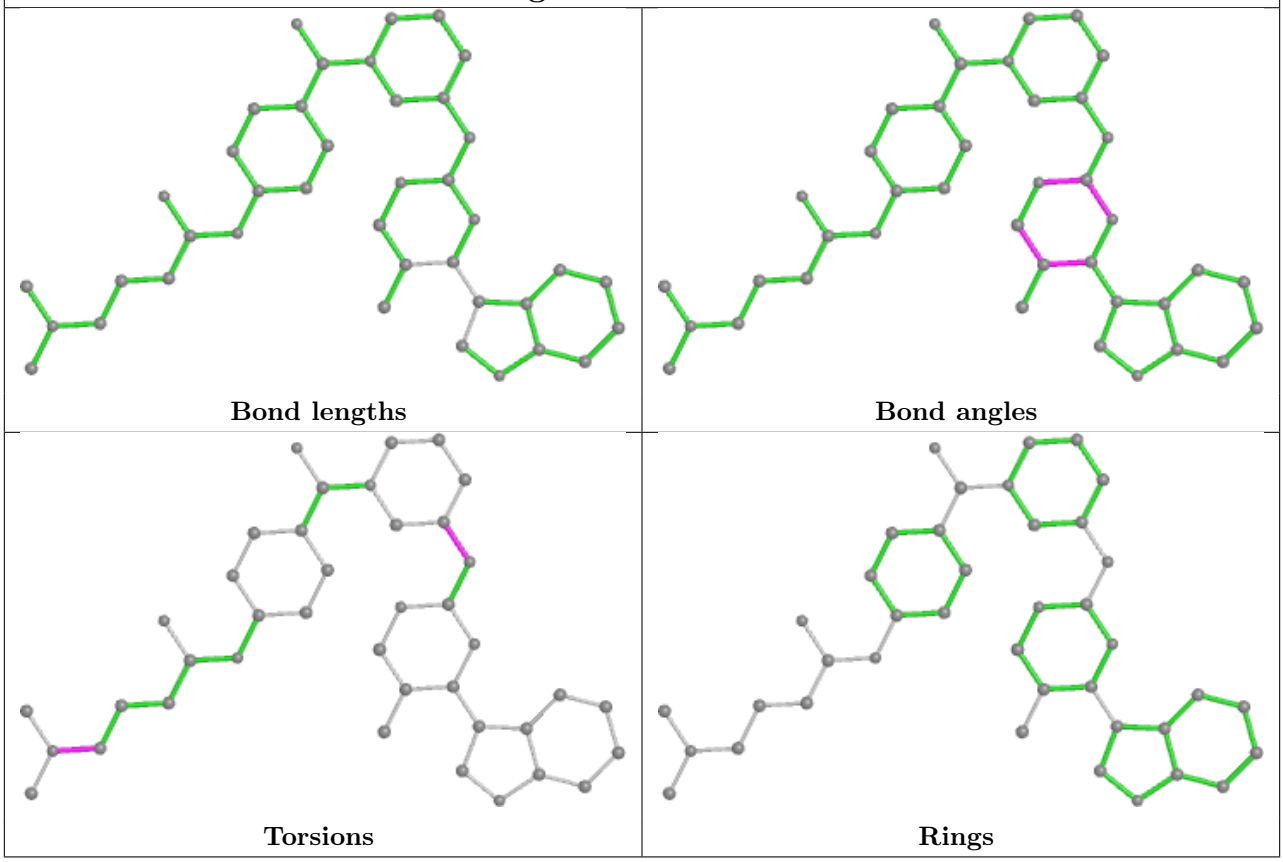
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	2000	5I1	2	0
3	A	2000	5I1	8	0

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 then 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 5I1 C 2000



Ligand 5I1 A 2000



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	324/347 (93%)	0.77	38 (11%)	4	7	57, 76, 100, 114	0
1	C	322/347 (92%)	0.56	21 (6%)	18	27	43, 60, 87, 107	0
2	B	244/268 (91%)	0.38	11 (4%)	33	46	44, 63, 89, 100	0
2	D	247/268 (92%)	0.41	8 (3%)	47	59	40, 54, 84, 100	0
All	All	1137/1230 (92%)	0.55	78 (6%)	16	24	40, 65, 95, 114	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	780	PHE	8.7
1	A	695	ILE	6.5
1	A	741	GLU	6.5
1	A	864	SER	6.2
1	A	784	LYS	6.0
1	C	778	LEU	5.9
1	C	784	LYS	5.4
1	A	783	ASP	5.2
2	B	231	TYR	4.9
1	C	777	ALA	4.9
1	A	727	THR	4.9
1	A	696	ASP	4.5
1	A	867	SER	4.4
1	C	806	LEU	4.3
1	C	779	ASP	4.3
2	B	232	ARG	4.1
1	A	728	GLY	4.1
2	D	264	GLN	4.1
2	B	227	SER	4.0
1	A	740	ASN	4.0
1	C	782	LYS	4.0

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Mol	Chain	Res	Type	RSRZ
2	D	267	HIS	4.0
1	C	1013	LEU	3.9
2	B	124	LEU	3.8
1	C	781	LYS	3.8
1	A	703	ASP	3.7
1	A	697	TRP	3.7
1	A	865	GLU	3.6
2	B	230	MET	3.6
1	C	695	ILE	3.5
1	A	868	ARG	3.4
2	B	229	PRO	3.4
2	D	173	GLY	3.4
1	C	745	PHE	3.4
1	A	700	ARG	3.3
1	A	785	GLY	3.3
1	A	1022	SER	3.2
1	A	1020	LEU	3.2
1	A	817	PHE	3.2
1	A	742	LYS	3.1
2	B	173	GLY	3.1
1	C	804	SER	3.0
1	A	698	GLY	3.0
1	A	738	LEU	2.9
1	A	699	LYS	2.9
1	C	713	GLU	2.8
1	A	726	ASP	2.8
2	D	27	ASP	2.8
2	B	171	LEU	2.8
1	A	941	VAL	2.7
1	C	739	ASP	2.6
1	A	708	ILE	2.6
1	A	1007	PRO	2.6
1	A	1010	ASP	2.6
1	A	966	PHE	2.5
1	A	1021	TRP	2.5
2	D	259	SER	2.5
2	D	261	GLY	2.5
1	C	726	ASP	2.4
1	A	782	LYS	2.4
1	C	1015	GLN	2.4
1	A	964	PHE	2.4
1	A	950	MET	2.4

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Mol	Chain	Res	Type	RSRZ
2	D	230	MET	2.3
1	A	917	ASN	2.3
1	C	776	ASP	2.3
1	A	1023	LYS	2.3
2	D	265	MET	2.3
2	B	58	PHE	2.3
1	C	746	PRO	2.3
1	A	714	GLY	2.3
1	C	738	LEU	2.2
1	C	865	GLU	2.2
1	A	956	TYR	2.2
2	B	228	LYS	2.1
1	C	783	ASP	2.1
1	A	1011	LEU	2.1
2	B	123	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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
1	TPO	A	871	11/12	0.93	0.13	71,73,82,89	0
1	TPO	C	871	11/12	0.96	0.13	54,59,64,68	0

6.3 Carbohydrates (i)

There are no monosaccharides 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	5I1	A	2000	40/40	0.85	0.20	69,75,89,90	0

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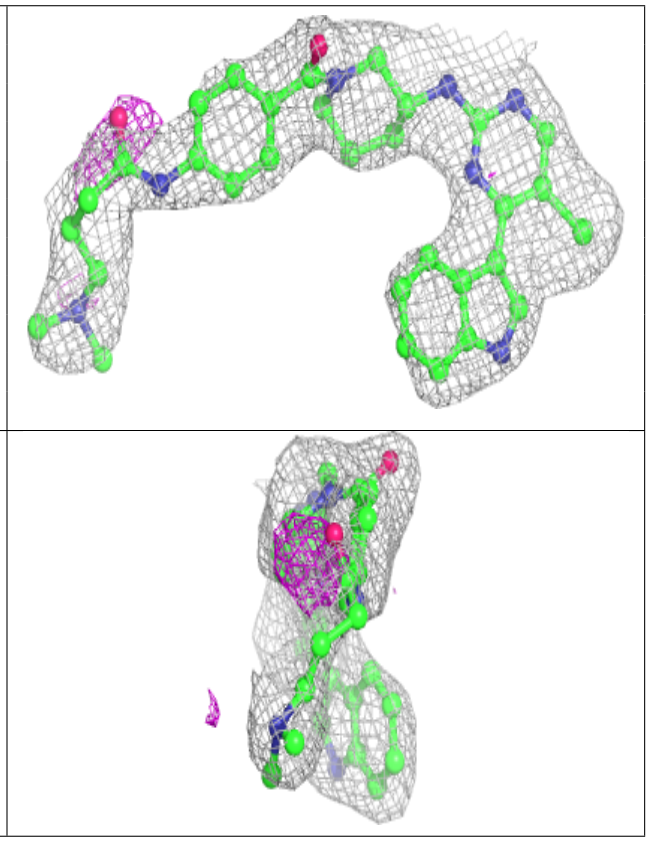
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	5I1	C	2000	40/40	0.92	0.21	50,67,99,101	0

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.

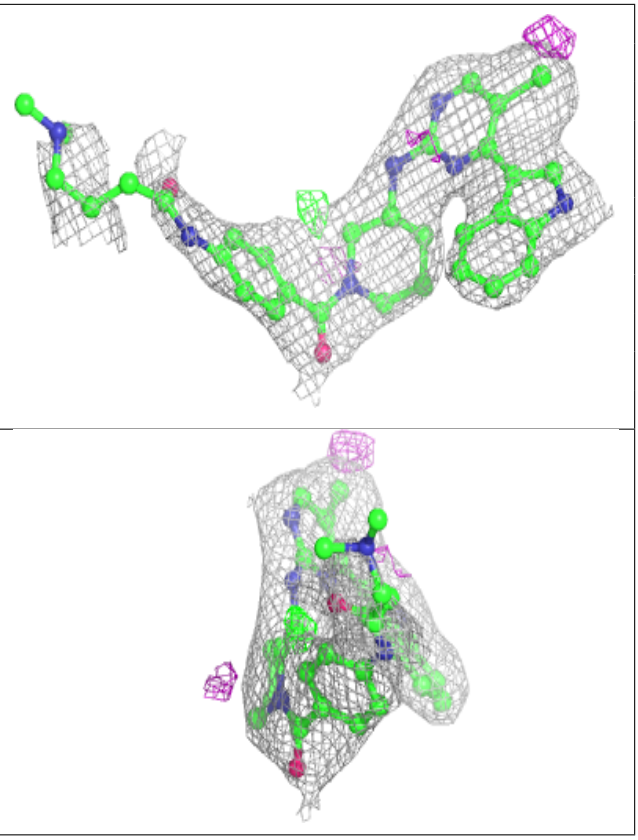
Electron density around 5I1 A 2000:

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 5I1 C 2000:

$2mF_o$-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.