



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:09 PM UTC

PDB ID : 6P7E / pdb_00006p7e
Title : Structure of T7 DNA Polymerase Bound to a Primer/Template DNA and a Peptide that Mimics the C-terminal Tail of the Primase-Helicase
Authors : Foster, B.M.; Rosenberg, D.; Salvo, H.; Stephens, K.L.; Bintz, B.J.; Hammel, M.; Ellenberger, T.; Gainey, M.D.; Wallen, J.R.
Deposited on : 2019-06-05
Resolution : 3.00 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

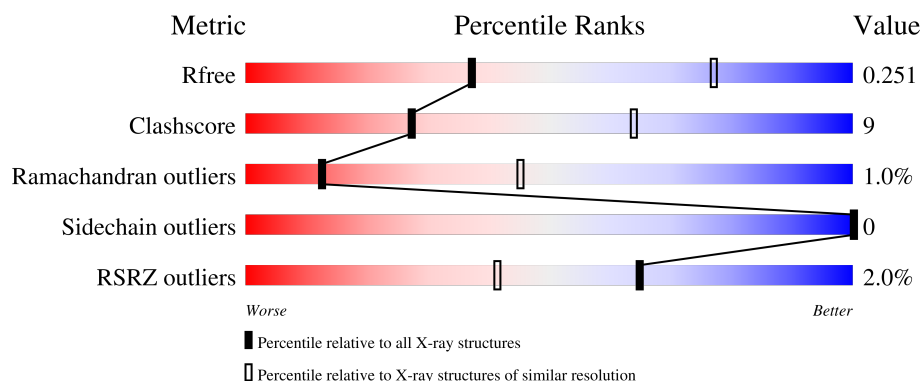
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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.



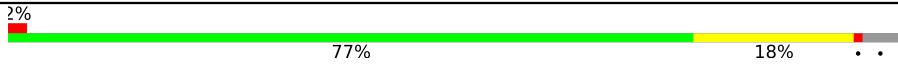
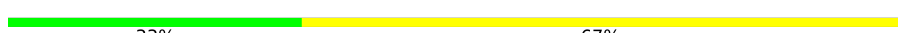
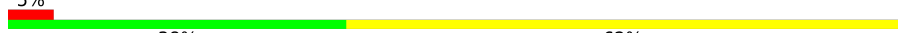
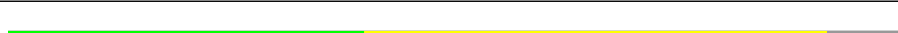
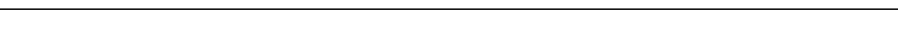
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	704	1% 80% 18%
1	B	704	80% 19%
1	C	704	2% 74% 20% 5%
1	D	704	3% 71% 20% 9%

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Mol	Chain	Length	Quality of chain
2	E	109	
2	F	109	
2	G	109	
2	H	109	
3	I	21	
3	K	21	
3	M	21	
3	O	21	
4	J	25	
4	L	25	
4	N	25	
4	P	25	
5	U	4	
5	V	4	
6	W	3	

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 28831 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 DNA-directed DNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	694	Total	C	N	O	S	0	0	0
			5531	3524	963	1020	24			
1	B	697	Total	C	N	O	S	0	0	0
			5555	3537	969	1025	24			
1	C	672	Total	C	N	O	S	0	0	0
			5368	3417	936	992	23			
1	D	644	Total	C	N	O	S	0	0	0
			5140	3272	894	953	21			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	5	ALA	ASP	engineered mutation	UNP P00581
A	7	ALA	GLU	engineered mutation	UNP P00581
B	5	ALA	ASP	engineered mutation	UNP P00581
B	7	ALA	GLU	engineered mutation	UNP P00581
C	5	ALA	ASP	engineered mutation	UNP P00581
C	7	ALA	GLU	engineered mutation	UNP P00581
D	5	ALA	ASP	engineered mutation	UNP P00581
D	7	ALA	GLU	engineered mutation	UNP P00581

- Molecule 2 is a protein called TrxA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	105	Total	C	N	O	S	0	0	0
			802	518	129	152	3			
2	F	105	Total	C	N	O	S	0	0	0
			802	518	129	152	3			
2	G	105	Total	C	N	O	S	0	0	0
			802	518	129	152	3			
2	H	105	Total	C	N	O	S	0	0	0
			802	518	129	152	3			

- Molecule 3 is a DNA chain called DNA (5'-D(P*GP*GP*CP*AP*GP*GP*TP*GP*GP*TP*CP*TP*TP*GP*CP*CP*GP*GP*TP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	I	21	Total	C	N	O	P	0	0	0
			437	206	82	128	21			
3	K	21	Total	C	N	O	P	0	0	0
			437	206	82	128	21			
3	M	21	Total	C	N	O	P	0	0	0
			437	206	82	128	21			
3	O	21	Total	C	N	O	P	0	0	0
			437	206	82	128	21			

- Molecule 4 is a DNA chain called DNA (25-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	J	25	Total	C	N	O	P	0	0	0
			501	237	93	146	25			
4	L	23	Total	C	N	O	P	0	0	0
			463	219	87	134	23			
4	N	24	Total	C	N	O	P	0	0	0
			482	228	90	140	24			
4	P	23	Total	C	N	O	P	0	0	0
			463	219	87	134	23			

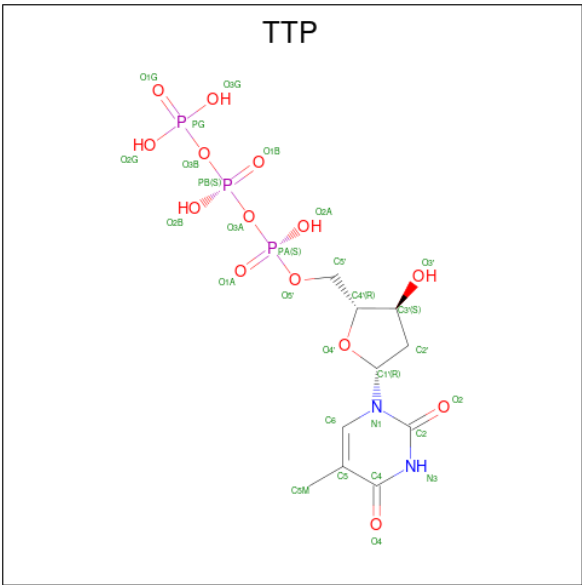
- Molecule 5 is a protein called ASP-THR-ASP-PHE peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	U	4	Total	C	N	O	0	0	0
			35	21	4	10			
5	V	4	Total	C	N	O	0	0	0
			35	21	4	10			

- Molecule 6 is a protein called THR-ASP-PHE peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	W	3	Total	C	N	O	0	0	0
			27	17	3	7			

- Molecule 7 is THYMIDINE-5'-TRIPHOSPHATE (CCD ID: TTP) (formula: C₁₀H₁₇N₂O₁₄P₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
7	A	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
7	B	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
7	B	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
7	C	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
7	C	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
7	D	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
7	D	1	Total	C	N	O	P	0	0
			29	10	2	14	3		

- Molecule 8 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

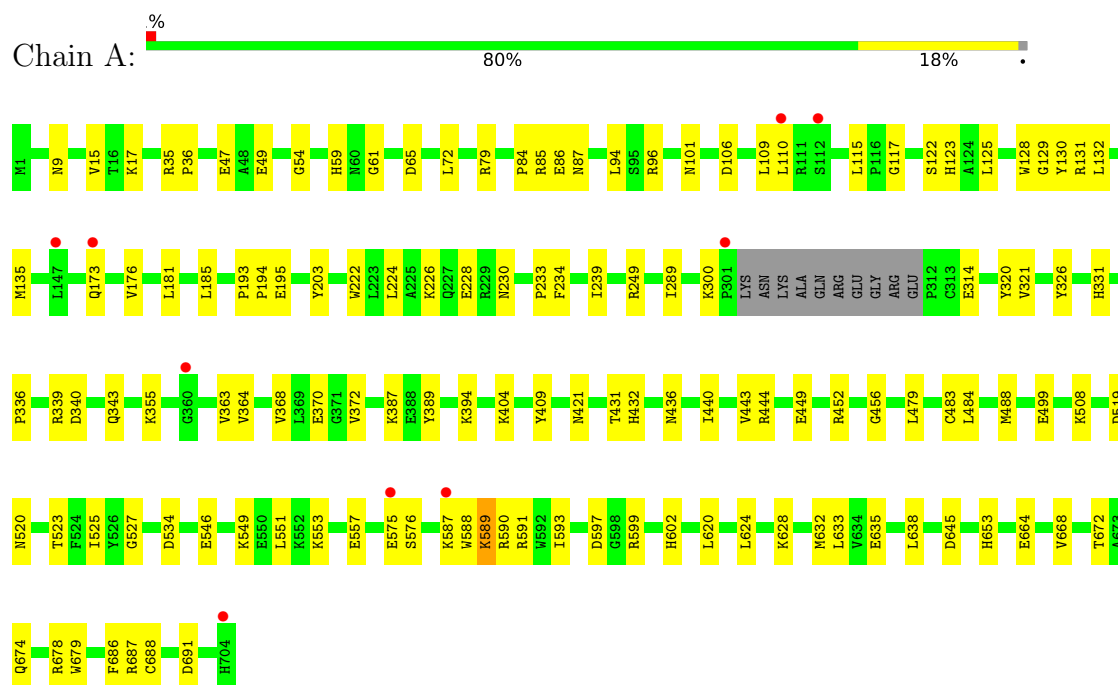
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	15	Total 15	O 15	0	0
9	B	7	Total 7	O 7	0	0
9	C	1	Total 1	O 1	0	0
9	D	3	Total 3	O 3	0	0
9	J	2	Total 2	O 2	0	0
9	K	2	Total 2	O 2	0	0
9	L	1	Total 1	O 1	0	0

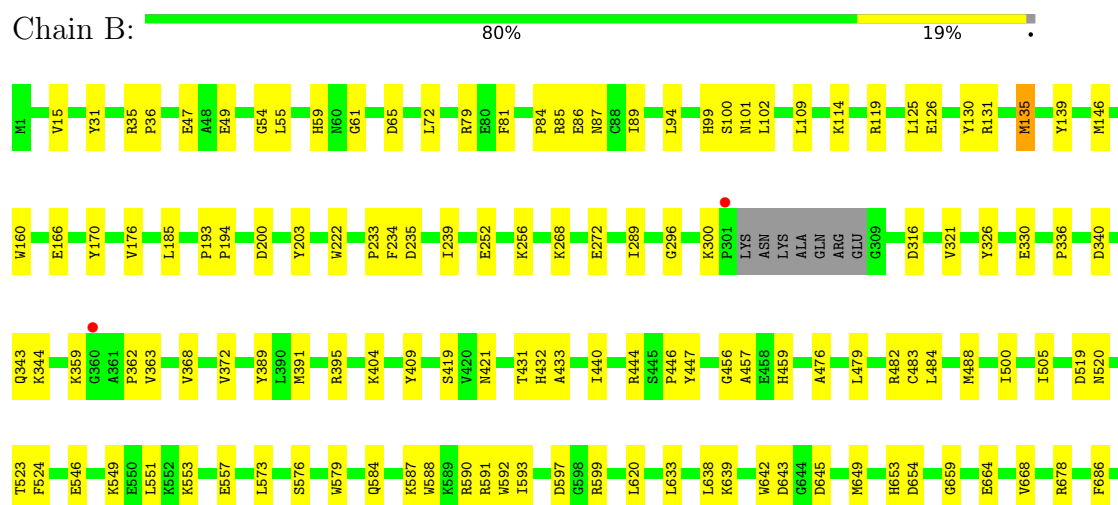
3 Residue-property plots

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: DNA-directed DNA polymerase

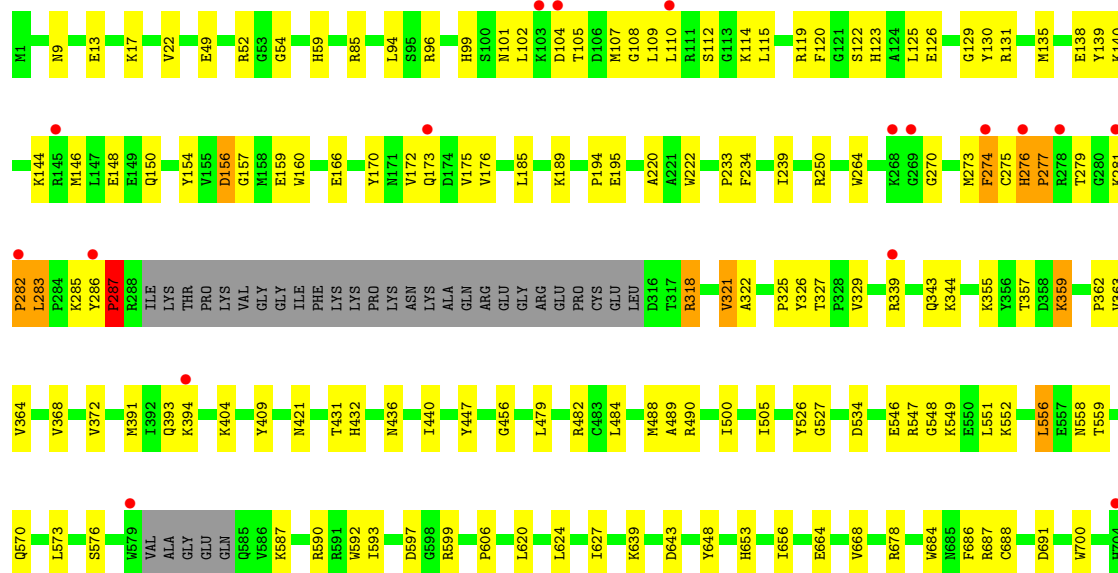
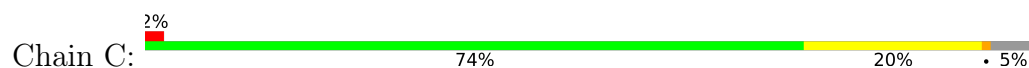


• Molecule 1: DNA-directed DNA polymerase

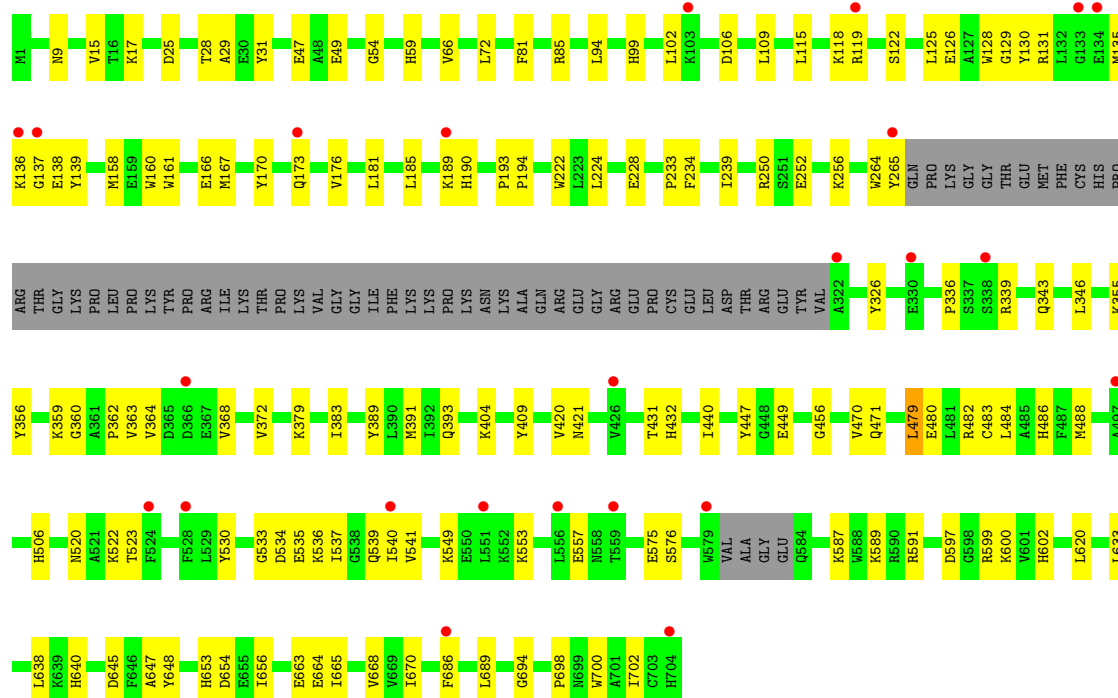




• Molecule 1: DNA-directed DNA polymerase



• Molecule 1: DNA-directed DNA polymerase



- Molecule 2: TrxA

Chain E: 



- Molecule 2: TrxA

Chain F: 



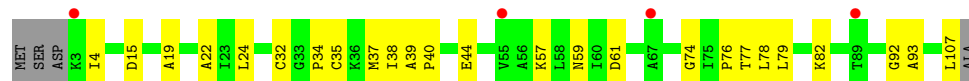
- Molecule 2: TrxA

Chain G: 



- Molecule 2: TrxA

Chain H: 



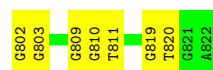
- Molecule 3: DNA (5'-D(P*GP*GP*CP*AP*GP*GP*TP*GP*GP*TP*CP*TP*TP*GP*CP*C
P*GP*GP*TP*GP*A)-3')

Chain I: 



- Molecule 3: DNA (5'-D(P*GP*GP*CP*AP*GP*GP*TP*GP*GP*TP*CP*TP*TP*GP*CP*C
P*GP*GP*TP*GP*A)-3')

Chain K: 



- Molecule 3: DNA (5'-D(P*GP*GP*CP*AP*GP*GP*TP*GP*GP*TP*CP*TP*TP*GP*CP*C
P*GP*GP*TP*GP*A)-3')

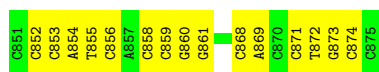
Chain M: 



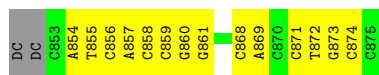
- Molecule 3: DNA (5'-D(P*GP*GP*CP*AP*GP*GP*TP*GP*GP*TP*CP*TP*TP*GP*CP*C
P*GP*GP*TP*GP*A)-3')



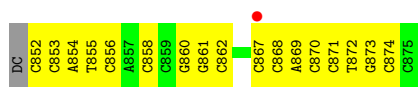
- Molecule 4: DNA (25-MER)



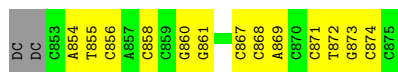
- Molecule 4: DNA (25-MER)



- Molecule 4: DNA (25-MER)



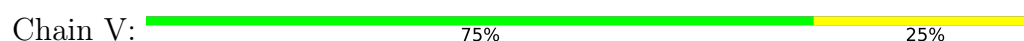
- Molecule 4: DNA (25-MER)



- Molecule 5: ASP-THR-ASP-PHE peptide



- Molecule 5: ASP-THR-ASP-PHE peptide





- Molecule 6: THR-ASP-PHE peptide

Chain W:  100%

There are no outlier residues recorded for this chain.

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	100.72Å 102.70Å 148.88Å 91.36° 96.83° 113.11°	Depositor
Resolution (Å)	49.13 – 3.00 49.13 – 3.00	Depositor EDS
% Data completeness (in resolution range)	98.8 (49.13-3.00) 98.7 (49.13-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.40 (at 3.01Å)	Xtriage
Refinement program	PHENIX (1.14_3260)	Depositor
R, R_{free}	0.214 , 0.248 0.215 , 0.251	Depositor DCC
R_{free} test set	5358 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	58.6	Xtriage
Anisotropy	0.106	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 47.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	28831	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.53% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: TTP, MG

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	0.26	0/5669	0.72	2/7668 (0.0%)
1	B	0.25	0/5693	0.72	0/7700
1	C	0.25	0/5501	0.75	3/7441 (0.0%)
1	D	0.23	0/5264	0.71	0/7120
2	E	0.25	0/817	0.77	1/1108 (0.1%)
2	F	0.33	0/817	0.71	0/1108
2	G	0.20	0/817	0.68	0/1108
2	H	0.21	0/817	0.68	0/1108
3	I	0.36	0/490	0.63	0/756
3	K	0.37	0/490	0.69	0/756
3	M	0.33	0/490	0.64	0/756
3	O	0.35	0/490	0.62	0/756
4	J	0.40	0/560	0.61	0/858
4	L	0.36	0/518	0.57	0/794
4	N	0.34	0/539	0.53	0/826
4	P	0.35	0/518	0.59	0/794
5	U	0.27	0/35	0.41	0/45
5	V	0.22	0/35	0.51	0/45
6	W	0.13	0/27	0.25	0/34
All	All	0.27	0/29587	0.71	6/40781 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	156	ASP	CA-C-N	6.39	133.93	121.41
1	C	156	ASP	C-N-CA	6.39	133.93	121.41
1	A	589	LYS	N-CA-C	-6.26	103.99	111.69
1	A	331	HIS	N-CA-C	-6.02	99.42	109.24
1	C	556	LEU	N-CA-C	-5.47	102.80	110.35
2	E	19	ALA	N-CA-C	5.45	122.40	110.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	287	PRO	Peptide

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	5531	0	5432	84	0
1	B	5555	0	5453	87	0
1	C	5368	0	5251	113	0
1	D	5140	0	5024	97	0
2	E	802	0	816	15	0
2	F	802	0	816	22	0
2	G	802	0	816	26	0
2	H	802	0	816	17	0
3	I	437	0	235	5	0
3	K	437	0	235	6	0
3	M	437	0	235	9	0
3	O	437	0	235	8	0
4	J	501	0	278	12	0
4	L	463	0	256	10	0
4	N	482	0	267	12	0
4	P	463	0	256	11	0
5	U	35	0	26	4	0
5	V	35	0	26	1	0
6	W	27	0	22	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	58	0	26	0	0
7	B	58	0	26	1	0
7	C	58	0	26	1	0
7	D	58	0	26	3	0
8	A	3	0	0	0	0
8	B	3	0	0	0	0
8	C	3	0	0	0	0
8	D	3	0	0	0	0
9	A	15	0	0	0	0
9	B	7	0	0	0	0
9	C	1	0	0	0	0
9	D	3	0	0	0	0
9	J	2	0	0	0	0
9	K	2	0	0	0	0
9	L	1	0	0	0	0
All	All	28831	0	26599	502	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 9.

All (502) 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:C:273:MET:HG3	1:C:281:LYS:HZ1	1.34	0.93
3:O:803:DG:H2"	3:O:804:DC:H5"	1.52	0.92
1:C:250:ARG:CD	1:C:393:GLN:OE1	2.21	0.88
1:B:391:MET:HE1	1:B:447:TYR:HD2	1.40	0.87
1:C:273:MET:HG2	1:C:274:PHE:H	1.40	0.86
4:P:867:DC:H2"	4:P:868:DC:H5"	1.59	0.85
1:C:250:ARG:HD2	1:C:393:GLN:OE1	1.78	0.84
2:H:4:ILE:HD11	2:H:57:LYS:HG3	1.60	0.83
1:C:115:LEU:HD21	1:C:131:ARG:HG3	1.63	0.79
4:N:867:DC:H2"	4:N:868:DC:H5"	1.64	0.79
1:C:281:LYS:HE2	2:G:34:PRO:HD3	1.65	0.77
1:A:339:ARG:NH1	1:A:364:VAL:O	2.17	0.77
1:D:119:ARG:NH2	1:D:126:GLU:OE1	2.18	0.77
1:B:296:GLY:HA2	1:B:321:VAL:HG23	1.66	0.77
4:N:852:DC:H2"	4:N:853:DC:H5"	1.65	0.77
1:B:94:LEU:HB3	1:B:185:LEU:HD13	1.67	0.77
1:A:300:LYS:HG3	4:J:874:DC:H5"	1.67	0.76
1:B:391:MET:HE1	1:B:447:TYR:CD2	2.22	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:590:ARG:NH2	5:U:4:PHE:O	2.19	0.74
4:J:873:DG:H2''	4:J:874:DC:O5'	1.87	0.74
3:O:805:DA:N6	4:P:872:DT:O4	2.20	0.74
1:D:129:GLY:HA3	1:D:135:MET:HE3	1.69	0.74
1:C:250:ARG:HD3	1:C:393:GLN:OE1	1.87	0.73
1:D:94:LEU:HB3	1:D:185:LEU:HD13	1.70	0.73
1:B:272:GLU:HB2	1:B:289:ILE:HG22	1.71	0.73
1:C:189:LYS:HD2	1:C:194:PRO:HG3	1.71	0.73
4:P:873:DG:H2''	4:P:874:DC:O5'	1.88	0.73
4:N:868:DC:H2''	4:N:869:DA:C8	2.24	0.72
3:I:806:DG:H2''	3:I:807:DG:O5'	1.89	0.71
4:N:873:DG:H2''	4:N:874:DC:O5'	1.90	0.71
1:A:122:SER:O	1:A:131:ARG:NH1	2.22	0.71
1:C:286:TYR:HE1	1:C:318:ARG:HG3	1.56	0.71
1:C:343:GLN:HG3	1:C:362:PRO:HG2	1.72	0.70
1:D:250:ARG:HD2	1:D:393:GLN:OE1	1.92	0.70
1:B:483:CYS:HA	1:B:686:PHE:HZ	1.57	0.70
1:C:556:LEU:O	1:C:558:ASN:N	2.24	0.70
4:P:868:DC:H2''	4:P:869:DA:C8	2.28	0.69
1:C:220:ALA:HA	1:C:627:ILE:HD11	1.74	0.69
1:C:122:SER:O	1:C:131:ARG:NH1	2.24	0.69
1:C:94:LEU:HB3	1:C:185:LEU:HD13	1.75	0.68
1:C:321:VAL:HG23	2:G:90:LYS:HD2	1.75	0.68
1:A:499:GLU:OE2	1:A:508:LYS:NZ	2.26	0.68
1:B:340:ASP:OD2	1:B:344:LYS:NZ	2.27	0.67
1:C:282:PRO:HD3	1:C:326:TYR:O	1.94	0.67
1:C:556:LEU:C	1:C:558:ASN:H	2.02	0.67
3:K:802:DG:H4'	3:K:803:DG:OP1	1.93	0.67
1:C:130:TYR:CD1	1:C:135:MET:HG3	2.30	0.67
1:D:534:ASP:OD2	1:D:549:LYS:HG2	1.94	0.67
1:A:599:ARG:HD3	1:A:620:LEU:HD11	1.78	0.65
1:C:500:ILE:HG13	1:C:505:ILE:HD13	1.78	0.65
1:D:648:TYR:CE2	1:D:656:ILE:HD11	2.30	0.65
1:B:35:ARG:HB3	1:B:36:PRO:HD2	1.79	0.65
1:A:249:ARG:HH21	1:B:79:ARG:HG3	1.62	0.65
1:D:698:PRO:HD2	1:D:702:ILE:HG21	1.78	0.64
1:C:281:LYS:CE	2:G:34:PRO:HD3	2.28	0.64
1:A:109:LEU:HD22	1:A:131:ARG:HG2	1.80	0.64
1:A:195:GLU:H	1:A:195:GLU:CD	2.06	0.64
1:D:356:TYR:HA	1:D:362:PRO:HA	1.78	0.63
1:C:275:CYS:C	1:C:277:PRO:HD3	2.24	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:136:LYS:HD3	1:D:173:GLN:HG3	1.79	0.63
1:D:137:GLY:O	1:D:170:TYR:HE1	1.81	0.63
1:C:273:MET:HG3	1:C:281:LYS:NZ	2.13	0.63
1:D:339:ARG:NH1	1:D:364:VAL:O	2.31	0.63
1:C:114:LYS:NZ	1:C:130:TYR:O	2.32	0.62
2:G:96:LYS:HE2	2:G:100:LYS:HE2	1.81	0.62
1:C:326:TYR:HB3	2:G:92:GLY:HA2	1.81	0.62
1:C:101:ASN:O	1:C:105:THR:OG1	2.17	0.62
1:D:130:TYR:CD1	1:D:135:MET:HG3	2.34	0.62
4:L:873:DG:H2''	4:L:874:DC:O5'	1.98	0.62
1:D:484:LEU:O	1:D:488:MET:HG2	1.98	0.62
4:J:860:DG:H2''	4:J:861:DG:C8	2.35	0.62
1:D:533:GLY:N	4:P:854:DA:OP2	2.23	0.62
1:D:368:VAL:O	1:D:372:VAL:HG23	1.99	0.62
1:B:130:TYR:CD1	1:B:135:MET:HG3	2.34	0.62
1:D:25:ASP:O	1:D:29:ALA:N	2.33	0.61
1:A:483:CYS:HA	1:A:686:PHE:HZ	1.65	0.61
4:L:854:DA:H2'	4:L:855:DT:C6	2.36	0.61
4:N:860:DG:H2''	4:N:861:DG:C8	2.36	0.61
1:C:108:GLY:O	1:C:112:SER:OG	2.07	0.60
1:B:368:VAL:O	1:B:372:VAL:HG23	2.01	0.60
2:F:94:LEU:HD12	2:F:99:LEU:HD22	1.83	0.60
2:G:32:CYS:SG	2:G:34:PRO:HD2	2.41	0.60
1:C:234:PHE:CZ	1:C:239:ILE:HG13	2.36	0.60
1:A:484:LEU:O	1:A:488:MET:HG2	2.01	0.60
1:C:678:ARG:NH1	1:C:691:ASP:OD1	2.34	0.60
1:D:599:ARG:HD3	1:D:620:LEU:HD11	1.84	0.60
1:B:500:ILE:HG13	1:B:505:ILE:HD13	1.85	0.59
1:B:109:LEU:HD13	1:B:131:ARG:HG2	1.85	0.59
1:D:189:LYS:CE	1:D:194:PRO:HG3	2.33	0.59
2:H:15:ASP:O	2:H:19:ALA:HB2	2.02	0.59
1:A:129:GLY:HA3	1:A:135:MET:HE3	1.85	0.59
1:A:576:SER:OG	1:A:587:LYS:HD3	2.03	0.59
1:C:368:VAL:O	1:C:372:VAL:HG23	2.02	0.59
2:G:102:PHE:O	2:G:106:ASN:ND2	2.35	0.59
3:M:819:DG:C8	3:M:820:DT:H72	2.38	0.59
4:N:855:DT:H2'	4:N:856:DC:C6	2.38	0.59
1:B:484:LEU:O	1:B:488:MET:HG2	2.02	0.58
1:B:15:VAL:HG22	1:B:72:LEU:HD21	1.84	0.58
1:A:321:VAL:HG23	2:E:90:LYS:HE3	1.86	0.58
2:G:74:GLY:O	2:G:77:THR:OG1	2.22	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:320:TYR:OH	2:E:73:ARG:NH2	2.35	0.58
1:D:645:ASP:HB3	1:D:665:ILE:HD13	1.85	0.58
2:F:15:ASP:O	2:F:19:ALA:HB2	2.04	0.58
2:H:34:PRO:HA	2:H:37:MET:HE2	1.85	0.58
1:B:321:VAL:HG12	2:F:94:LEU:HD22	1.85	0.58
1:B:321:VAL:CG1	2:F:94:LEU:HD22	2.34	0.58
1:C:273:MET:HG2	1:C:274:PHE:N	2.17	0.57
2:F:94:LEU:HB2	2:F:99:LEU:HD23	1.86	0.57
1:D:122:SER:O	1:D:131:ARG:NH1	2.35	0.57
1:D:252:GLU:OE2	1:D:256:LYS:NZ	2.35	0.57
1:D:234:PHE:CZ	1:D:239:ILE:HG13	2.39	0.57
1:A:404:LYS:HA	1:A:409:TYR:HE2	1.70	0.57
1:B:114:LYS:NZ	1:B:131:ARG:O	2.38	0.57
1:D:355:LYS:HG3	1:D:363:VAL:HB	1.86	0.57
1:A:49:GLU:OE2	1:A:54:GLY:HA3	2.05	0.57
1:B:59:HIS:CD2	1:B:125:LEU:HG	2.40	0.56
1:A:687:ARG:HH22	5:U:3:ASP:CG	2.13	0.56
4:L:860:DG:H2''	4:L:861:DG:C8	2.40	0.56
4:P:860:DG:H2''	4:P:861:DG:C8	2.39	0.56
1:B:457:ALA:HA	1:B:649:MET:HE1	1.87	0.56
1:C:287:PRO:HB3	1:C:321:VAL:HA	1.87	0.56
1:C:287:PRO:HB3	1:C:322:ALA:H	1.69	0.56
1:C:363:VAL:O	1:C:368:VAL:HG21	2.05	0.56
2:E:19:ALA:O	2:E:21:GLY:N	2.31	0.56
2:F:94:LEU:HB2	2:F:99:LEU:CD2	2.36	0.56
1:B:99:HIS:C	1:B:101:ASN:H	2.13	0.56
1:C:484:LEU:O	1:C:488:MET:HG2	2.06	0.56
1:C:146:MET:HE1	1:C:166:GLU:HG2	1.87	0.56
2:G:4:ILE:HG21	2:G:57:LYS:HG3	1.88	0.56
1:A:628:LYS:HG3	1:A:679:TRP:CE3	2.40	0.55
1:C:590:ARG:HD2	1:C:592:TRP:O	2.05	0.55
1:B:139:TYR:HB2	1:B:170:TYR:CG	2.41	0.55
1:C:59:HIS:CD2	1:C:125:LEU:HG	2.41	0.55
1:B:590:ARG:HD2	1:B:592:TRP:O	2.07	0.55
2:G:17:LEU:HA	2:G:84:GLY:HA2	1.86	0.55
1:C:109:LEU:HD13	1:C:131:ARG:HG2	1.89	0.55
1:B:642:TRP:HH2	1:B:649:MET:HE2	1.71	0.55
1:A:132:LEU:HD11	1:A:181:LEU:HD13	1.88	0.55
1:A:432:HIS:O	4:J:858:DC:H4'	2.06	0.54
1:C:233:PRO:HB2	1:C:456:GLY:O	2.07	0.54
1:D:420:VAL:HG12	1:D:432:HIS:CD2	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:593:ILE:HG23	5:V:4:PHE:HB2	1.89	0.54
1:C:421:ASN:HB3	1:C:431:THR:OG1	2.08	0.54
1:C:556:LEU:C	1:C:558:ASN:N	2.66	0.54
2:E:77:THR:HG22	2:E:79:LEU:HD13	1.89	0.54
1:B:86:GLU:HG3	1:B:203:TYR:CE1	2.43	0.54
1:D:115:LEU:HD13	1:D:119:ARG:O	2.08	0.54
1:D:189:LYS:HE2	1:D:194:PRO:HG3	1.90	0.54
1:C:85:ARG:HG3	1:C:222:TRP:CG	2.42	0.54
1:C:150:GLN:OE1	1:C:150:GLN:N	2.41	0.54
2:F:90:LYS:HG2	2:F:94:LEU:HD11	1.90	0.54
1:A:368:VAL:O	1:A:372:VAL:HG23	2.09	0.53
1:C:599:ARG:HD3	1:C:620:LEU:HD11	1.90	0.53
1:D:326:TYR:HB3	2:H:92:GLY:HA2	1.89	0.53
1:A:590:ARG:HH22	5:U:4:PHE:C	2.17	0.53
3:O:809:DG:H4'	3:O:810:DG:OP1	2.09	0.53
1:C:339:ARG:NH1	1:C:364:VAL:O	2.41	0.53
3:I:810:DG:H2'	3:I:811:DT:C6	2.44	0.53
1:C:264:TRP:NE1	1:C:344:LYS:HE2	2.24	0.53
3:M:809:DG:H2''	3:M:810:DG:O5'	2.08	0.53
4:L:855:DT:H2'	4:L:856:DC:C6	2.44	0.53
4:P:871:DC:H2'	4:P:872:DT:C6	2.43	0.53
1:C:489:ALA:HB3	1:C:687:ARG:HE	1.73	0.52
4:N:871:DC:H2'	4:N:872:DT:C6	2.44	0.52
1:D:49:GLU:OE2	1:D:54:GLY:HA3	2.09	0.52
1:A:233:PRO:HB2	1:A:456:GLY:O	2.09	0.52
2:E:15:ASP:O	2:E:19:ALA:HB2	2.09	0.52
4:J:859:DC:H2''	4:J:860:DG:H8	1.75	0.52
4:P:854:DA:H2'	4:P:855:DT:C6	2.44	0.52
1:B:268:LYS:HE3	1:B:330:GLU:OE1	2.09	0.52
1:D:106:ASP:OD1	1:D:109:LEU:HD12	2.10	0.52
4:N:854:DA:H2'	4:N:855:DT:C6	2.44	0.52
1:C:648:TYR:CE2	1:C:656:ILE:HD11	2.45	0.52
2:G:34:PRO:HB3	2:G:93:ALA:HB2	1.92	0.52
1:B:573:LEU:HD11	1:B:593:ILE:HD11	1.90	0.52
1:C:432:HIS:O	4:N:858:DC:H4'	2.10	0.52
1:A:336:PRO:HB2	1:A:389:TYR:CD1	2.45	0.52
1:C:195:GLU:CD	1:C:195:GLU:H	2.17	0.52
2:H:77:THR:HG22	2:H:79:LEU:HD13	1.92	0.52
1:B:146:MET:HE1	1:B:166:GLU:HG2	1.92	0.51
1:D:480:GLU:OE1	1:D:530:TYR:OH	2.16	0.51
2:G:49:TYR:HB3	2:G:53:LEU:HG	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:871:DC:H2'	4:L:872:DT:C6	2.46	0.51
1:A:226:LYS:NZ	1:A:230:ASN:OD1	2.41	0.51
1:B:252:GLU:OE2	1:B:256:LYS:NZ	2.43	0.51
1:D:99:HIS:HB3	1:D:102:LEU:HD21	1.93	0.51
1:C:139:TYR:HB2	1:C:170:TYR:HD1	1.75	0.51
1:C:500:ILE:HG13	1:C:505:ILE:CD1	2.40	0.51
1:D:421:ASN:HB3	1:D:431:THR:OG1	2.10	0.51
1:B:404:LYS:HA	1:B:409:TYR:HE2	1.76	0.51
1:B:546:GLU:H	1:B:546:GLU:CD	2.19	0.51
1:B:49:GLU:OE2	1:B:54:GLY:HA3	2.10	0.51
1:B:326:TYR:HB3	2:F:92:GLY:HA2	1.93	0.51
1:B:336:PRO:HG2	1:B:389:TYR:CE1	2.46	0.51
1:B:233:PRO:HB2	1:B:456:GLY:O	2.10	0.51
1:B:599:ARG:HD3	1:B:620:LEU:HD11	1.92	0.51
2:E:32:CYS:SG	2:E:34:PRO:HD2	2.51	0.51
2:F:77:THR:HG22	2:F:79:LEU:HD13	1.93	0.51
1:A:84:PRO:HG2	1:A:87:ASN:OD1	2.11	0.50
1:A:195:GLU:CD	1:A:195:GLU:N	2.69	0.50
1:C:273:MET:CG	1:C:274:PHE:H	2.18	0.50
2:G:77:THR:HG22	2:G:79:LEU:HD13	1.94	0.50
1:A:575:GLU:HB2	1:A:589:LYS:HG3	1.92	0.50
1:C:326:TYR:HA	2:G:34:PRO:HG2	1.92	0.50
1:D:173:GLN:O	1:D:176:VAL:HG22	2.11	0.50
1:D:470:VAL:HG21	1:D:663:GLU:HG2	1.93	0.50
2:F:103:LEU:O	2:F:107:LEU:N	2.44	0.50
2:G:38:ILE:HD12	2:G:93:ALA:HA	1.93	0.50
1:B:483:CYS:HA	1:B:686:PHE:CZ	2.41	0.50
1:C:546:GLU:CD	1:C:546:GLU:H	2.20	0.50
1:C:576:SER:OG	1:C:587:LYS:HD2	2.12	0.50
4:J:854:DA:H2'	4:J:855:DT:C6	2.47	0.50
2:F:17:LEU:HA	2:F:84:GLY:HA2	1.92	0.50
2:H:74:GLY:O	2:H:77:THR:OG1	2.28	0.50
1:A:588:TRP:CE2	1:A:591:ARG:HD2	2.46	0.50
1:C:404:LYS:HA	1:C:409:TYR:HE2	1.77	0.49
2:H:22:ALA:HB2	2:H:107:LEU:HD21	1.93	0.49
1:A:94:LEU:HB3	1:A:185:LEU:HD13	1.94	0.49
1:C:129:GLY:HA3	1:C:135:MET:HE3	1.92	0.49
1:D:483:CYS:HA	1:D:686:PHE:HZ	1.75	0.49
1:D:540:ILE:HG13	1:D:541:VAL:HG23	1.94	0.49
4:L:859:DC:H2''	4:L:860:DG:C8	2.47	0.49
4:L:859:DC:H2''	4:L:860:DG:H8	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:326:TYR:HB3	2:E:92:GLY:HA2	1.93	0.49
1:A:633:LEU:HD22	1:A:638:LEU:HD12	1.94	0.49
1:B:649:MET:HE3	1:B:659:GLY:HA3	1.94	0.49
1:C:355:LYS:HG3	1:C:363:VAL:HB	1.94	0.49
1:C:527:GLY:HA2	4:N:854:DA:C4	2.47	0.49
1:C:329:VAL:HG11	2:G:31:TRP:HH2	1.77	0.49
1:B:444:ARG:HD2	1:B:519:ASP:OD2	2.12	0.49
1:C:49:GLU:OE2	1:C:54:GLY:HA3	2.13	0.49
1:D:47:GLU:HG3	1:D:81:PHE:HE1	1.78	0.49
1:D:166:GLU:OE1	1:D:166:GLU:N	2.45	0.49
1:B:363:VAL:O	1:B:368:VAL:HG21	2.11	0.49
1:D:189:LYS:HE3	1:D:194:PRO:HG3	1.94	0.49
1:D:553:LYS:O	1:D:557:GLU:HG2	2.13	0.49
1:C:101:ASN:OD1	1:C:105:THR:HG23	2.13	0.49
1:A:597:ASP:OD1	1:A:599:ARG:HD2	2.12	0.49
1:C:157:GLY:HA2	1:C:159:GLU:OE1	2.12	0.49
1:D:575:GLU:HB3	1:D:589:LYS:HE2	1.95	0.49
2:F:74:GLY:O	2:F:77:THR:OG1	2.31	0.49
3:K:802:DG:H2''	3:K:803:DG:C8	2.48	0.49
3:O:806:DG:H2''	3:O:807:DG:O5'	2.13	0.49
1:D:379:LYS:O	1:D:383:ILE:HG13	2.13	0.49
1:A:15:VAL:HG22	1:A:72:LEU:HD21	1.95	0.48
1:A:421:ASN:HB3	1:A:431:THR:OG1	2.12	0.48
1:D:404:LYS:HA	1:D:409:TYR:HE2	1.78	0.48
1:D:420:VAL:HG12	1:D:432:HIS:HD2	1.78	0.48
1:D:591:ARG:HG2	1:D:602:HIS:CE1	2.47	0.48
2:F:12:PHE:CD2	2:F:66:THR:HG21	2.47	0.48
1:B:160:TRP:CZ2	7:B:801:TTP:H2'2	2.47	0.48
1:C:276:HIS:N	1:C:277:PRO:HD3	2.28	0.48
1:B:84:PRO:HG2	1:B:87:ASN:OD1	2.13	0.48
1:B:432:HIS:O	4:L:858:DC:H4'	2.13	0.48
2:H:59:ASN:OD1	2:H:61:ASP:HB2	2.13	0.48
1:C:482:ARG:HG2	1:C:500:ILE:HD13	1.95	0.48
1:D:126:GLU:CD	1:D:138:GLU:HG2	2.38	0.48
4:L:868:DC:H2''	4:L:869:DA:C8	2.48	0.48
1:C:488:MET:HE1	1:C:559:THR:HB	1.96	0.48
1:B:55:LEU:HD13	1:B:89:ILE:HD11	1.95	0.48
2:E:23:ILE:HD13	2:E:54:THR:HB	1.94	0.48
2:G:40:PRO:O	2:G:44:GLU:HG3	2.14	0.48
1:D:535:GLU:O	1:D:539:GLN:HG3	2.12	0.48
1:A:173:GLN:O	1:A:176:VAL:HG22	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:39:ALA:HB3	2:E:40:PRO:HD3	1.96	0.48
3:K:810:DG:H2'	3:K:811:DT:C6	2.48	0.48
1:A:686:PHE:HD2	1:A:688:CYS:O	1.97	0.48
1:D:190:HIS:CG	1:D:600:LYS:HE3	2.48	0.48
1:D:440:ILE:HD13	1:D:700:TRP:CZ3	2.49	0.48
1:A:363:VAL:O	1:A:368:VAL:HG21	2.14	0.47
1:C:283:LEU:HD23	1:C:325:PRO:HB3	1.95	0.47
1:C:624:LEU:HD12	1:C:684:TRP:CZ2	2.49	0.47
1:D:506:HIS:CG	1:D:522:LYS:HG2	2.49	0.47
2:F:39:ALA:HB3	2:F:40:PRO:HD3	1.95	0.47
1:B:391:MET:HE3	1:B:446:PRO:HB2	1.95	0.47
1:B:597:ASP:OD2	1:B:599:ARG:NH1	2.45	0.47
1:C:286:TYR:N	1:C:322:ALA:HB2	2.29	0.47
1:C:490:ARG:HG3	1:C:490:ARG:HH11	1.79	0.47
2:H:40:PRO:O	2:H:44:GLU:HG3	2.14	0.47
1:B:576:SER:HB3	1:B:587:LYS:HD3	1.95	0.47
1:C:126:GLU:CD	1:C:138:GLU:HG2	2.39	0.47
1:D:160:TRP:CZ2	7:D:801:TTP:H2'2	2.48	0.47
1:B:524:PHE:CE1	1:B:551:LEU:HD13	2.48	0.47
1:D:537:ILE:HA	1:D:540:ILE:HG12	1.96	0.47
1:C:139:TYR:HB2	1:C:170:TYR:CD1	2.49	0.47
1:C:172:VAL:O	1:C:176:VAL:HG13	2.14	0.47
2:G:39:ALA:HB3	2:G:40:PRO:HD3	1.97	0.47
1:A:632:MET:HA	1:A:635:GLU:HG2	1.96	0.47
1:B:686:PHE:CE2	1:B:690:LEU:HD21	2.49	0.47
1:C:357:THR:O	1:C:359:LYS:N	2.44	0.47
1:A:234:PHE:CZ	1:A:239:ILE:HG13	2.50	0.47
4:J:852:DC:H4'	4:J:853:DC:OP2	2.14	0.47
4:J:871:DC:H2'	4:J:872:DT:C6	2.49	0.47
1:B:234:PHE:CZ	1:B:239:ILE:HG13	2.50	0.47
1:B:359:LYS:O	1:B:359:LYS:HG2	2.15	0.47
4:J:859:DC:H2''	4:J:860:DG:C8	2.50	0.47
1:A:678:ARG:NH1	1:A:691:ASP:OD1	2.48	0.47
1:B:639:LYS:NZ	1:B:643:ASP:OD1	2.48	0.47
1:D:343:GLN:HG3	1:D:362:PRO:CG	2.45	0.47
1:D:480:GLU:CD	7:D:805:TTP:H2'1	2.40	0.47
3:I:813:DT:H2''	3:I:814:DT:H5''	1.96	0.47
1:A:128:TRP:CD2	1:A:181:LEU:HD11	2.50	0.47
1:C:9:ASN:HA	1:C:160:TRP:O	2.14	0.46
1:C:22:VAL:HB	1:C:175:VAL:HG21	1.96	0.46
1:D:99:HIS:O	1:D:102:LEU:HG	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:482:ARG:NH1	1:D:689:LEU:O	2.45	0.46
2:H:82:LYS:HE3	2:H:107:LEU:HD23	1.97	0.46
1:A:432:HIS:HB2	1:A:436:ASN:OD1	2.15	0.46
1:A:546:GLU:CD	1:A:546:GLU:H	2.22	0.46
1:A:130:TYR:CG	1:A:135:MET:HG3	2.50	0.46
1:C:624:LEU:HD12	1:C:684:TRP:CH2	2.51	0.46
2:E:27:PHE:CE1	2:E:79:LEU:HD22	2.50	0.46
1:A:59:HIS:CD2	1:A:125:LEU:HG	2.50	0.46
1:B:235:ASP:HB2	1:B:459:HIS:CE1	2.50	0.46
1:C:96:ARG:HA	1:C:123:HIS:CE1	2.50	0.46
2:H:39:ALA:HB3	2:H:40:PRO:HD3	1.97	0.46
4:J:868:DC:H2''	4:J:869:DA:C8	2.51	0.46
1:A:340:ASP:O	1:A:343:GLN:HG3	2.15	0.46
1:A:444:ARG:HD2	1:A:519:ASP:OD2	2.15	0.46
1:C:440:ILE:HD13	1:C:700:TRP:CZ3	2.51	0.46
1:A:101:ASN:HD22	1:A:602:HIS:CG	2.34	0.46
1:D:537:ILE:O	1:D:540:ILE:HG12	2.15	0.46
1:D:264:TRP:HD1	1:D:265:TYR:CD1	2.34	0.46
1:C:639:LYS:NZ	1:C:643:ASP:OD1	2.49	0.46
2:H:35:CYS:HA	2:H:76:PRO:HG3	1.98	0.46
1:A:370:GLU:OE1	1:A:387:LYS:HE3	2.15	0.45
1:B:678:ARG:NH1	1:B:691:ASP:OD1	2.49	0.45
1:A:527:GLY:HA2	4:J:854:DA:C4	2.51	0.45
1:D:189:LYS:HD2	1:D:189:LYS:N	2.31	0.45
1:D:336:PRO:HG2	1:D:389:TYR:CE1	2.51	0.45
2:G:7:LEU:HD13	2:G:12:PHE:CD2	2.51	0.45
3:O:821:DG:H2'	3:O:822:DA:C8	2.51	0.45
1:C:110:LEU:HD11	1:C:120:PHE:CD1	2.52	0.45
1:C:287:PRO:CB	1:C:321:VAL:HA	2.46	0.45
1:C:547:ARG:HG3	1:C:551:LEU:HD12	1.98	0.45
2:G:34:PRO:HA	2:G:37:MET:HE2	1.99	0.45
3:I:819:DG:C8	3:I:820:DT:H72	2.51	0.45
1:B:419:SER:HB3	1:B:433:ALA:O	2.16	0.45
2:F:48:GLU:OE1	2:F:100:LYS:NZ	2.35	0.45
2:G:34:PRO:HB2	2:G:76:PRO:HD3	1.98	0.45
1:A:85:ARG:HG3	1:A:222:TRP:CG	2.52	0.45
1:C:432:HIS:HB2	1:C:436:ASN:OD1	2.16	0.45
1:D:479:LEU:HD22	1:D:654:ASP:HB3	1.99	0.45
1:D:85:ARG:HG3	1:D:222:TRP:CG	2.52	0.45
2:F:32:CYS:SG	2:F:34:PRO:HD2	2.57	0.45
2:H:34:PRO:HB2	2:H:76:PRO:HD3	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:GLU:OE2	1:A:79:ARG:NH2	2.50	0.45
1:C:173:GLN:O	1:C:176:VAL:HG22	2.17	0.45
1:B:482:ARG:HG2	1:B:500:ILE:HD13	1.99	0.44
1:B:200:ASP:OD2	1:C:148:GLU:HB3	2.17	0.44
1:C:573:LEU:HD11	1:C:593:ILE:HD11	1.99	0.44
3:M:814:DT:H2''	3:M:815:DG:C8	2.52	0.44
1:D:355:LYS:HE2	1:D:363:VAL:HG11	1.99	0.44
1:C:49:GLU:OE1	1:C:52:ARG:NH1	2.51	0.44
1:D:597:ASP:OD1	1:D:599:ARG:HD2	2.17	0.44
1:A:110:LEU:HD23	1:A:115:LEU:HB3	1.99	0.44
1:B:391:MET:HE2	1:B:395:ARG:CG	2.48	0.44
1:D:15:VAL:HG22	1:D:72:LEU:HD21	1.98	0.44
1:D:128:TRP:CD1	1:D:181:LEU:HD11	2.53	0.44
2:G:37:MET:HE3	2:G:93:ALA:HB3	1.99	0.44
4:J:855:DT:H2'	4:J:856:DC:C6	2.53	0.44
3:K:809:DG:H2''	3:K:810:DG:C8	2.53	0.44
1:A:9:ASN:ND2	1:A:17:LYS:HB2	2.32	0.44
1:B:99:HIS:HB3	1:B:102:LEU:HD21	1.99	0.44
2:G:37:MET:HE3	2:G:93:ALA:CB	2.48	0.44
2:H:38:ILE:HD12	2:H:93:ALA:HA	1.99	0.44
3:O:802:DG:H2''	3:O:803:DG:C8	2.53	0.44
1:A:632:MET:HE2	1:A:672:THR:HG23	2.00	0.44
1:A:86:GLU:HG3	1:A:203:TYR:CE1	2.52	0.44
1:B:421:ASN:HB3	1:B:431:THR:OG1	2.18	0.44
1:D:432:HIS:O	4:P:858:DC:H4'	2.18	0.44
2:F:103:LEU:O	2:F:107:LEU:HB2	2.17	0.44
3:M:821:DG:H2'	3:M:822:DA:C8	2.53	0.43
1:A:355:LYS:HD3	1:A:368:VAL:CG1	2.48	0.43
1:B:597:ASP:OD1	1:B:599:ARG:HD2	2.18	0.43
2:F:20:ASP:O	2:F:20:ASP:CG	2.60	0.43
1:B:99:HIS:C	1:B:101:ASN:N	2.77	0.43
1:B:549:LYS:HE2	1:B:549:LYS:HB3	1.90	0.43
1:D:486:HIS:HB2	1:D:686:PHE:CE2	2.53	0.43
7:D:805:TTP:N3	4:P:854:DA:N1	2.53	0.43
1:A:289:ILE:HD12	1:A:289:ILE:HA	1.89	0.43
1:B:520:ASN:HA	1:B:523:THR:HG22	2.01	0.43
1:D:233:PRO:HB2	1:D:456:GLY:O	2.18	0.43
1:C:327:THR:HG22	2:G:75:ILE:HG13	1.99	0.43
1:C:394:LYS:HE2	3:M:820:DT:H5'	2.00	0.43
2:F:3:LYS:HE2	2:F:3:LYS:HB3	1.72	0.43
1:A:551:LEU:HA	1:A:551:LEU:HD23	1.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:GLU:O	1:B:130:TYR:HD1	2.01	0.43
1:B:638:LEU:HD22	1:B:645:ASP:CB	2.48	0.43
1:D:536:LYS:O	1:D:540:ILE:HG23	2.18	0.43
1:D:633:LEU:HD22	1:D:638:LEU:HD12	2.00	0.43
1:A:597:ASP:OD2	1:A:599:ARG:NH1	2.49	0.43
1:B:119:ARG:NH2	1:B:126:GLU:OE1	2.51	0.43
2:F:38:ILE:HD12	2:F:93:ALA:HA	2.00	0.43
1:C:570:GLN:HE22	1:C:606:PRO:HB3	1.84	0.43
3:M:813:DT:H2''	3:M:814:DT:H5''	2.01	0.43
1:A:593:ILE:HG23	5:U:4:PHE:HB2	2.01	0.43
1:B:343:GLN:HG3	1:B:362:PRO:CG	2.49	0.43
1:C:285:LYS:O	1:C:286:TYR:HB3	2.19	0.43
1:D:25:ASP:HB3	1:D:28:THR:OG1	2.19	0.43
3:O:814:DT:H2''	3:O:815:DG:C8	2.53	0.43
1:B:553:LYS:O	1:B:557:GLU:HG3	2.18	0.43
1:C:99:HIS:O	1:C:102:LEU:HG	2.19	0.43
1:D:118:LYS:HG3	1:D:119:ARG:HG3	2.00	0.42
2:E:90:LYS:HD2	2:E:90:LYS:HA	1.88	0.42
1:B:404:LYS:HA	1:B:409:TYR:CE2	2.53	0.42
1:A:443:VAL:HG13	1:A:449:GLU:HG3	2.01	0.42
1:B:579:TRP:CE2	1:B:584:GLN:HG3	2.55	0.42
1:C:534:ASP:CG	1:C:549:LYS:HG2	2.44	0.42
1:D:31:TYR:CZ	1:D:176:VAL:HG12	2.54	0.42
1:D:597:ASP:OD2	1:D:599:ARG:NH1	2.49	0.42
1:A:106:ASP:OD2	1:A:131:ARG:NH2	2.53	0.42
1:B:85:ARG:HG3	1:B:222:TRP:CG	2.55	0.42
1:B:476:ALA:HB3	1:B:654:ASP:HB2	2.01	0.42
4:P:855:DT:H2'	4:P:856:DC:C6	2.54	0.42
1:C:140:LYS:HE2	1:C:144:LYS:HE3	2.00	0.42
1:D:193:PRO:HA	1:D:194:PRO:HD3	1.92	0.42
1:A:289:ILE:HD11	2:E:34:PRO:HG2	2.01	0.42
1:A:483:CYS:HA	1:A:686:PHE:CZ	2.49	0.42
1:A:525:ILE:HD12	1:A:525:ILE:HA	1.94	0.42
1:D:638:LEU:HD22	1:D:645:ASP:CB	2.50	0.42
3:M:810:DG:H2'	3:M:811:DT:C6	2.55	0.42
1:C:13:GLU:CD	1:C:13:GLU:H	2.28	0.42
1:C:282:PRO:HD2	1:C:283:LEU:H	1.84	0.42
1:A:193:PRO:HA	1:A:194:PRO:HD3	1.91	0.42
1:A:638:LEU:HD22	1:A:645:ASP:CB	2.50	0.42
1:C:548:GLY:O	1:C:552:LYS:HG3	2.20	0.42
1:C:664:GLU:O	1:C:668:VAL:HG23	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:9:ASN:ND2	1:D:17:LYS:HB2	2.34	0.42
1:D:391:MET:HE1	1:D:447:TYR:HE2	1.85	0.42
3:O:810:DG:H2'	3:O:811:DT:C6	2.55	0.42
1:A:130:TYR:CD1	1:A:135:MET:HG3	2.55	0.42
1:A:664:GLU:O	1:A:668:VAL:HG23	2.19	0.42
1:B:296:GLY:O	1:B:316:ASP:HB3	2.20	0.42
1:C:287:PRO:HB3	1:C:322:ALA:N	2.33	0.42
2:H:37:MET:HE3	2:H:93:ALA:CB	2.50	0.42
1:B:431:THR:HG22	4:L:857:DA:H4'	2.01	0.42
3:K:819:DG:C2'	3:K:820:DT:H71	2.50	0.42
1:C:391:MET:HE1	1:C:447:TYR:HE2	1.84	0.41
1:D:158:MET:HA	1:D:161:TRP:CE2	2.54	0.41
1:D:520:ASN:HA	1:D:523:THR:HG22	2.01	0.41
1:D:640:HIS:CD2	1:D:647:ALA:HA	2.55	0.41
1:B:59:HIS:CG	1:B:125:LEU:HG	2.55	0.41
1:C:109:LEU:HD23	1:C:109:LEU:HA	1.87	0.41
1:D:109:LEU:HB3	1:D:115:LEU:HD23	2.02	0.41
3:K:802:DG:H2''	3:K:803:DG:H8	1.85	0.41
1:B:47:GLU:HG3	1:B:81:PHE:HE1	1.86	0.41
1:C:17:LYS:NZ	1:D:449:GLU:OE2	2.52	0.41
1:C:119:ARG:NH2	1:C:126:GLU:OE1	2.52	0.41
1:C:329:VAL:HG11	2:G:31:TRP:CH2	2.54	0.41
2:E:8:THR:CG2	2:E:11:SER:H	2.33	0.41
2:F:27:PHE:CE1	2:F:79:LEU:HD22	2.56	0.41
1:D:139:TYR:HB2	1:D:170:TYR:CG	2.55	0.41
2:H:24:LEU:HD11	2:H:78:LEU:HB3	2.02	0.41
1:A:96:ARG:HA	1:A:123:HIS:CE1	2.56	0.41
1:A:624:LEU:HD12	1:A:624:LEU:HA	1.92	0.41
1:A:674:GLN:O	1:A:678:ARG:HG2	2.21	0.41
2:E:107:LEU:HD23	2:E:107:LEU:HA	1.85	0.41
1:A:224:LEU:O	1:A:228:GLU:HG3	2.20	0.41
1:C:526:TYR:CD2	7:C:805:TTP:H2'2	2.55	0.41
1:D:534:ASP:OD1	1:D:534:ASP:N	2.54	0.41
2:E:9:ASP:OD2	2:E:66:THR:OG1	2.32	0.41
1:B:457:ALA:CA	1:B:649:MET:HE1	2.50	0.41
1:D:160:TRP:CE3	1:D:167:MET:HE2	2.55	0.41
1:D:456:GLY:HA2	1:D:471:GLN:OE1	2.21	0.41
4:N:870:DC:H2''	4:N:871:DC:C6	2.55	0.41
1:A:520:ASN:HA	1:A:523:THR:HG22	2.02	0.41
1:B:664:GLU:O	1:B:668:VAL:HG23	2.20	0.41
1:D:224:LEU:O	1:D:228:GLU:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:99:LEU:HD23	2:G:99:LEU:HA	1.80	0.41
1:A:61:GLY:HA2	1:A:65:ASP:HB2	2.03	0.41
1:A:628:LYS:HE3	1:A:679:TRP:CE2	2.56	0.41
1:B:61:GLY:HA2	1:B:65:ASP:HB2	2.03	0.41
1:B:633:LEU:HD22	1:B:638:LEU:HD12	2.03	0.41
1:C:104:ASP:HA	1:C:107:MET:HG2	2.03	0.41
1:C:140:LYS:HE3	1:C:154:TYR:OH	2.21	0.41
1:D:59:HIS:CD2	1:D:125:LEU:HG	2.55	0.41
4:N:861:DG:H2''	4:N:862:DC:C6	2.56	0.41
1:A:440:ILE:O	1:A:452:ARG:NH2	2.52	0.41
1:B:193:PRO:HA	1:B:194:PRO:HD3	1.92	0.41
1:B:336:PRO:HG2	1:B:389:TYR:CZ	2.56	0.41
1:C:597:ASP:OD1	1:C:599:ARG:HD2	2.21	0.41
2:F:36:LYS:HE2	2:F:36:LYS:HB2	1.71	0.41
1:A:35:ARG:HB3	1:A:36:PRO:HD2	2.04	0.40
1:A:534:ASP:CG	1:A:549:LYS:HG2	2.46	0.40
1:B:440:ILE:HD12	1:B:700:TRP:CZ3	2.56	0.40
1:C:686:PHE:HD2	1:C:688:CYS:O	2.03	0.40
1:D:664:GLU:O	1:D:668:VAL:HG23	2.20	0.40
1:B:588:TRP:CE2	1:B:591:ARG:HD2	2.56	0.40
1:C:17:LYS:HZ2	1:D:449:GLU:CD	2.29	0.40
1:D:670:ILE:HG23	1:D:694:GLY:HA3	2.04	0.40
2:H:32:CYS:SG	2:H:34:PRO:HD2	2.60	0.40
3:M:802:DG:H2''	3:M:803:DG:H8	1.84	0.40
1:A:394:LYS:HE3	3:I:819:DG:H1'	2.03	0.40
1:B:31:TYR:OH	1:B:176:VAL:HG22	2.20	0.40
1:C:173:GLN:HA	1:C:176:VAL:HG22	2.02	0.40
3:M:806:DG:H2''	3:M:807:DG:O5'	2.22	0.40
1:D:346:LEU:HD21	1:D:383:ILE:HG12	2.03	0.40
1:D:549:LYS:HE2	1:D:549:LYS:HB3	1.97	0.40
1:D:576:SER:OG	1:D:587:LYS:HD2	2.21	0.40
1:A:553:LYS:O	1:A:557:GLU:HG3	2.20	0.40
1:D:66:VAL:HG11	1:D:222:TRP:CH2	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	690/704 (98%)	666 (96%)	20 (3%)	4 (1%)	21	56
1	B	693/704 (98%)	666 (96%)	22 (3%)	5 (1%)	18	53
1	C	666/704 (95%)	631 (95%)	21 (3%)	14 (2%)	5	27
1	D	638/704 (91%)	619 (97%)	15 (2%)	4 (1%)	21	56
2	E	103/109 (94%)	99 (96%)	2 (2%)	2 (2%)	6	30
2	F	103/109 (94%)	100 (97%)	3 (3%)	0	100	100
2	G	103/109 (94%)	94 (91%)	8 (8%)	1 (1%)	12	45
2	H	103/109 (94%)	99 (96%)	4 (4%)	0	100	100
5	U	2/4 (50%)	2 (100%)	0	0	100	100
5	V	2/4 (50%)	1 (50%)	1 (50%)	0	100	100
6	W	1/3 (33%)	1 (100%)	0	0	100	100
All	All	3104/3263 (95%)	2978 (96%)	96 (3%)	30 (1%)	12	45

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	135	MET
1	C	276	HIS
1	C	282	PRO
1	C	287	PRO
1	C	359	LYS
1	A	117	GLY
1	C	156	ASP
1	C	270	GLY
1	C	274	PHE
1	C	277	PRO
1	C	318	ARG
1	D	359	LYS
2	E	19	ALA

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Mol	Chain	Res	Type
1	C	479	LEU
1	D	479	LEU
1	A	479	LEU
1	B	100	SER
1	B	300	LYS
1	B	479	LEU
2	E	20	ASP
1	A	314	GLU
1	A	653	HIS
1	B	653	HIS
1	C	279	THR
1	C	283	LEU
1	C	321	VAL
1	C	653	HIS
1	D	360	GLY
1	D	653	HIS
2	G	105	ALA

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	574/582 (99%)	574 (100%)	0	100	100
1	B	576/582 (99%)	576 (100%)	0	100	100
1	C	556/582 (96%)	556 (100%)	0	100	100
1	D	531/582 (91%)	531 (100%)	0	100	100
2	E	85/88 (97%)	85 (100%)	0	100	100
2	F	85/88 (97%)	85 (100%)	0	100	100
2	G	85/88 (97%)	85 (100%)	0	100	100
2	H	85/88 (97%)	85 (100%)	0	100	100
5	U	4/4 (100%)	4 (100%)	0	100	100
5	V	4/4 (100%)	4 (100%)	0	100	100
6	W	3/3 (100%)	3 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	2588/2691 (96%)	2588 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	101	ASN
1	A	171	ASN
1	A	173	GLN
1	A	227	GLN
1	A	331	HIS
1	A	347	GLN
1	A	432	HIS
1	A	493	ASN
1	A	602	HIS
1	B	150	GLN
1	B	227	GLN
1	B	432	HIS
1	B	704	HIS
1	C	173	GLN
1	C	227	GLN
1	C	704	HIS
1	D	101	ASN
1	D	450	GLN
1	D	570	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](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 20 ligands modelled in this entry, 12 are monoatomic - leaving 8 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
7	TTP	D	805	8	29,30,30	1.41	5 (17%)	43,47,47	1.85	7 (16%)
7	TTP	A	801	-	29,30,30	1.42	8 (27%)	43,47,47	1.71	6 (13%)
7	TTP	B	805	8	29,30,30	1.42	7 (24%)	43,47,47	1.89	7 (16%)
7	TTP	B	801	-	29,30,30	1.45	9 (31%)	43,47,47	2.00	10 (23%)
7	TTP	D	801	-	29,30,30	1.52	9 (31%)	43,47,47	1.84	9 (20%)
7	TTP	C	801	-	29,30,30	1.42	8 (27%)	43,47,47	1.84	7 (16%)
7	TTP	A	805	8	29,30,30	1.35	8 (27%)	43,47,47	1.83	7 (16%)
7	TTP	C	805	8	29,30,30	1.37	7 (24%)	43,47,47	1.77	9 (20%)

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
7	TTP	D	805	8	-	2/22/34/34	0/2/2/2
7	TTP	A	801	-	-	3/22/34/34	0/2/2/2
7	TTP	B	805	8	-	3/22/34/34	0/2/2/2
7	TTP	B	801	-	-	7/22/34/34	0/2/2/2
7	TTP	D	801	-	-	4/22/34/34	0/2/2/2
7	TTP	C	801	-	-	6/22/34/34	0/2/2/2
7	TTP	A	805	8	-	2/22/34/34	0/2/2/2
7	TTP	C	805	8	-	2/22/34/34	0/2/2/2

All (61) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	D	801	TTP	PA-O3A	3.23	1.63	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	D	805	TTP	C6-C5	3.17	1.39	1.34
7	D	801	TTP	PB-O3A	3.16	1.62	1.59
7	B	805	TTP	PA-O3A	3.16	1.62	1.59
7	B	801	TTP	C6-C5	3.03	1.39	1.34
7	C	805	TTP	C6-C5	3.00	1.39	1.34
7	A	801	TTP	PB-O3A	2.84	1.62	1.59
7	B	805	TTP	C6-C5	2.83	1.39	1.34
7	B	805	TTP	PB-O3A	2.81	1.62	1.59
7	C	801	TTP	C6-C5	2.77	1.39	1.34
7	A	801	TTP	C6-C5	2.75	1.39	1.34
7	B	801	TTP	PB-O3A	2.69	1.62	1.59
7	D	801	TTP	PB-O3B	2.68	1.62	1.59
7	D	805	TTP	C4-C5	2.68	1.49	1.44
7	A	801	TTP	PA-O3A	2.63	1.62	1.59
7	B	801	TTP	C4-N3	-2.63	1.33	1.38
7	D	805	TTP	PB-O3A	2.63	1.62	1.59
7	D	801	TTP	C6-C5	2.63	1.38	1.34
7	C	801	TTP	PB-O3B	2.61	1.62	1.59
7	C	805	TTP	C4-C5	2.60	1.49	1.44
7	B	801	TTP	PB-O3B	2.55	1.62	1.59
7	C	801	TTP	PA-O3A	2.54	1.62	1.59
7	C	801	TTP	C4-N3	-2.50	1.34	1.38
7	A	805	TTP	C4-N3	-2.50	1.34	1.38
7	C	805	TTP	C4-N3	-2.50	1.34	1.38
7	A	805	TTP	C6-C5	2.49	1.38	1.34
7	D	801	TTP	C4-N3	-2.49	1.34	1.38
7	A	805	TTP	PA-O3A	2.49	1.62	1.59
7	D	805	TTP	C4-N3	-2.47	1.34	1.38
7	A	801	TTP	C4-N3	-2.45	1.34	1.38
7	B	805	TTP	C4-C5	2.41	1.48	1.44
7	C	801	TTP	C4-C5	2.40	1.48	1.44
7	A	801	TTP	C2-N1	2.37	1.42	1.38
7	D	801	TTP	C4-C5	2.34	1.48	1.44
7	B	801	TTP	C4-C5	2.31	1.48	1.44
7	B	801	TTP	PA-O3A	2.25	1.61	1.59
7	D	801	TTP	C6-N1	-2.23	1.34	1.38
7	B	801	TTP	C2-N3	-2.22	1.34	1.38
7	A	805	TTP	PB-O3A	2.22	1.61	1.59
7	A	805	TTP	C2-N1	2.21	1.41	1.38
7	A	805	TTP	C6-N1	-2.21	1.34	1.38
7	B	805	TTP	C2-N1	2.19	1.41	1.38
7	C	801	TTP	C6-N1	-2.18	1.34	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	C	801	TTP	PB-O3A	2.16	1.61	1.59
7	C	805	TTP	C2-N1	2.16	1.41	1.38
7	D	805	TTP	PA-O3A	2.15	1.61	1.59
7	B	805	TTP	C4-N3	-2.15	1.34	1.38
7	A	801	TTP	C6-N1	-2.14	1.34	1.38
7	C	801	TTP	C2-N3	-2.12	1.34	1.38
7	B	801	TTP	C6-N1	-2.12	1.34	1.38
7	A	805	TTP	C2-N3	-2.09	1.34	1.38
7	D	801	TTP	C2-N3	-2.09	1.34	1.38
7	A	805	TTP	C4-C5	2.07	1.48	1.44
7	B	801	TTP	C2-N1	2.07	1.41	1.38
7	A	801	TTP	C4-C5	2.07	1.48	1.44
7	C	805	TTP	C6-N1	-2.05	1.34	1.38
7	C	805	TTP	PB-O3B	2.03	1.61	1.59
7	D	801	TTP	C2-N1	2.02	1.41	1.38
7	B	805	TTP	C6-N1	-2.01	1.34	1.38
7	C	805	TTP	PB-O3A	2.01	1.61	1.59
7	A	801	TTP	PB-O3B	2.00	1.61	1.59

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	801	TTP	N3-C2-N1	5.48	122.03	114.89
7	C	805	TTP	N3-C2-N1	5.23	121.70	114.89
7	C	801	TTP	C4-N3-C2	-5.15	120.58	127.34
7	D	805	TTP	C4-N3-C2	-5.11	120.64	127.34
7	B	801	TTP	C4-N3-C2	-5.09	120.67	127.34
7	B	805	TTP	C4-N3-C2	-5.02	120.75	127.34
7	D	805	TTP	N3-C2-N1	4.96	121.34	114.89
7	D	801	TTP	C4-N3-C2	-4.93	120.87	127.34
7	A	805	TTP	C4-N3-C2	-4.91	120.91	127.34
7	C	801	TTP	N3-C2-N1	4.89	121.26	114.89
7	C	805	TTP	C4-N3-C2	-4.82	121.03	127.34
7	A	801	TTP	N3-C2-N1	4.77	121.11	114.89
7	D	801	TTP	N3-C2-N1	4.77	121.10	114.89
7	B	805	TTP	N3-C2-N1	4.76	121.08	114.89
7	A	805	TTP	N3-C2-N1	4.71	121.02	114.89
7	D	801	TTP	C5-C4-N3	4.64	119.35	115.32
7	C	801	TTP	C5-C4-N3	4.58	119.30	115.32
7	B	805	TTP	O4-C4-C5	-4.54	119.72	124.92
7	D	805	TTP	C5-C4-N3	4.49	119.22	115.32
7	B	805	TTP	C5-C4-N3	4.41	119.16	115.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	801	TTP	C4-N3-C2	-4.40	121.57	127.34
7	A	805	TTP	O4-C4-C5	-4.27	120.03	124.92
7	A	805	TTP	C5-C4-N3	4.24	119.01	115.32
7	A	801	TTP	O4-C4-C5	-4.14	120.18	124.92
7	B	801	TTP	C5-C4-N3	4.12	118.91	115.32
7	D	801	TTP	O4-C4-C5	-3.99	120.35	124.92
7	D	805	TTP	C5-C6-N1	-3.97	119.00	123.31
7	B	801	TTP	O4-C4-C5	-3.97	120.38	124.92
7	C	801	TTP	O4-C4-C5	-3.85	120.51	124.92
7	C	805	TTP	C5-C4-N3	3.84	118.66	115.32
7	B	801	TTP	O4'-C1'-N1	3.83	114.66	107.86
7	D	805	TTP	O4-C4-C5	-3.83	120.54	124.92
7	A	801	TTP	C5-C4-N3	3.76	118.59	115.32
7	B	805	TTP	C2'-C1'-N1	-3.73	104.50	113.81
7	C	801	TTP	C5-C6-N1	-3.65	119.35	123.31
7	C	805	TTP	O4-C4-C5	-3.48	120.94	124.92
7	A	805	TTP	C5-C6-N1	-3.41	119.61	123.31
7	D	805	TTP	O2-C2-N1	-3.29	118.51	122.80
7	B	801	TTP	C5-C6-N1	-3.26	119.77	123.31
7	A	805	TTP	C2'-C1'-N1	-3.25	105.69	113.81
7	C	801	TTP	O2-C2-N1	-3.08	118.78	122.80
7	B	805	TTP	O2-C2-N1	-3.08	118.79	122.80
7	C	805	TTP	C5-C6-N1	-3.04	120.01	123.31
7	B	805	TTP	C5-C6-N1	-3.04	120.02	123.31
7	B	801	TTP	O2-C2-N1	-2.76	119.20	122.80
7	C	805	TTP	O2-C2-N1	-2.75	119.22	122.80
7	D	801	TTP	C5-C6-N1	-2.74	120.34	123.31
7	D	805	TTP	C2'-C1'-N1	-2.73	106.98	113.81
7	D	801	TTP	C5M-C5-C4	2.73	121.69	118.78
7	D	801	TTP	C2'-C1'-N1	-2.60	107.30	113.81
7	A	805	TTP	O2-C2-N1	-2.49	119.56	122.80
7	B	801	TTP	C2'-C1'-N1	-2.48	107.62	113.81
7	D	801	TTP	O2-C2-N1	-2.28	119.82	122.80
7	A	801	TTP	C5-C6-N1	-2.28	120.84	123.31
7	C	805	TTP	O4'-C1'-N1	2.27	111.89	107.86
7	B	801	TTP	C3'-C2'-C1'	2.24	108.08	102.60
7	C	805	TTP	C2'-C1'-N1	-2.15	108.44	113.81
7	A	801	TTP	C1'-N1-C2	2.14	121.84	117.66
7	D	801	TTP	O2B-PB-O3A	2.14	113.05	107.27
7	B	801	TTP	C4'-O4'-C1'	2.08	114.46	109.51
7	C	801	TTP	O2A-PA-O3A	2.03	112.75	107.27
7	C	805	TTP	O2B-PB-O3A	2.00	112.69	107.27

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	B	801	TTP	C5'-O5'-PA-O1A
7	B	801	TTP	C5'-O5'-PA-O2A
7	B	801	TTP	C5'-O5'-PA-O3A
7	C	801	TTP	O4'-C4'-C5'-O5'
7	D	801	TTP	O4'-C4'-C5'-O5'
7	C	801	TTP	C3'-C4'-C5'-O5'
7	D	801	TTP	C3'-C4'-C5'-O5'
7	B	801	TTP	O4'-C4'-C5'-O5'
7	B	801	TTP	C3'-C4'-C5'-O5'
7	B	801	TTP	PA-O3A-PB-O1B
7	A	801	TTP	C5'-O5'-PA-O2A
7	A	801	TTP	C5'-O5'-PA-O3A
7	A	805	TTP	PA-O3A-PB-O2B
7	B	805	TTP	PA-O3A-PB-O2B
7	C	801	TTP	PG-O3B-PB-O2B
7	D	801	TTP	PA-O3A-PB-O2B
7	B	805	TTP	PB-O3B-PG-O2G
7	A	801	TTP	PB-O3A-PA-O2A
7	A	805	TTP	PA-O3A-PB-O1B
7	B	801	TTP	PA-O3A-PB-O2B
7	B	805	TTP	PA-O3A-PB-O1B
7	C	801	TTP	PB-O3A-PA-O1A
7	C	801	TTP	PB-O3A-PA-O2A
7	C	805	TTP	PA-O3A-PB-O1B
7	C	805	TTP	PA-O3A-PB-O2B
7	D	801	TTP	PA-O3A-PB-O1B
7	D	805	TTP	PA-O3A-PB-O1B
7	D	805	TTP	PA-O3A-PB-O2B
7	C	801	TTP	PG-O3B-PB-O1B

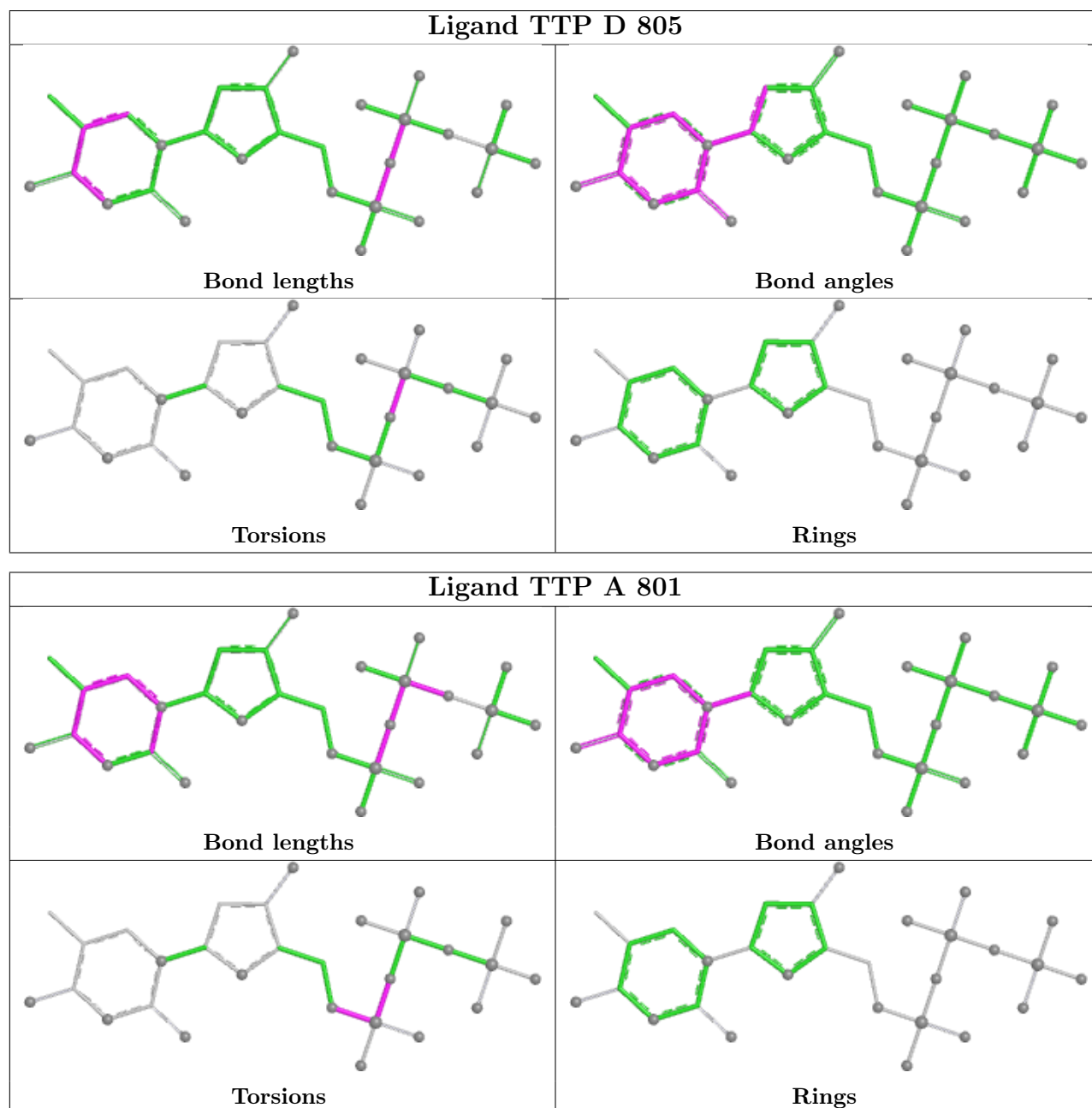
There are no ring outliers.

4 monomers are involved in 5 short contacts:

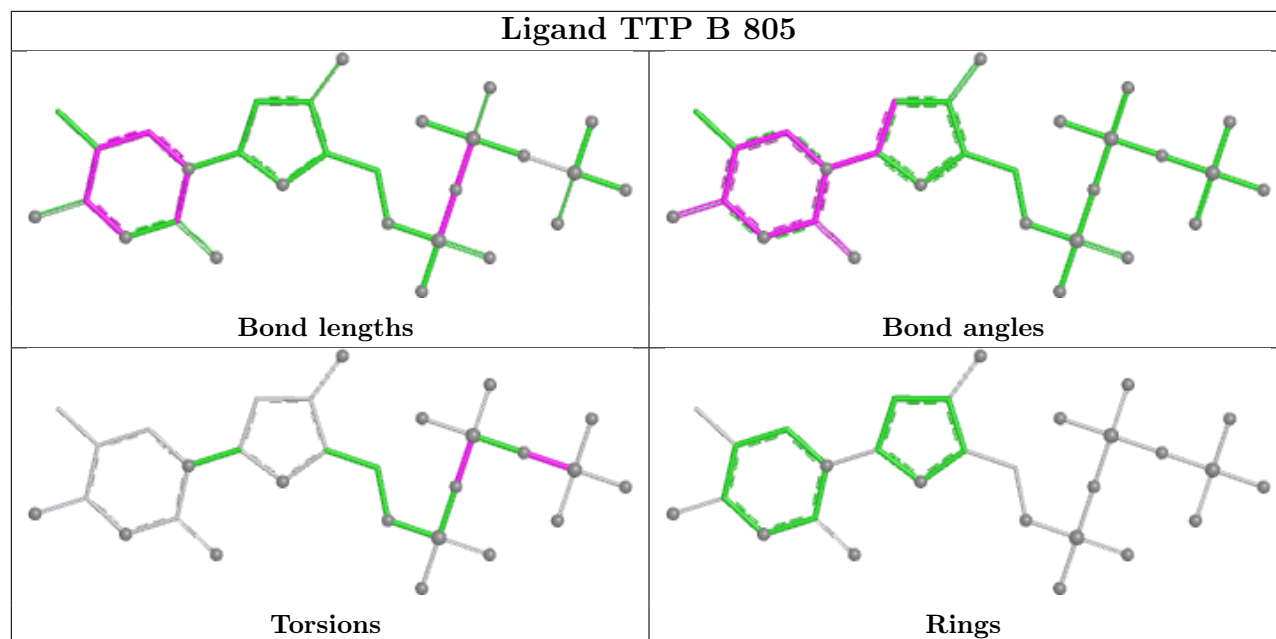
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	D	805	TTP	2	0
7	B	801	TTP	1	0
7	D	801	TTP	1	0
7	C	805	TTP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

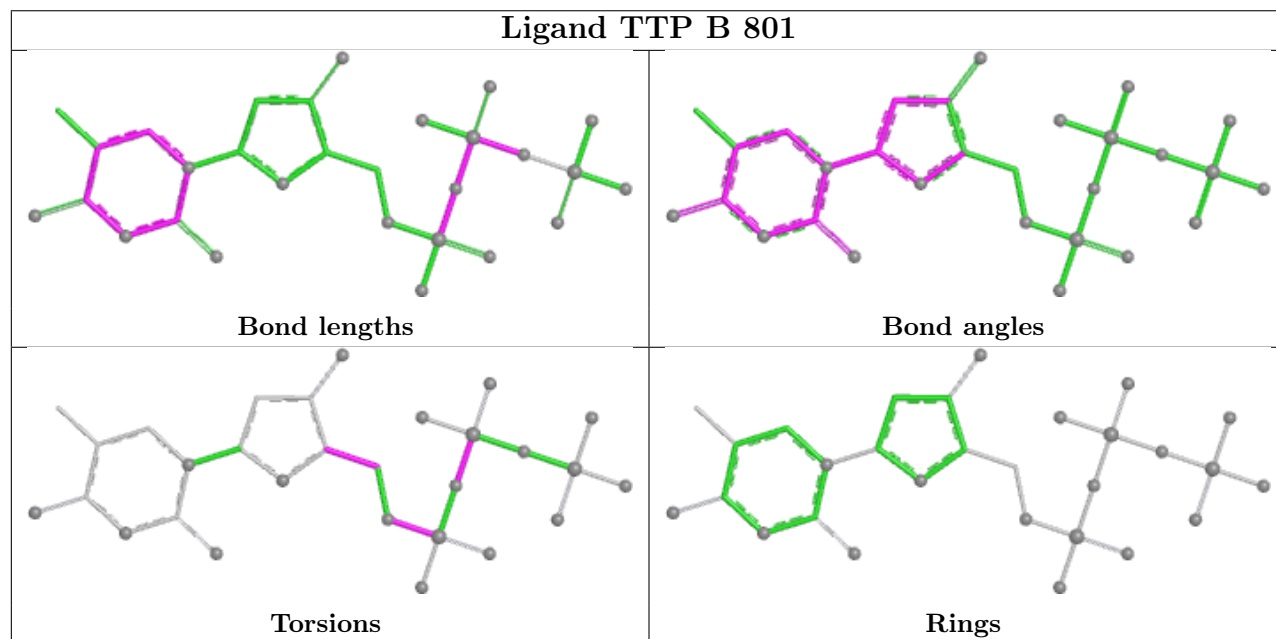
bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 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.



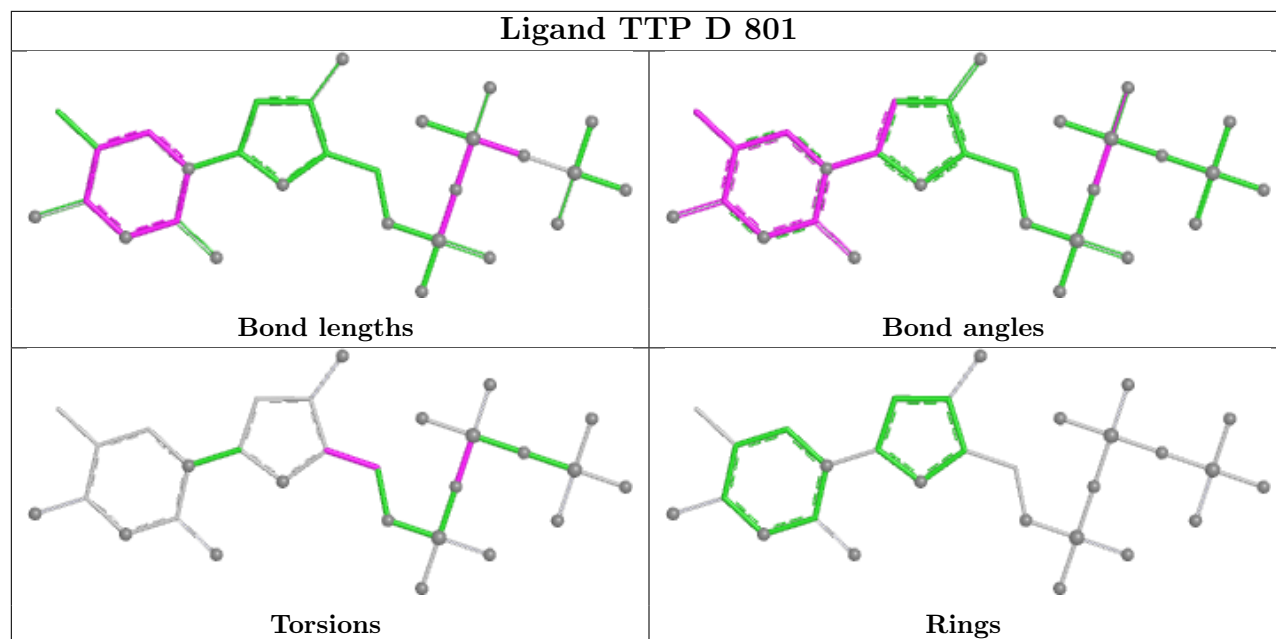
Ligand TTP B 805



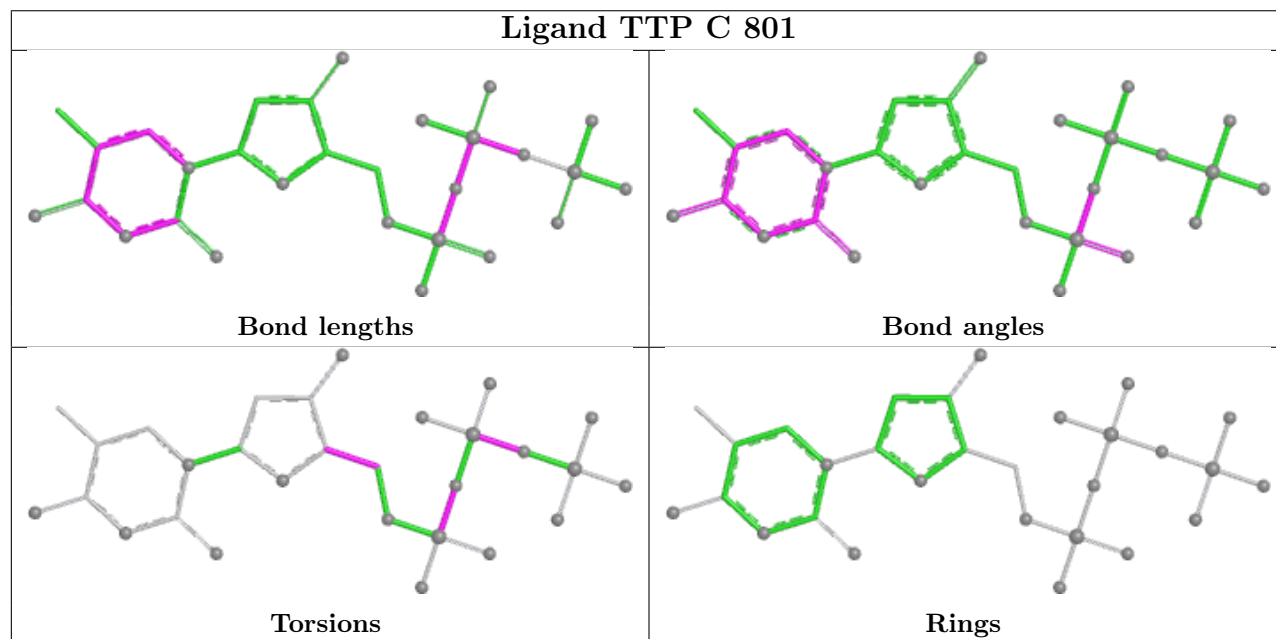
Ligand TTP B 801

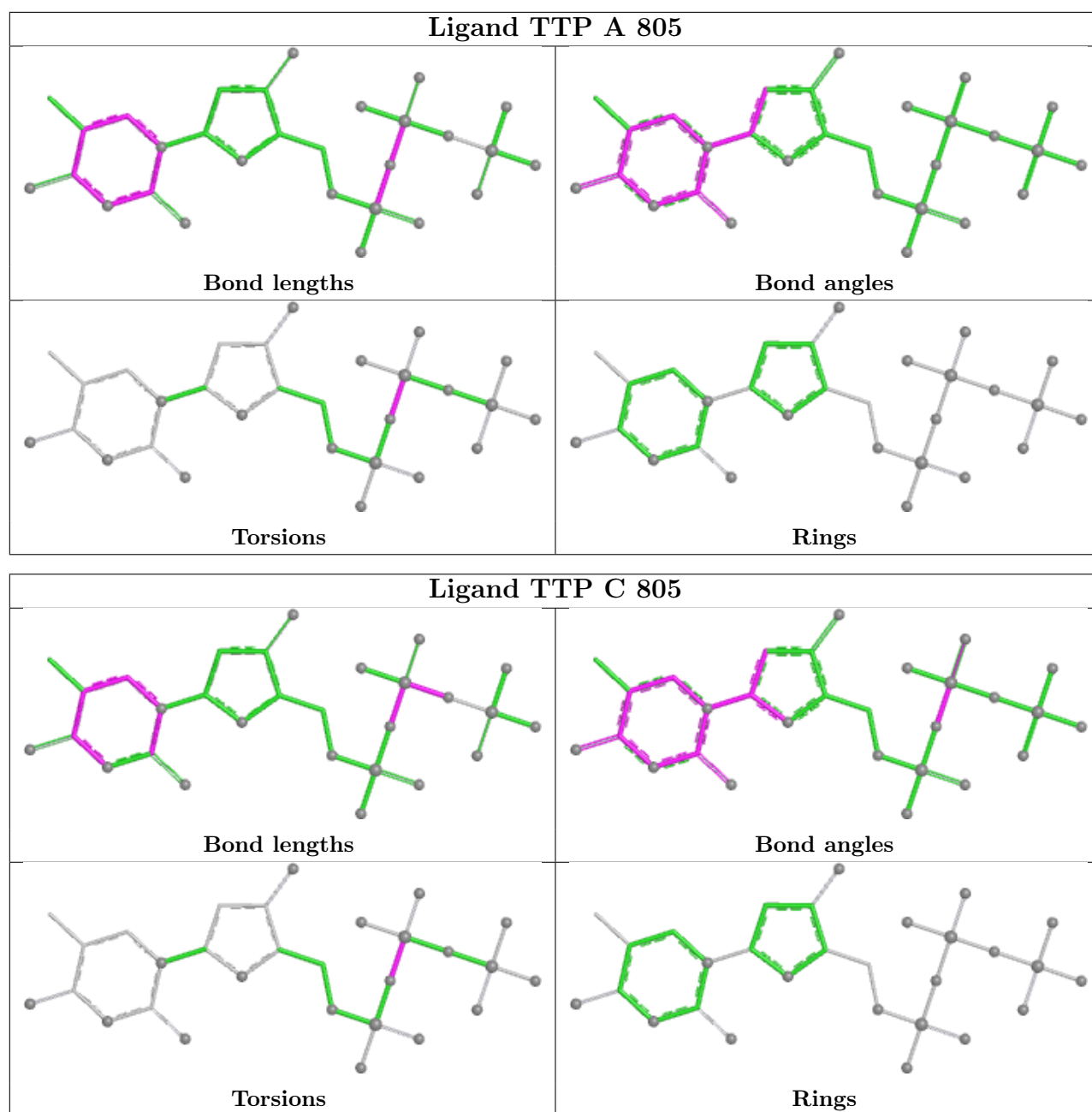


Ligand TTP D 801



Ligand TTP C 801





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	694/704 (98%)	-0.15	9 (1%) 75 53	18, 42, 78, 117	0
1	B	697/704 (99%)	-0.24	3 (0%) 88 76	23, 40, 76, 123	0
1	C	672/704 (95%)	0.23	17 (2%) 58 35	36, 60, 105, 136	0
1	D	644/704 (91%)	0.40	24 (3%) 45 25	36, 74, 110, 135	0
2	E	105/109 (96%)	0.03	2 (1%) 66 43	38, 58, 85, 95	0
2	F	105/109 (96%)	0.22	0 100 100	45, 67, 97, 106	0
2	G	105/109 (96%)	0.85	5 (4%) 35 19	63, 107, 131, 141	0
2	H	105/109 (96%)	0.95	4 (3%) 44 24	81, 122, 141, 151	0
3	I	21/21 (100%)	0.04	0 100 100	26, 51, 68, 81	0
3	K	21/21 (100%)	0.19	0 100 100	26, 60, 74, 88	0
3	M	21/21 (100%)	0.38	0 100 100	38, 70, 82, 84	0
3	O	21/21 (100%)	0.72	1 (4%) 35 19	61, 78, 88, 90	0
4	J	25/25 (100%)	-0.10	0 100 100	22, 51, 66, 86	0
4	L	23/25 (92%)	-0.11	0 100 100	24, 45, 75, 79	0
4	N	24/25 (96%)	0.40	1 (4%) 40 22	38, 73, 94, 110	0
4	P	23/25 (92%)	0.42	0 100 100	56, 75, 94, 100	0
5	U	4/4 (100%)	0.48	0 100 100	54, 69, 95, 109	0
5	V	4/4 (100%)	0.55	0 100 100	54, 73, 86, 95	0
6	W	3/3 (100%)	0.32	0 100 100	74, 74, 81, 84	0
All	All	3317/3447 (96%)	0.12	66 (1%) 65 41	18, 56, 113, 151	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	301	PRO	4.1

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Mol	Chain	Res	Type	RSRZ
1	C	286	TYR	3.9
1	D	322	ALA	3.8
1	C	276	HIS	3.8
1	C	579	TRP	3.5
1	D	556	LEU	3.3
1	D	540	ILE	3.2
1	B	704	HIS	3.1
1	C	704	HIS	3.1
1	D	265	TYR	3.1
1	A	147	LEU	3.0
2	G	25	VAL	3.0
1	D	528	PHE	3.0
1	D	133	GLY	2.9
1	D	551	LEU	2.9
1	C	394	LYS	2.9
1	D	103	LYS	2.8
1	A	704	HIS	2.8
1	D	579	TRP	2.7
2	G	3	LYS	2.7
1	C	281	LYS	2.7
2	G	12	PHE	2.6
2	H	67	ALA	2.6
1	A	587	LYS	2.5
1	D	136	LYS	2.5
1	C	145	ARG	2.4
1	A	575	GLU	2.4
2	G	5	ILE	2.4
1	D	559	THR	2.3
1	C	268	LYS	2.3
1	A	112	SER	2.3
2	H	89	THR	2.3
1	D	173	GLN	2.3
1	D	704	HIS	2.3
2	G	107	LEU	2.3
1	A	173	GLN	2.3
1	B	360	GLY	2.3
1	C	278	ARG	2.2
1	A	360	GLY	2.2
1	C	173	GLN	2.2
1	C	104	ASP	2.2
4	N	867	DC	2.2
1	C	274	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
2	E	19	ALA	2.2
1	C	103	LYS	2.2
1	D	524	PHE	2.2
1	D	686	PHE	2.2
1	D	134	GLU	2.1
1	A	301	PRO	2.1
1	C	282	PRO	2.1
2	H	3	LYS	2.1
1	D	137	GLY	2.1
3	O	802	DG	2.1
2	H	55	VAL	2.1
1	C	339	ARG	2.1
1	D	338	SER	2.1
2	E	3	LYS	2.1
1	C	269	GLY	2.1
1	D	366	ASP	2.1
1	D	426	VAL	2.0
1	D	119	ARG	2.0
1	D	189	LYS	2.0
1	D	330	GLU	2.0
1	A	110	LEU	2.0
1	D	497	ALA	2.0
1	C	110	LEU	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.

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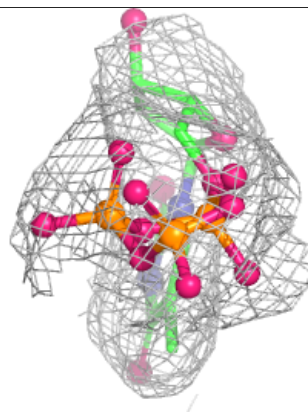
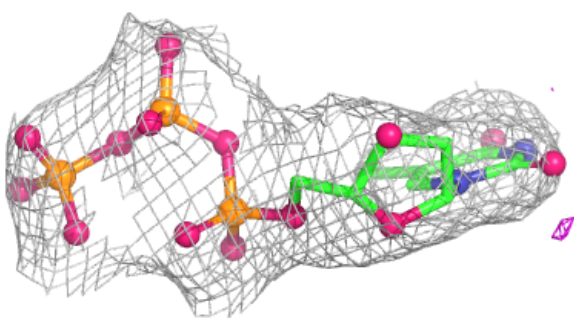
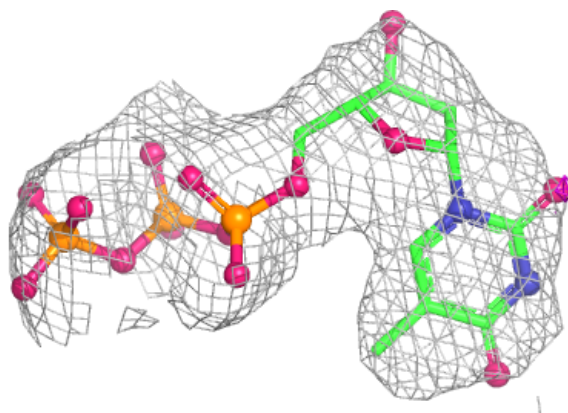
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	MG	C	803	1/1	0.62	0.20	79,79,79,79	0
8	MG	A	803	1/1	0.68	0.21	56,56,56,56	0
7	TTP	C	801	29/29	0.82	0.11	53,93,147,157	0
8	MG	C	804	1/1	0.82	0.36	61,61,61,61	0
7	TTP	D	801	29/29	0.85	0.09	45,73,125,136	0
8	MG	B	803	1/1	0.87	0.23	64,64,64,64	0
7	TTP	A	801	29/29	0.89	0.10	44,60,137,146	0
8	MG	B	804	1/1	0.89	0.22	38,38,38,38	0
8	MG	D	803	1/1	0.89	0.11	90,90,90,90	0
7	TTP	B	801	29/29	0.90	0.10	38,69,137,142	0
7	TTP	D	805	29/29	0.92	0.08	47,67,92,96	0
8	MG	D	804	1/1	0.92	0.12	68,68,68,68	0
7	TTP	C	805	29/29	0.96	0.08	40,50,59,62	0
7	TTP	B	805	29/29	0.97	0.06	20,28,39,45	0
8	MG	A	804	1/1	0.97	0.16	33,33,33,33	0
7	TTP	A	805	29/29	0.98	0.06	22,32,45,54	0
8	MG	D	802	1/1	0.98	0.03	85,85,85,85	0
8	MG	A	802	1/1	0.99	0.06	38,38,38,38	0
8	MG	C	802	1/1	1.00	0.08	43,43,43,43	0
8	MG	B	802	1/1	1.00	0.03	36,36,36,36	0

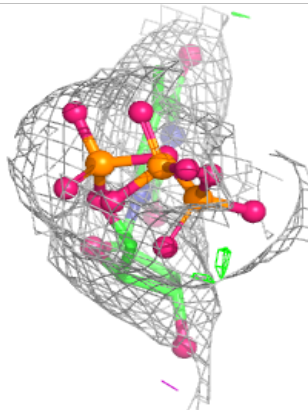
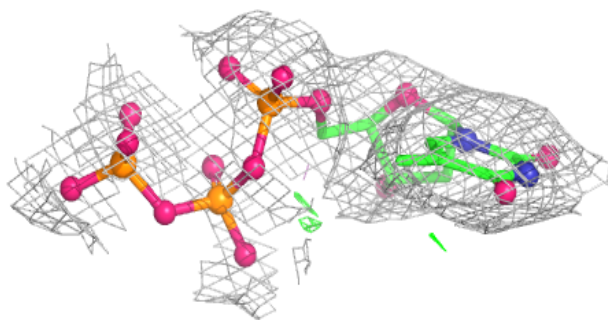
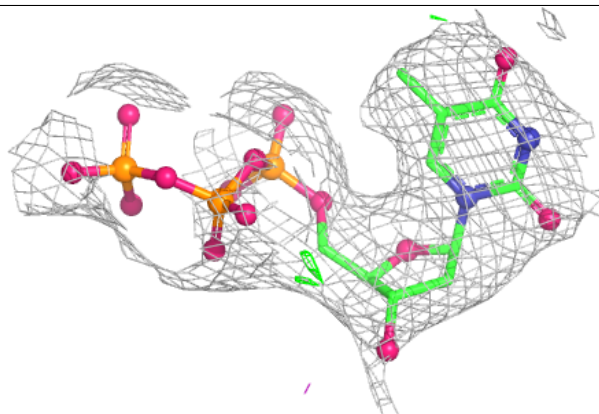
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around TTP C 801:

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

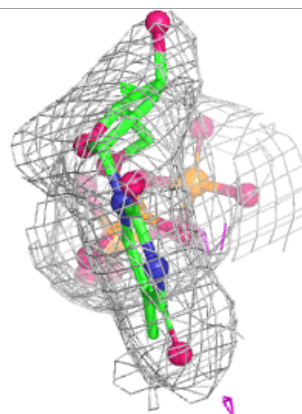
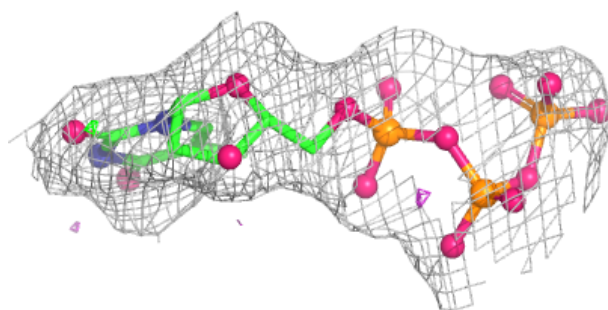
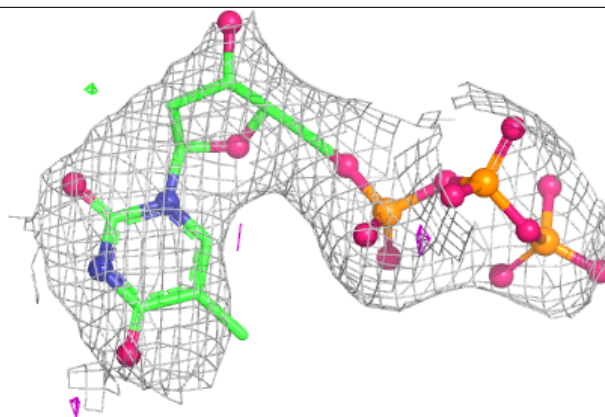
**Electron density around TTP D 801:**

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



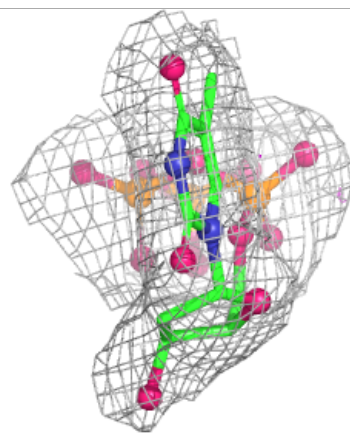
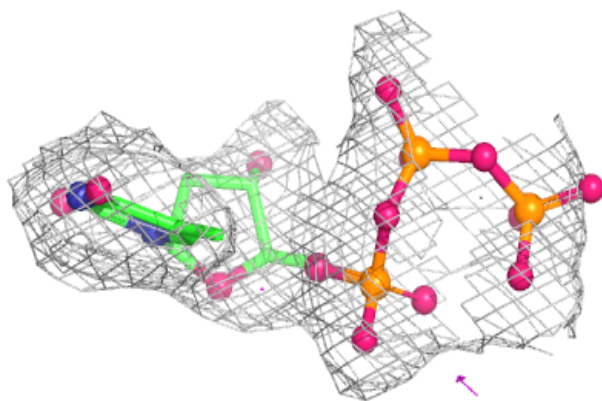
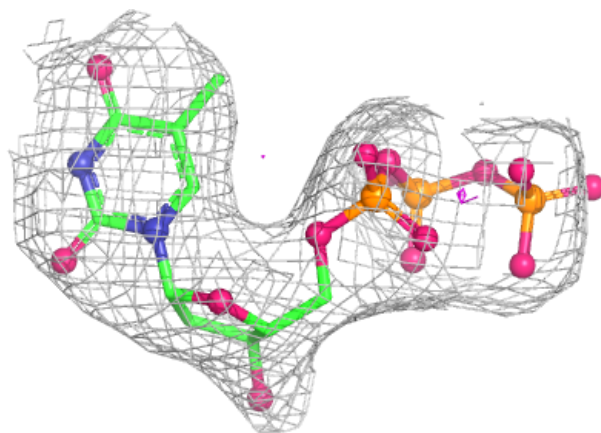
Electron density around TTP A 801:

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



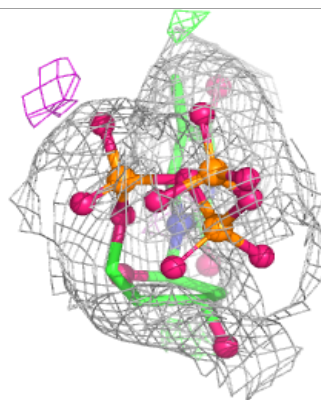
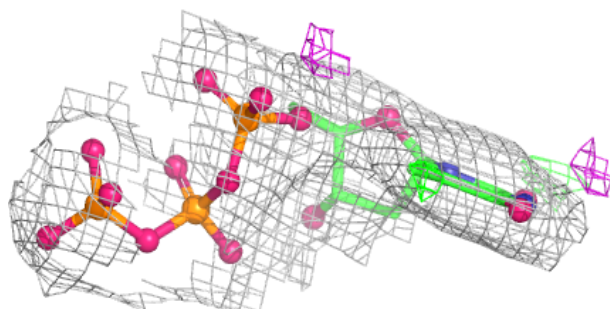
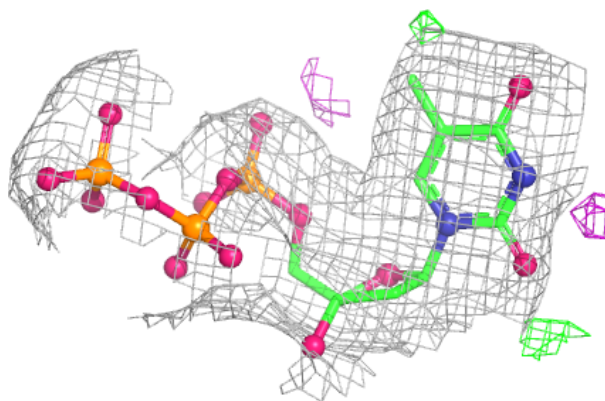
Electron density around TTP B 801:

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

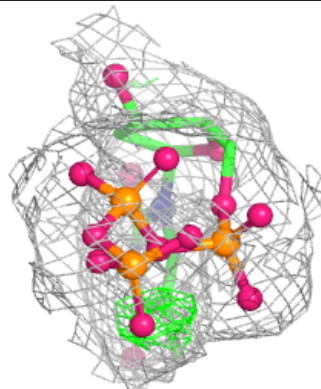
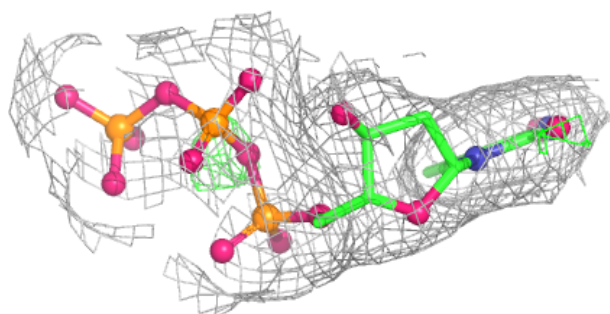
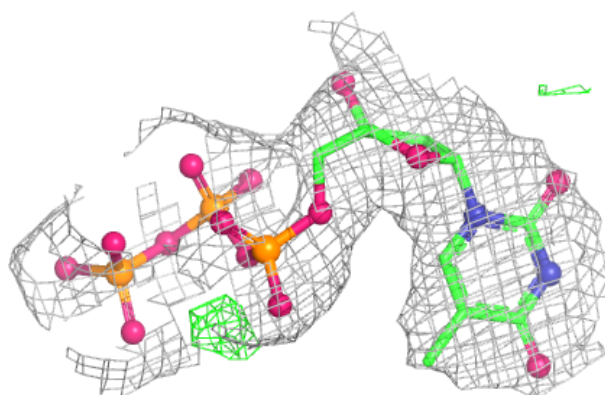


Electron density around TTP D 805:

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

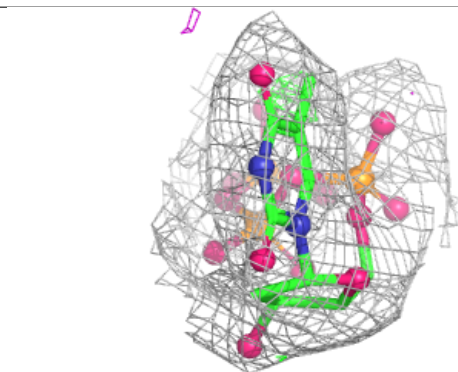
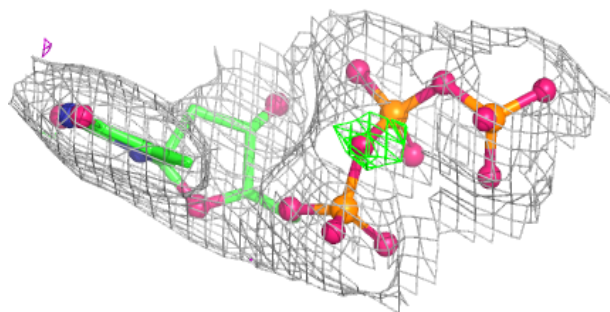
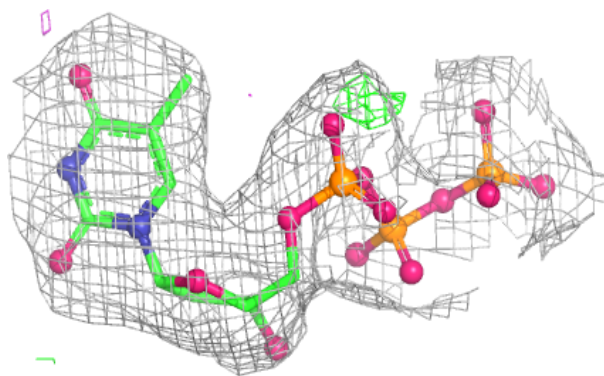
**Electron density around TTP C 805:**

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

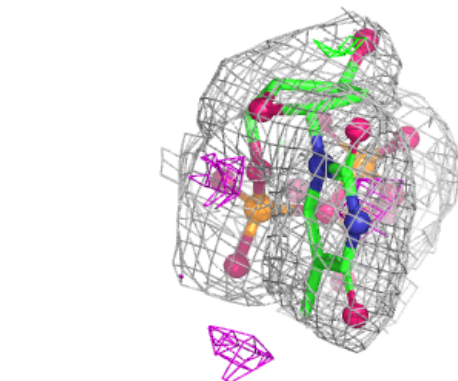
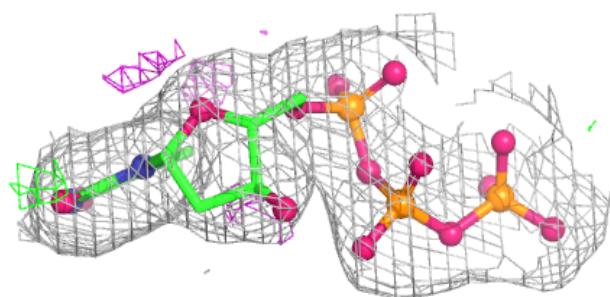
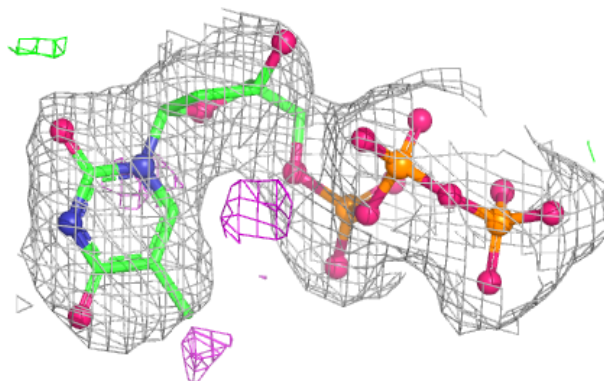


Electron density around TTP B 805:

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

**Electron density around TTP A 805:**

$2mF_o-DF_c$ (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.