



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:34 AM UTC

PDB ID : 4IR1 / pdb_00004ir1
Title : Polymerase-DNA Complex
Authors : Sharma, A.; Nair, D.T.
Deposited on : 2013-01-14
Resolution : 2.38 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

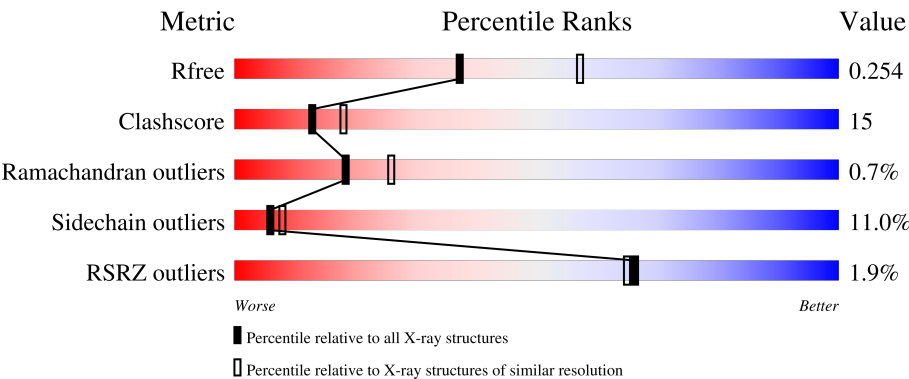
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance ⓘ

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

The reported resolution of this entry is 2.38 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	7164 (2.40-2.36)
Clashscore	190562	7722 (2.40-2.36)
Ramachandran outliers	187476	7626 (2.40-2.36)
Sidechain outliers	187428	7627 (2.40-2.36)
RSRZ outliers	180081	7170 (2.40-2.36)

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 $\leq 5\%$. 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	352	3% 59% 33% 5%
1	F	352	% 70% 24%
2	B	18	44% 50% 6%
2	G	18	6% 67% 33%
3	H	14	57% 43%

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Mol	Chain	Length	Quality of chain
4	C	17	 A horizontal bar chart showing the quality of chain C. The bar is divided into three segments: green (53%), yellow (41%), and orange (6%).

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7080 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 polymerase IV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	F	342	Total	C	N	O	S	0	0	0
			2687	1695	494	484	14			
1	A	342	Total	C	N	O	S	0	0	0
			2687	1695	494	484	14			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	0	GLY	-	expression tag	UNP Q47155
F	1	SER	-	expression tag	UNP Q47155
A	0	GLY	-	expression tag	UNP Q47155
A	1	SER	-	expression tag	UNP Q47155

- Molecule 2 is a DNA chain called DNA (5'-D(*TP*CP*TP*AP*GP*GP*GP*TP*CP*CP*TP*AP*GP*GP*AP*CP*CP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	18	Total	C	N	O	P	0	0	0
			364	174	66	107	17			
2	B	18	Total	C	N	O	P	0	0	0
			364	174	66	107	17			

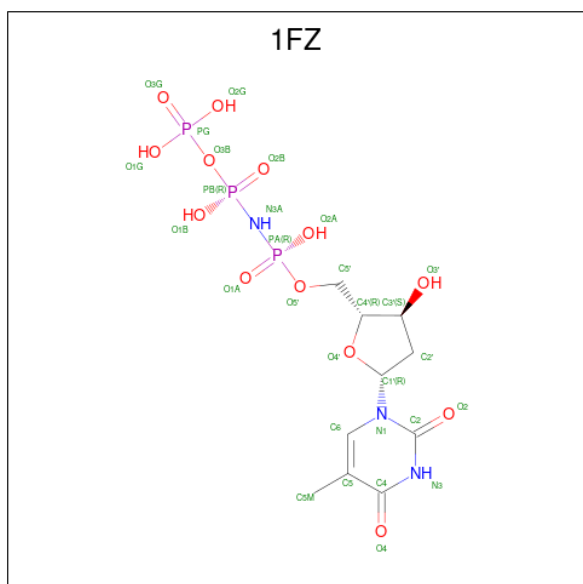
- Molecule 3 is a DNA chain called DNA (5'-D(*GP*GP*GP*TP*CP*CP*TP*AP*GP*GP*AP*CP*CP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	14	Total	C	N	O	P	0	0	0
			284	135	54	82	13			

- Molecule 4 is a DNA chain called DNA (5'-D(*CP*TP*AP*GP*GP*GP*TP*CP*CP*TP*AP*GP*GP*AP*CP*CP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	C	17	Total	C	N	O	P	0	0	0
			344	164	64	100	16			

- Molecule 5 is 5'-O-[(R)-hydroxy{[(R)-hydroxy(phosphonooxy)phosphoryl]amino}phosphoryl]thymidine (CCD ID: 1FZ) (formula: $C_{10}H_{18}N_3O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	F	1	Total 29	C 10	N 3	O 13	P 3	0	0
5	A	1	Total 29	C 10	N 3	O 13	P 3	0	0

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	F	2	Total Mg 2 2	0	0
6	A	2	Total Mg 2 2	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	F	138	Total O 138 138	0	0
7	A	54	Total O 54 54	0	0

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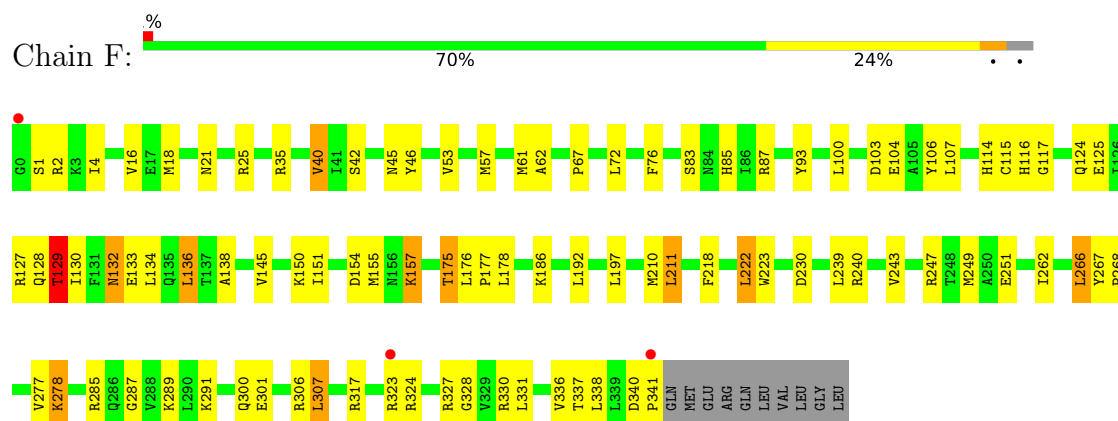
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	G	35	Total 35	O 35	0	0
7	H	27	Total 27	O 27	0	0
7	B	20	Total 20	O 20	0	0
7	C	14	Total 14	O 14	0	0

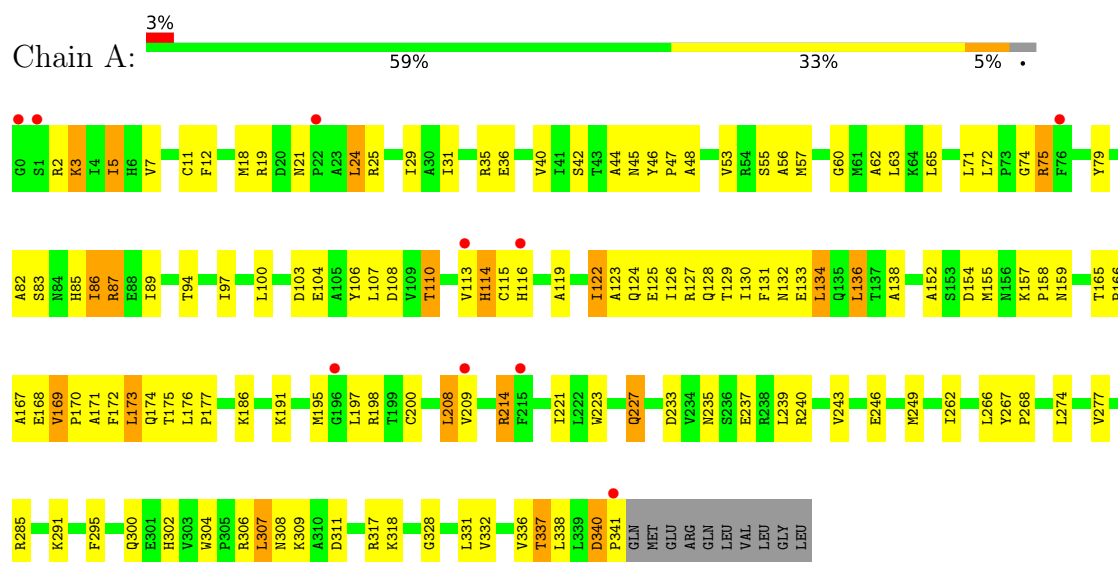
3 Residue-property plots [i](#)

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: DNA polymerase IV



• Molecule 1: DNA polymerase IV



• Molecule 2: DNA (5'-D(*TP*CP*TP*AP*GP*GP*GP*TP*CP*CP*TP*AP*GP*GP*AP*CP*CP*C)-3')





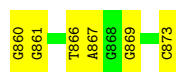
- Molecule 2: DNA (5'-D(*TP*CP*TP*AP*GP*GP*GP*TP*CP*CP*TP*AP*GP*GP*AP*CP*CP*C)-3')

Chain B: 44% 50% 6%



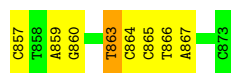
- Molecule 3: DNA (5'-D(*GP*GP*GP*TP*CP*CP*TP*AP*GP*GP*AP*CP*CP*C)-3')

Chain H: 57% 43%



- Molecule 4: DNA (5'-D(*CP*TP*AP*GP*GP*GP*TP*CP*CP*TP*AP*GP*GP*AP*CP*CP*C)-3')

Chain C: 53% 41% 6%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	85.75Å 57.25Å 111.30Å 90.00° 90.44° 90.00°	Depositor
Resolution (Å)	47.66 – 2.38 47.66 – 2.38	Depositor EDS
% Data completeness (in resolution range)	99.8 (47.66-2.38) 99.8 (47.66-2.38)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 2.37Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.214 , 0.269 (Not available) , 0.254	Depositor DCC
R_{free} test set	2199 reflections (3.22%)	wwPDB-VP
Wilson B-factor (Å ²)	49.0	Xtriage
Anisotropy	0.209	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 41.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.023 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7080	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.10% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 1FZ, 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	0.71	1/2738 (0.0%)	1.02	4/3699 (0.1%)
1	F	0.84	0/2738	1.06	7/3699 (0.2%)
2	B	0.39	0/407	1.01	3/626 (0.5%)
2	G	0.45	0/407	0.85	0/626
3	H	0.49	0/318	0.86	0/489
4	C	0.40	0/385	0.96	3/592 (0.5%)
All	All	0.72	1/6993 (0.0%)	1.01	17/9731 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	42	SER	C-N	6.56	1.41	1.33

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	169	VAL	CA-C-N	-10.30	107.62	118.85
1	A	169	VAL	C-N-CA	-10.30	107.62	118.85
2	B	839	DT	C2'-C3'-O3'	9.21	125.31	111.50
1	F	132	ASN	N-CA-C	-6.87	101.46	110.53
1	F	46	TYR	CA-C-N	-5.98	112.70	119.28
1	F	46	TYR	C-N-CA	-5.98	112.70	119.28
2	B	838	DC	C2'-C3'-O3'	-5.89	102.67	111.50
4	C	859	DA	C2'-C3'-O3'	-5.88	102.68	111.50
1	F	76	PHE	N-CA-C	5.79	118.36	111.71
4	C	860	DG	C4'-C3'-O3'	-5.67	101.50	110.00
1	F	129	THR	N-CA-CB	5.58	119.92	110.49
2	B	839	DT	C5'-C4'-C3'	-5.57	106.55	114.90
1	A	173	LEU	N-CA-C	5.44	117.21	111.28
1	A	122	ILE	N-CA-C	-5.33	105.33	110.72

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
4	C	863	DT	C2'-C3'-O3'	5.21	119.31	111.50
1	F	21	ASN	CA-C-N	5.08	125.37	119.47
1	F	21	ASN	C-N-CA	5.08	125.37	119.47

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2687	0	2739	100	0
1	F	2687	0	2739	73	0
2	B	364	0	204	8	0
2	G	364	0	204	7	0
3	H	284	0	158	6	0
4	C	344	0	192	11	0
5	A	29	0	14	3	0
5	F	29	0	14	0	0
6	A	2	0	0	0	0
6	F	2	0	0	0	0
7	A	54	0	0	4	0
7	B	20	0	0	0	0
7	C	14	0	0	1	0
7	F	138	0	0	11	0
7	G	35	0	0	3	0
7	H	27	0	0	1	0
All	All	7080	0	6264	197	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 15.

All (197) 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:340:ASP:HB2	1:A:341:PRO:HD3	1.29	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:18:MET:HE3	1:F:45:ASN:HD22	1.13	1.08
1:F:18:MET:CE	1:F:45:ASN:HD22	1.70	1.02
1:A:317:ARG:HD3	7:A:515:HOH:O	1.57	1.02
1:F:57:MET:HE3	1:F:62:ALA:HA	1.43	1.00
1:F:285:ARG:NH1	3:H:866:DT:OP2	1.97	0.96
1:A:57:MET:HE3	1:A:62:ALA:HA	1.48	0.94
1:A:285:ARG:HB2	1:A:337:THR:HG23	1.52	0.90
1:A:285:ARG:HB2	1:A:337:THR:CG2	2.02	0.90
1:A:195:MET:HE1	1:A:214:ARG:HG2	1.55	0.87
4:C:864:DC:H2''	4:C:865:DC:H5''	1.57	0.87
1:F:57:MET:HE3	1:F:62:ALA:CA	2.04	0.87
1:A:18:MET:HE3	1:A:45:ASN:HD22	1.41	0.86
1:F:125:GLU:O	1:F:128:GLN:O	1.94	0.85
1:A:18:MET:CE	1:A:45:ASN:HD22	1.91	0.83
1:F:57:MET:HG3	1:F:61:MET:HE2	1.61	0.83
1:A:57:MET:HE3	1:A:62:ALA:CA	2.09	0.82
1:A:340:ASP:CB	1:A:341:PRO:HD3	2.10	0.82
1:F:128:GLN:HA	7:F:501:HOH:O	1.82	0.78
1:F:18:MET:HE3	1:F:45:ASN:ND2	1.95	0.78
1:F:150:LYS:NZ	3:H:873:DC:OP1	2.17	0.78
1:F:57:MET:CE	1:F:62:ALA:HA	2.15	0.77
1:F:129:THR:O	1:F:132:ASN:O	2.03	0.76
1:A:130:ILE:O	1:A:134:LEU:HB2	1.88	0.73
1:A:173:LEU:O	1:A:200:CYS:HB2	1.89	0.72
1:A:132:ASN:O	1:A:133:GLU:HB2	1.89	0.72
1:A:340:ASP:HB2	1:A:341:PRO:CD	2.17	0.70
1:F:18:MET:HE2	1:F:25:ARG:O	1.93	0.69
1:A:128:GLN:O	1:A:129:THR:HB	1.91	0.68
1:F:150:LYS:HE3	7:F:583:HOH:O	1.92	0.68
1:A:331:LEU:C	1:A:331:LEU:HD23	2.19	0.67
1:A:166:PRO:O	1:A:169:VAL:HG12	1.95	0.67
1:F:129:THR:N	7:F:501:HOH:O	1.90	0.66
1:F:157:LYS:HD3	7:F:579:HOH:O	1.94	0.66
1:F:300:GLN:HB2	7:F:553:HOH:O	1.95	0.66
4:C:857:DC:H5'	4:C:857:DC:H6	1.61	0.65
1:A:5:ILE:HG23	1:A:107:LEU:HB2	1.77	0.65
1:A:57:MET:CE	1:A:62:ALA:HA	2.24	0.64
1:F:340:ASP:O	1:F:341:PRO:C	2.40	0.63
1:F:287:GLY:HA3	1:F:301:GLU:HG2	1.80	0.63
1:F:249:MET:HE3	1:F:249:MET:HA	1.82	0.62
1:A:308:ASN:ND2	1:A:311:ASP:H	1.97	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:308:ASN:HD22	1:A:311:ASP:H	1.48	0.62
1:A:19:ARG:HD3	1:A:136:LEU:HD13	1.83	0.61
1:A:157:LYS:HE2	5:A:875:1FZ:PG	2.40	0.61
1:A:291:LYS:HB3	1:A:331:LEU:HB3	1.83	0.60
1:A:29:ILE:HG12	1:A:45:ASN:ND2	2.16	0.60
1:F:104:GLU:OE2	1:F:150:LYS:NZ	2.35	0.59
1:F:128:GLN:CA	7:F:501:HOH:O	2.47	0.58
1:F:18:MET:HE1	1:F:45:ASN:HD22	1.67	0.58
1:F:40:VAL:HG13	2:G:840:DA:H5'	1.86	0.58
1:A:3:LYS:H	1:A:110:THR:HG22	1.68	0.58
1:F:145:VAL:HB	1:F:230:ASP:HB3	1.86	0.58
1:F:127:ARG:HD2	1:F:138:ALA:O	2.03	0.57
1:A:82:ALA:O	1:A:86:ILE:HG23	2.04	0.56
1:A:53:VAL:HA	1:A:57:MET:HE1	1.88	0.56
1:A:304:TRP:HD1	1:A:306:ARG:O	1.89	0.56
1:F:240:ARG:NH1	2:G:843:DG:OP1	2.39	0.55
1:A:249:MET:SD	1:A:262:ILE:HD13	2.46	0.55
1:F:128:GLN:O	1:F:129:THR:OG1	2.24	0.55
1:F:18:MET:HE2	1:F:25:ARG:HA	1.90	0.54
1:A:267:TYR:OH	1:A:307:LEU:HD13	2.06	0.54
1:F:124:GLN:OE1	1:F:127:ARG:NH2	2.40	0.54
1:A:169:VAL:HG13	1:A:170:PRO:HD3	1.90	0.54
1:F:278:LYS:HE2	7:F:572:HOH:O	2.06	0.54
1:F:18:MET:CE	1:F:45:ASN:ND2	2.55	0.54
1:F:40:VAL:HG21	2:G:840:DA:C5	2.43	0.54
1:A:85:HIS:O	1:A:89:ILE:HG13	2.08	0.53
4:C:865:DC:H2'	4:C:866:DT:C6	2.43	0.53
1:F:285:ARG:HB2	1:F:337:THR:HB	1.91	0.53
1:A:18:MET:HE1	1:A:45:ASN:HB2	1.90	0.53
4:C:865:DC:C2'	4:C:866:DT:C6	2.91	0.53
1:A:74:GLY:HA2	1:A:79:TYR:OH	2.09	0.53
1:A:227:GLN:NE2	7:A:526:HOH:O	2.40	0.53
1:A:75:ARG:HB2	1:A:75:ARG:HH11	1.74	0.53
1:F:85:HIS:CD2	1:F:134:LEU:HD11	2.44	0.52
1:A:167:ALA:O	1:A:170:PRO:HD2	2.10	0.52
1:A:291:LYS:HE2	1:A:295:PHE:O	2.10	0.52
1:F:85:HIS:CG	1:F:134:LEU:HD11	2.45	0.52
4:C:857:DC:H5'	4:C:857:DC:C6	2.44	0.52
4:C:867:DA:N3	7:C:902:HOH:O	2.34	0.51
1:A:55:SER:O	1:A:56:ALA:HB3	2.10	0.51
1:F:93:TYR:OH	1:F:125:GLU:HG2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:THR:CG2	1:A:107:LEU:HD23	2.40	0.51
1:A:208:LEU:HD23	1:A:223:TRP:CD1	2.46	0.51
1:A:21:ASN:HB3	1:A:24:LEU:HD22	1.93	0.51
3:H:860:DG:N7	7:H:915:HOH:O	2.33	0.51
1:A:46:TYR:CE1	1:A:158:PRO:HG3	2.46	0.51
1:A:246:GLU:HG2	2:B:842:DG:OP2	2.12	0.50
1:F:240:ARG:HH21	1:F:277:VAL:HG21	1.77	0.49
1:F:328:GLY:HA3	7:F:507:HOH:O	2.12	0.49
1:A:31:ILE:HD12	1:A:31:ILE:N	2.28	0.49
4:C:865:DC:H2'	4:C:866:DT:C7	2.43	0.49
1:A:18:MET:CE	1:A:45:ASN:ND2	2.70	0.48
1:A:71:LEU:C	1:A:72:LEU:HD12	2.39	0.48
1:A:18:MET:HE1	1:A:45:ASN:CB	2.43	0.48
1:A:11:CYS:O	1:A:12:PHE:C	2.57	0.48
1:F:330:ARG:HD3	7:G:917:HOH:O	2.14	0.48
1:A:123:ALA:O	1:A:127:ARG:HG3	2.14	0.48
1:A:152:ALA:HA	1:A:155:MET:HG3	1.96	0.48
1:F:57:MET:SD	1:F:61:MET:HE3	2.53	0.48
1:A:128:GLN:HA	7:A:531:HOH:O	2.14	0.48
1:F:18:MET:HE2	1:F:25:ARG:C	2.39	0.47
1:A:53:VAL:HA	1:A:57:MET:CE	2.43	0.47
1:A:235:ASN:C	1:A:237:GLU:H	2.22	0.47
1:F:178:LEU:HD13	1:F:192:LEU:HD13	1.96	0.47
1:F:132:ASN:O	1:F:133:GLU:HB2	2.14	0.47
1:F:324:ARG:HD2	1:F:327:ARG:O	2.14	0.47
1:F:306:ARG:O	1:F:307:LEU:C	2.58	0.47
1:A:128:GLN:O	1:A:129:THR:CB	2.62	0.47
1:A:195:MET:CE	1:A:214:ARG:HG2	2.35	0.47
1:F:2:ARG:NH2	1:F:4:ILE:HD11	2.30	0.47
1:F:115:CYS:C	1:F:117:GLY:N	2.72	0.47
1:A:40:VAL:HG13	2:B:840:DA:H5'	1.97	0.46
1:F:291:LYS:HB3	1:F:331:LEU:HB3	1.98	0.46
1:A:89:ILE:HD13	1:A:129:THR:HG22	1.97	0.46
1:F:106:TYR:O	1:F:107:LEU:HD12	2.15	0.46
1:A:122:ILE:HD12	1:A:123:ALA:N	2.31	0.46
1:A:127:ARG:HD2	1:A:138:ALA:O	2.15	0.46
1:A:317:ARG:NH1	3:H:869:DG:OP1	2.49	0.46
1:A:267:TYR:CG	1:A:309:LYS:HG3	2.51	0.46
1:A:274:LEU:HD22	1:A:307:LEU:HG	1.98	0.45
1:A:302:HIS:NE2	1:A:311:ASP:OD2	2.41	0.45
1:A:45:ASN:CG	1:A:47:PRO:HD2	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:THR:HG22	1:A:107:LEU:HD23	1.98	0.45
1:A:122:ILE:O	1:A:126:ILE:HG13	2.16	0.45
1:F:154:ASP:HA	1:F:157:LYS:HD2	1.99	0.45
1:F:218:PHE:CE2	1:F:222:LEU:HD22	2.52	0.45
1:F:67:PRO:HD2	7:F:544:HOH:O	2.16	0.45
1:F:57:MET:HG3	1:F:61:MET:CE	2.39	0.45
2:B:851:DA:H1'	2:B:852:DC:H5'	1.99	0.45
2:B:851:DA:H2''	2:B:852:DC:H5'	1.98	0.45
1:F:53:VAL:HA	1:F:57:MET:HE1	1.98	0.45
1:A:166:PRO:O	1:A:169:VAL:CG1	2.62	0.45
1:A:158:PRO:O	1:A:159:ASN:C	2.60	0.44
2:G:839:DT:H1'	7:G:925:HOH:O	2.16	0.44
1:A:328:GLY:HA3	7:A:507:HOH:O	2.17	0.44
1:A:157:LYS:HE3	1:A:157:LYS:HB3	1.82	0.44
1:F:289:LYS:NZ	7:F:547:HOH:O	2.50	0.44
1:A:44:ALA:HB1	1:A:48:ALA:HB3	1.99	0.44
1:A:175:THR:O	1:A:177:PRO:HD3	2.17	0.44
1:F:175:THR:O	1:F:177:PRO:HD3	2.18	0.44
1:A:267:TYR:HB3	1:A:268:PRO:HD3	1.99	0.44
1:A:300:GLN:HE22	1:A:318:LYS:NZ	2.15	0.44
2:B:848:DA:H2''	2:B:849:DG:C8	2.53	0.44
4:C:865:DC:H2'	4:C:866:DT:C5	2.53	0.44
1:A:2:ARG:HG3	1:A:2:ARG:HH11	1.83	0.43
1:A:100:LEU:HD11	1:A:106:TYR:CE1	2.53	0.43
1:A:108:ASP:OD2	1:A:110:THR:HG23	2.17	0.43
1:A:267:TYR:CZ	1:A:307:LEU:HD13	2.53	0.43
1:F:267:TYR:HB3	1:F:268:PRO:HD3	1.99	0.43
1:A:5:ILE:HD11	1:A:7:VAL:HG22	1.99	0.43
1:A:35:ARG:HD2	1:A:60:GLY:HA2	1.99	0.43
1:A:304:TRP:CD1	1:A:306:ARG:O	2.70	0.43
2:B:844:DT:H2''	2:B:845:DC:C5'	2.48	0.43
1:A:7:VAL:O	1:A:104:GLU:HA	2.19	0.43
1:A:18:MET:HG2	1:A:25:ARG:HA	2.00	0.43
2:G:843:DG:H2''	2:G:844:DT:H5'	2.01	0.43
1:F:106:TYR:C	1:F:107:LEU:HD12	2.44	0.43
1:F:115:CYS:C	1:F:117:GLY:H	2.27	0.43
2:G:838:DC:H4'	2:G:839:DT:O5'	2.18	0.43
1:A:114:HIS:HB3	1:A:115:CYS:H	1.71	0.43
4:C:865:DC:H2''	4:C:866:DT:C6	2.53	0.43
1:F:211:LEU:HD23	1:F:223:TRP:HB2	2.00	0.43
1:F:300:GLN:HA	3:H:867:DA:OP2	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:C:863:DT:H2''	4:C:864:DC:C6	2.54	0.42
1:A:108:ASP:OD2	1:A:108:ASP:C	2.62	0.42
3:H:860:DG:H2''	3:H:861:DG:OP2	2.20	0.42
1:A:165:THR:OG1	1:A:168:GLU:HG3	2.19	0.42
1:F:83:SER:O	1:F:87:ARG:HG3	2.19	0.42
1:F:130:ILE:O	1:F:134:LEU:HB2	2.20	0.42
1:A:308:ASN:ND2	1:A:311:ASP:N	2.66	0.42
1:F:331:LEU:HD23	1:F:331:LEU:C	2.45	0.41
1:A:63:LEU:HD23	1:A:63:LEU:HA	1.85	0.41
1:F:129:THR:O	1:F:133:GLU:HB2	2.20	0.41
1:F:262:ILE:HG22	1:F:266:LEU:HD22	2.01	0.41
1:A:119:ALA:O	1:A:122:ILE:HG13	2.19	0.41
1:F:323:ARG:HD3	7:F:635:HOH:O	2.21	0.41
1:A:125:GLU:O	1:A:128:GLN:O	2.38	0.41
1:F:100:LEU:HD11	1:F:106:TYR:CE1	2.55	0.41
1:A:46:TYR:N	1:A:47:PRO:CD	2.83	0.41
1:A:308:ASN:HD22	1:A:311:ASP:N	2.16	0.41
1:F:16:VAL:HG22	1:F:136:LEU:HD21	2.03	0.41
1:F:151:ILE:O	1:F:155:MET:HG3	2.21	0.41
1:A:169:VAL:O	1:A:173:LEU:HB2	2.20	0.41
1:A:331:LEU:C	1:A:331:LEU:CD2	2.91	0.41
2:B:838:DC:H6	2:B:838:DC:H2'	1.62	0.41
4:C:865:DC:H2'	4:C:866:DT:H72	2.02	0.41
1:A:83:SER:O	1:A:87:ARG:HD3	2.21	0.41
1:A:171:ALA:O	1:A:172:PHE:C	2.63	0.41
2:G:851:DA:H2'	7:G:921:HOH:O	2.21	0.40
2:B:844:DT:H2''	2:B:845:DC:H5'	2.02	0.40
1:F:72:LEU:HD12	1:F:72:LEU:N	2.37	0.40
1:A:131:PHE:CD2	1:A:131:PHE:C	2.99	0.40
1:F:53:VAL:HG13	1:F:57:MET:HE2	2.03	0.40
1:A:11:CYS:HA	5:A:875:1FZ:O1B	2.22	0.40
5:A:875:1FZ:O5'	5:A:875:1FZ:H1	2.21	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	340/352 (97%)	314 (92%)	24 (7%)	2 (1%)	21	30
1	F	340/352 (97%)	321 (94%)	16 (5%)	3 (1%)	14	20
All	All	680/704 (97%)	635 (93%)	40 (6%)	5 (1%)	18	26

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	114	HIS
1	F	116	HIS
1	F	129	THR
1	A	36	GLU
1	A	340	ASP

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	286/297 (96%)	247 (86%)	39 (14%)	3	4
1	F	286/297 (96%)	262 (92%)	24 (8%)	10	15
All	All	572/594 (96%)	509 (89%)	63 (11%)	6	8

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	1	SER
1	F	35	ARG
1	F	40	VAL
1	F	42	SER
1	F	103	ASP
1	F	136	LEU
1	F	157	LYS

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Mol	Chain	Res	Type
1	F	175	THR
1	F	176	LEU
1	F	186	LYS
1	F	197	LEU
1	F	210	MET
1	F	211	LEU
1	F	222	LEU
1	F	239	LEU
1	F	243	VAL
1	F	247	ARG
1	F	251	GLU
1	F	266	LEU
1	F	278	LYS
1	F	307	LEU
1	F	317	ARG
1	F	336	VAL
1	F	338	LEU
1	A	3	LYS
1	A	5	ILE
1	A	24	LEU
1	A	65	LEU
1	A	75	ARG
1	A	86	ILE
1	A	87	ARG
1	A	97	ILE
1	A	103	ASP
1	A	110	THR
1	A	113	VAL
1	A	114	HIS
1	A	116	HIS
1	A	124	GLN
1	A	134	LEU
1	A	136	LEU
1	A	154	ASP
1	A	174	GLN
1	A	176	LEU
1	A	186	LYS
1	A	191	LYS
1	A	197	LEU
1	A	198	ARG
1	A	208	LEU
1	A	209	VAL

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Mol	Chain	Res	Type
1	A	214	ARG
1	A	221	ILE
1	A	227	GLN
1	A	233	ASP
1	A	239	LEU
1	A	240	ARG
1	A	243	VAL
1	A	266	LEU
1	A	277	VAL
1	A	307	LEU
1	A	332	VAL
1	A	336	VAL
1	A	337	THR
1	A	338	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	128	GLN
1	F	132	ASN
1	F	255	HIS
1	F	297	GLN
1	F	300	GLN
1	A	45	ASN
1	A	84	ASN
1	A	297	GLN
1	A	300	GLN
1	A	308	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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
5	1FZ	A	875	6	29,30,30	2.67	10 (34%)	41,47,47	2.10	12 (29%)
5	1FZ	F	876	6	29,30,30	2.36	8 (27%)	41,47,47	2.19	12 (29%)

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
5	1FZ	A	875	6	-	2/19/34/34	0/2/2/2
5	1FZ	F	876	6	-	5/19/34/34	0/2/2/2

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	875	1FZ	O2-C2	7.72	1.36	1.23
5	A	875	1FZ	O4-C4	6.84	1.36	1.23
5	F	876	1FZ	O2-C2	6.38	1.34	1.23
5	F	876	1FZ	O4-C4	6.04	1.35	1.23
5	A	875	1FZ	PA-O1A	5.29	1.54	1.46
5	F	876	1FZ	PA-O1A	4.16	1.52	1.46
5	A	875	1FZ	PB-O2B	3.74	1.51	1.46
5	F	876	1FZ	PB-O2B	3.49	1.51	1.46
5	A	875	1FZ	C4-C5	3.41	1.50	1.44
5	F	876	1FZ	C4-C5	3.08	1.49	1.44
5	F	876	1FZ	PB-O1B	-2.96	1.48	1.56
5	A	875	1FZ	C6-C5	2.90	1.39	1.34
5	A	875	1FZ	C2-N1	2.78	1.42	1.38
5	F	876	1FZ	C6-C5	2.64	1.38	1.34
5	A	875	1FZ	PB-O3B	2.47	1.62	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	F	876	1FZ	C6-N1	2.47	1.42	1.38
5	A	875	1FZ	C6-N1	2.26	1.41	1.38
5	A	875	1FZ	PB-O1B	-2.14	1.51	1.56

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	876	1FZ	N3-C2-N1	5.80	122.44	114.89
5	A	875	1FZ	N3-C2-N1	5.22	121.68	114.89
5	F	876	1FZ	C4-N3-C2	-4.71	121.16	127.34
5	F	876	1FZ	O4-C4-C5	-4.65	119.59	124.92
5	F	876	1FZ	C5-C4-N3	4.18	118.96	115.32
5	A	875	1FZ	C5-C4-N3	4.00	118.80	115.32
5	A	875	1FZ	C6-N1-C2	-3.98	117.34	121.30
5	F	876	1FZ	O2-C2-N1	-3.97	117.63	122.80
5	F	876	1FZ	O2A-PA-O1A	3.92	118.28	109.87
5	A	875	1FZ	O2A-PA-O1A	3.83	118.08	109.87
5	A	875	1FZ	C4-N3-C2	-3.81	122.35	127.34
5	A	875	1FZ	O4-C4-C5	-3.77	120.60	124.92
5	F	876	1FZ	O1B-PB-O2B	3.50	117.37	109.87
5	F	876	1FZ	C5-C6-N1	-3.43	119.59	123.31
5	A	875	1FZ	O2-C2-N1	-3.19	118.64	122.80
5	A	875	1FZ	O2G-PG-O1G	3.10	119.41	107.80
5	A	875	1FZ	O1A-PA-N3A	-3.08	107.24	111.77
5	F	876	1FZ	C6-N1-C2	-3.07	118.25	121.30
5	F	876	1FZ	O5'-PA-O1A	-2.90	105.68	114.27
5	A	875	1FZ	C5M-C5-C4	2.76	121.72	118.78
5	A	875	1FZ	O5'-PA-O1A	-2.52	106.78	114.27
5	A	875	1FZ	O3B-PB-N3A	-2.22	100.44	106.59
5	F	876	1FZ	O3B-PB-N3A	-2.09	100.78	106.59
5	F	876	1FZ	PA-O5'-C5'	2.08	126.99	119.24

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	F	876	1FZ	PB-N3A-PA-O1A
5	F	876	1FZ	PB-O3B-PG-O1G
5	F	876	1FZ	PB-O3B-PG-O2G
5	A	875	1FZ	PB-N3A-PA-O5'
5	F	876	1FZ	PB-O3B-PG-O3G
5	F	876	1FZ	PB-N3A-PA-O5'

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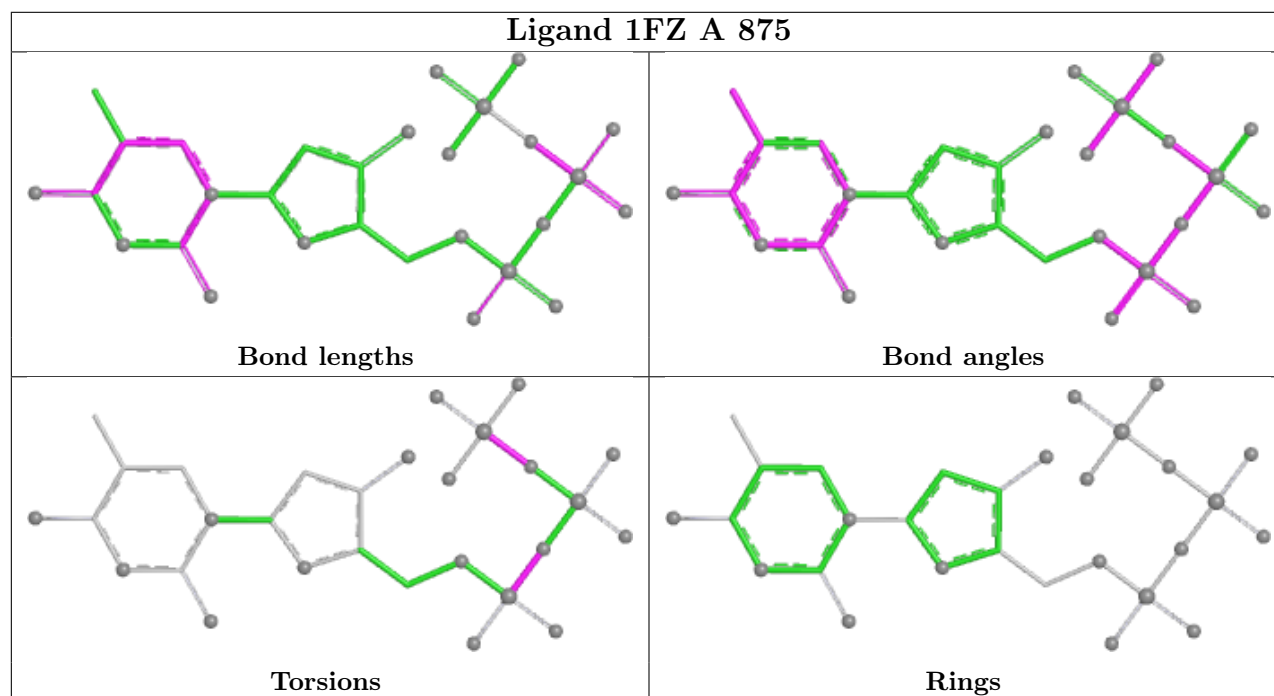
Mol	Chain	Res	Type	Atoms
5	A	875	1FZ	PB-O3B-PG-O3G

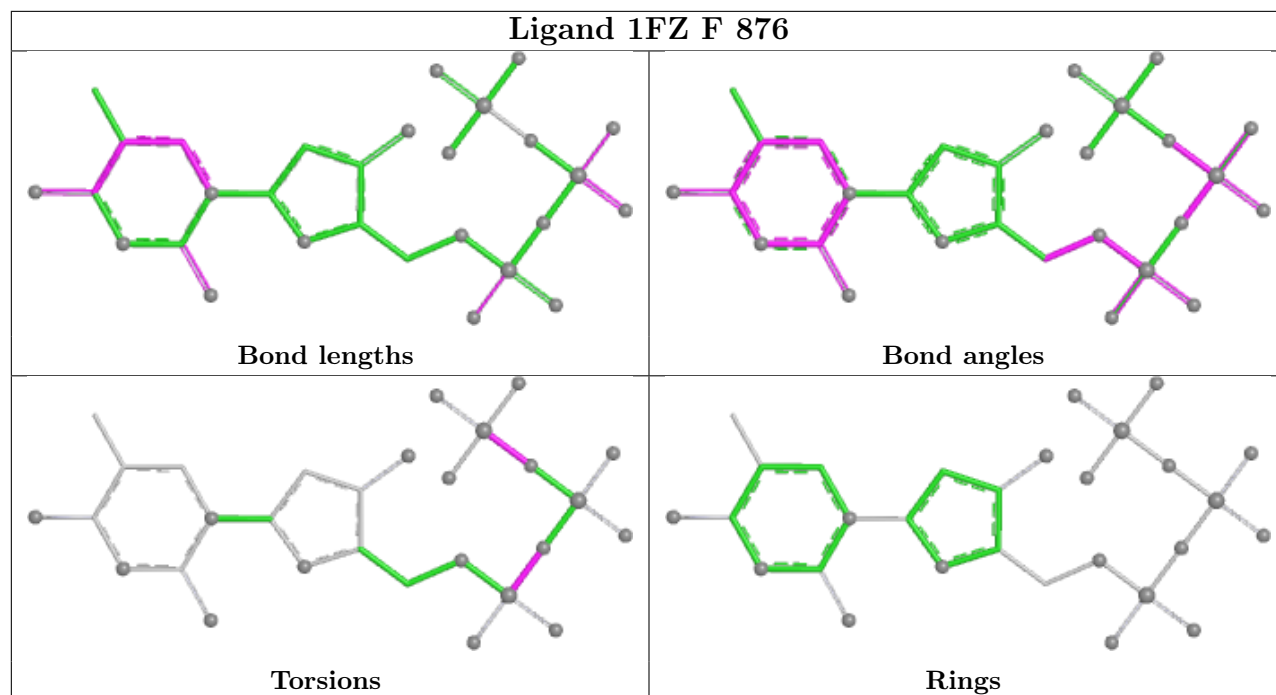
There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	875	1FZ	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	342/352 (97%)	0.39	10 (2%) 53 53	38, 63, 93, 130	0
1	F	342/352 (97%)	0.11	3 (0%) 81 80	32, 45, 63, 80	0
2	B	18/18 (100%)	-0.22	0 100 100	45, 61, 76, 85	0
2	G	18/18 (100%)	-0.20	1 (5%) 30 29	43, 49, 87, 117	0
3	H	14/14 (100%)	-0.14	0 100 100	42, 45, 81, 84	0
4	C	17/17 (100%)	-0.34	0 100 100	47, 51, 76, 81	0
All	All	751/771 (97%)	0.20	14 (1%) 66 65	32, 50, 87, 130	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	341	PRO	5.0
1	F	0	GLY	4.5
1	A	0	GLY	4.3
1	A	113	VAL	3.6
2	G	837	DT	3.2
1	A	209	VAL	3.1
1	A	116	HIS	3.1
1	F	341	PRO	3.0
1	A	1	SER	2.8
1	A	196	GLY	2.7
1	A	76	PHE	2.6
1	A	215	PHE	2.4
1	A	22	PRO	2.2
1	F	323	ARG	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

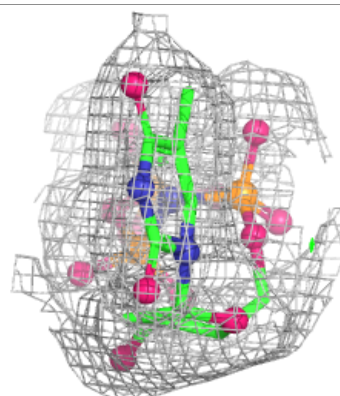
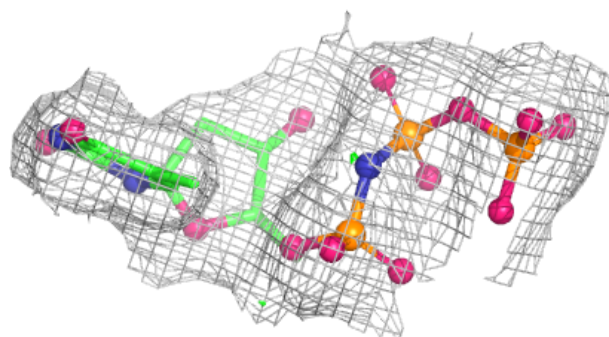
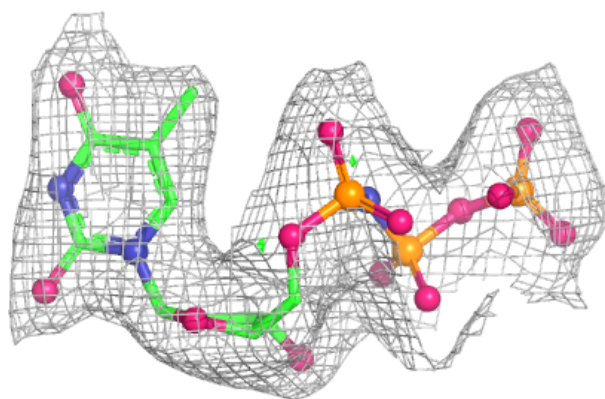
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	MG	F	904	1/1	0.93	0.17	44,44,44,44	0
6	MG	A	902	1/1	0.94	0.06	76,76,76,76	0
5	1FZ	A	875	29/29	0.95	0.07	48,49,52,52	0
5	1FZ	F	876	29/29	0.98	0.07	40,41,41,42	0
6	MG	A	901	1/1	0.99	0.06	49,49,49,49	0
6	MG	F	903	1/1	0.99	0.22	40,40,40,40	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.

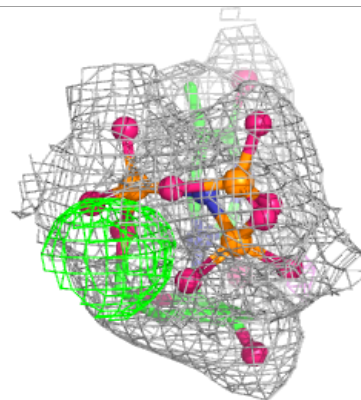
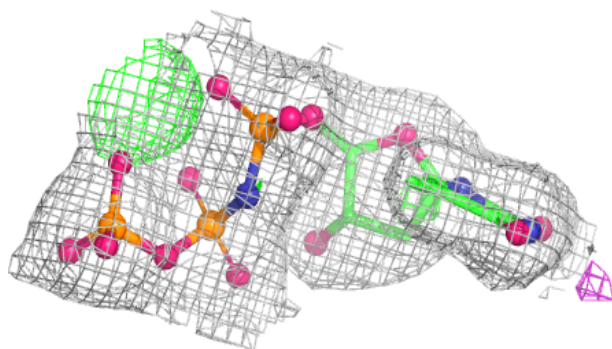
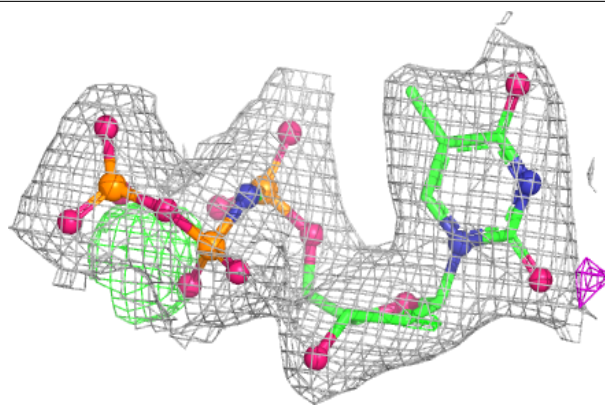
Electron density around 1FZ A 875:

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 1FZ F 876:

$2mF_o - DF_c$ (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.