Full wwPDB X-ray Structure Validation Report (i)

Jun 13, 2024 – 07:07 AM EDT

PDB ID	:	1P7D
Title	:	Crystal structure of the Lambda Integrase (residues 75-356) bound to DNA
Authors	:	Aihara, H.; Kwon, H.J.; Nunes-Duby, S.E.; Landy, A.; Ellenberger, T.
Deposited on	:	2003-05-01
Resolution	:	2.95 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

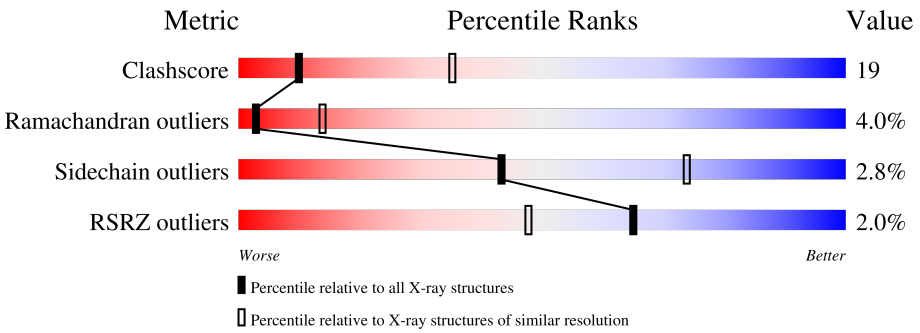
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.95 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))
Clashscore	141614	3462 (3.00-2.92)
Ramachandran outliers	138981	3340 (3.00-2.92)
Sidechain outliers	138945	3343 (3.00-2.92)
RSRZ outliers	127900	2986 (3.00-2.92)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	C	13	38% 54% 8%
1	E	13	54% 38% 8%
2	D	26	38% 62%
2	F	26	4% 35% 65%
3	A	283	% 67% 28% • •
3	B	283	3% 70% 25% 5% •

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 5884 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 DNA chain called 5'-D(*CP*AP*AP*TP*GP*CP*CP*AP*AP*CP*TP*TP *T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	C	13	259	126	45	76	12	0	0	0
1	E	13	259	126	45	76	12	0	0	0

- Molecule 2 is a DNA chain called 26-MER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	D	26	541	257	103	155	26	0	0	0
2	F	26	541	257	103	155	26	0	0	0

- Molecule 3 is a protein called Integrase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
3	A	280	2161	1358	389	402	1	11	0	0	0
3	B	280	2123	1332	383	397	1	10	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

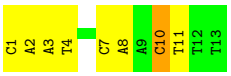
Chain	Residue	Modelled	Actual	Comment	Reference
A	74	MET	VAL	SEE REMARK 999	UNP P03700
A	342	PTR	TYR	MODIFIED RESIDUE	UNP P03700
B	74	MET	VAL	SEE REMARK 999	UNP P03700
B	342	PTR	TYR	MODIFIED RESIDUE	UNP P03700

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.

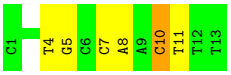
- Molecule 1: 5'-D(*CP*AP*AP*TP*GP*CP*CP*AP*AP*CP*TP*TP*T)-3'

Chain C:



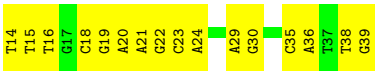
- Molecule 1: 5'-D(*CP*AP*AP*TP*GP*CP*CP*AP*AP*CP*TP*TP*T)-3'

Chain E:



- Molecule 2: 26-MER

Chain D:

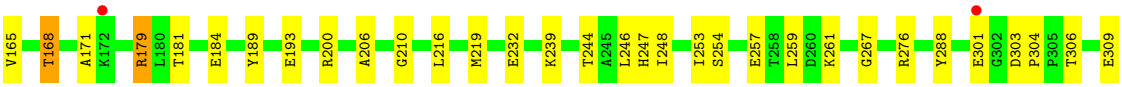


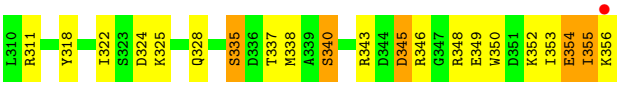
- Molecule 2: 26-MER

Chain F:

- Molecule 3: Integrase

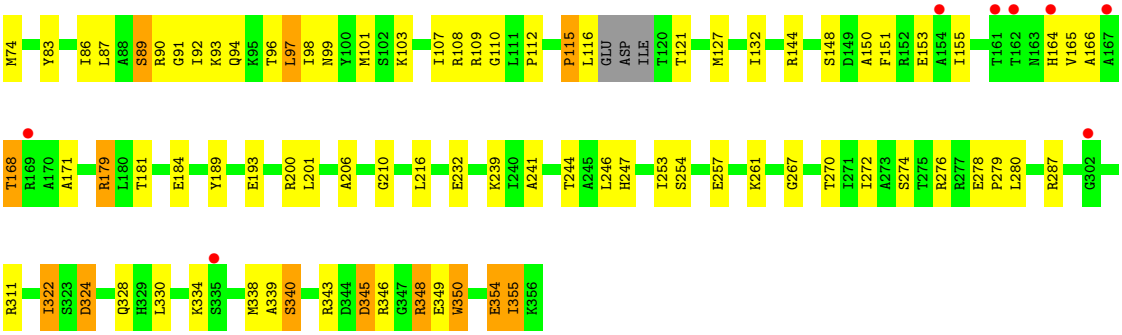
Chain A:





• Molecule 3: Integrase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	59.02Å 106.16Å 73.17Å 90.00° 92.10° 90.00°	Depositor
Resolution (Å)	50.00 – 2.95 35.08 – 2.81	Depositor EDS
% Data completeness (in resolution range)	99.7 (50.00-2.95) 99.3 (35.08-2.81)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.47 (at 2.81Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.231 , 0.259 0.226 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	57.6	Xtriage
Anisotropy	0.133	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 48.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.037 for h,-k,-l	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	5884	wwPDB-VP
Average B, all atoms (Å^2)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.80% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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

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	C	0.49	0/289	0.85	0/443
1	E	0.51	0/289	0.84	0/443
2	D	0.52	0/608	0.87	0/936
2	F	0.51	0/608	0.86	0/936
3	A	0.38	0/2179	0.63	0/2931
3	B	0.38	0/2141	0.63	0/2888
All	All	0.42	0/6114	0.71	0/8577

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
1	E	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	10	DC	Sidechain
1	E	10	DC	Sidechain

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	259	0	148	17	0
1	E	259	0	148	9	0
2	D	541	0	294	17	0
2	F	541	0	294	24	0
3	A	2161	0	2165	78	0
3	B	2123	0	2075	81	0
All	All	5884	0	5124	210	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 19.

All (210) 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)
2:D:21:DA:H2"	2:D:22:DG:H5"	1.08	1.06
3:A:328:GLN:HG2	3:A:343:ARG:HH21	1.31	0.95
3:B:328:GLN:HG2	3:B:343:ARG:HH21	1.31	0.94
3:B:345:ASP:OD2	3:B:349:GLU:HB2	1.68	0.93
2:D:21:DA:C2'	2:D:22:DG:H5"	1.97	0.93
3:B:328:GLN:HE22	3:B:340:SER:HB2	1.33	0.91
3:A:328:GLN:HE22	3:A:340:SER:HB2	1.35	0.91
3:A:352:LYS:HA	3:B:241:ALA:HB3	1.55	0.88
3:B:328:GLN:HE22	3:B:340:SER:CB	1.86	0.88
3:A:328:GLN:HE22	3:A:340:SER:CB	1.88	0.86
3:A:92:ILE:HG23	3:A:96:THR:HB	1.64	0.80
3:A:301:GLU:OE2	3:A:352:LYS:HD2	1.80	0.80
3:A:193:GLU:HA	3:A:200:ARG:NH1	1.97	0.79
3:B:92:ILE:HG23	3:B:96:THR:HB	1.64	0.78
3:A:254:SER:HB3	3:A:257:GLU:HB3	1.67	0.77
3:B:193:GLU:HA	3:B:200:ARG:NH1	2.00	0.76
3:A:116:LEU:HD12	3:A:120:THR:HG23	1.68	0.75
3:B:345:ASP:CG	3:B:349:GLU:HB2	2.07	0.75
2:F:36:DA:H1'	2:F:37:DT:H5'	1.72	0.72
3:B:189:TYR:CE2	3:B:200:ARG:HG2	2.26	0.71
3:B:206:ALA:HA	3:B:216:LEU:HD11	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:116:LEU:HD12	3:A:120:THR:CG2	2.22	0.70
3:A:345:ASP:HB3	3:A:349:GLU:OE2	1.91	0.69
3:B:148:SER:HB2	3:B:165:VAL:HG11	1.76	0.68
3:A:189:TYR:CE2	3:A:200:ARG:HG2	2.28	0.68
3:B:328:GLN:HG2	3:B:343:ARG:NH2	2.08	0.68
3:B:110:GLY:HA3	3:B:127:MET:CE	2.24	0.67
3:A:206:ALA:HA	3:A:216:LEU:HD11	1.76	0.67
3:A:324:ASP:HB2	3:A:343:ARG:CZ	2.25	0.67
3:A:110:GLY:HA3	3:A:127:MET:CE	2.25	0.66
3:A:324:ASP:HB2	3:A:343:ARG:NH2	2.11	0.66
3:B:110:GLY:HA3	3:B:127:MET:HE2	1.78	0.66
1:C:3:DA:H2"	1:C:4:DT:C5'	2.26	0.66
2:F:14:DT:H2'	2:F:15:DT:H72	1.77	0.66
3:B:165:VAL:O	3:B:168:THR:HG22	1.96	0.65
3:A:193:GLU:HA	3:A:200:ARG:HH11	1.61	0.65
3:B:108:ARG:HG3	3:B:108:ARG:HH11	1.61	0.65
2:D:23:DC:H2"	2:D:24:DA:C8	2.32	0.64
1:C:3:DA:H2"	1:C:4:DT:H5'	1.79	0.64
3:A:328:GLN:HG2	3:A:343:ARG:NH2	2.08	0.64
1:C:10:DC:H2'	1:C:11:DT:H72	1.80	0.64
3:A:116:LEU:CD1	3:A:120:THR:HG23	2.28	0.63
2:D:38:DT:H1'	2:D:39:DG:C8	2.33	0.63
3:A:345:ASP:OD1	3:A:348:ARG:HB2	1.99	0.63
3:B:132:ILE:HD11	3:B:171:ALA:HB2	1.82	0.62
3:B:193:GLU:HA	3:B:200:ARG:HH11	1.65	0.61
3:A:132:ILE:HD11	3:A:171:ALA:HB2	1.82	0.61
3:A:108:ARG:HG3	3:A:108:ARG:HH11	1.64	0.60
2:D:14:DT:OP1	3:A:340:SER:HA	2.02	0.60
2:F:39:DG:O3'	3:B:276:ARG:CZ	2.49	0.60
3:B:254:SER:HB3	3:B:257:GLU:HB3	1.83	0.60
1:C:10:DC:C6	1:C:11:DT:H72	2.36	0.60
3:B:179:ARG:O	3:B:348:ARG:HD2	2.01	0.60
3:B:276:ARG:HG2	3:B:276:ARG:HH11	1.67	0.60
1:E:10:DC:C6	1:E:11:DT:H72	2.36	0.59
3:B:328:GLN:HG3	3:B:338:MET:SD	2.43	0.59
2:F:38:DT:H2"	2:F:39:DG:OP2	2.02	0.59
3:A:276:ARG:HG2	3:A:276:ARG:HH11	1.66	0.59
3:A:189:TYR:CZ	3:A:200:ARG:HG2	2.38	0.59
3:A:328:GLN:HG3	3:A:338:MET:HG2	1.83	0.59
1:E:7:DC:H1'	1:E:8:DA:C5'	2.33	0.58
3:A:324:ASP:HB2	3:A:343:ARG:NH1	2.18	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:10:DC:H2'	1:E:11:DT:H72	1.83	0.58
2:D:29:DA:C2'	2:D:30:DG:C8	2.86	0.58
1:C:7:DC:H1'	1:C:8:DA:C5'	2.34	0.57
1:E:7:DC:H1'	1:E:8:DA:H5"	1.87	0.57
3:B:189:TYR:CZ	3:B:200:ARG:HG2	2.39	0.57
1:E:7:DC:H2"	1:E:8:DA:H5'	1.88	0.56
3:A:110:GLY:HA3	3:A:127:MET:HE2	1.86	0.56
3:B:115:PRO:HG2	3:B:116:LEU:H	1.70	0.56
2:F:25:DA:H2"	2:F:26:DA:OP2	2.06	0.56
3:A:110:GLY:HA3	3:A:127:MET:HE3	1.88	0.55
2:D:29:DA:H2"	2:D:30:DG:C8	2.41	0.55
3:A:115:PRO:HG2	3:A:116:LEU:H	1.71	0.55
2:F:29:DA:C2'	2:F:30:DG:C8	2.89	0.55
1:C:7:DC:H1'	1:C:8:DA:H5"	1.88	0.55
3:A:355:ILE:HG22	3:A:355:ILE:O	2.06	0.55
3:A:165:VAL:O	3:A:168:THR:HG22	2.07	0.54
3:B:181:THR:OG1	3:B:184:GLU:HG3	2.07	0.54
3:B:108:ARG:HG3	3:B:108:ARG:NH1	2.22	0.54
3:B:121:THR:HG1	3:B:164:HIS:HD1	1.55	0.54
3:B:89:SER:O	3:B:91:GLY:N	2.39	0.53
3:B:97:LEU:HD22	3:B:101:MET:HG2	1.90	0.53
2:D:14:DT:P	3:A:340:SER:HA	2.48	0.53
2:D:18:DC:H2"	2:D:19:DG:C8	2.43	0.53
3:B:354:GLU:O	3:B:355:ILE:HG13	2.09	0.53
2:D:38:DT:H1'	2:D:39:DG:N7	2.23	0.53
2:F:38:DT:H1'	2:F:39:DG:C8	2.43	0.53
2:F:15:DT:H3'	3:B:334:LYS:HD2	1.91	0.53
2:D:15:DT:H1'	2:D:16:DT:H5'	1.92	0.52
2:D:29:DA:H2'	2:D:30:DG:C8	2.44	0.52
3:B:166:ALA:C	3:B:168:THR:H	2.13	0.52
3:A:99:ASN:O	3:A:103:LYS:HG2	2.09	0.52
3:A:350:TRP:CE3	3:B:239:LYS:HB3	2.44	0.52
1:C:10:DC:H2'	1:C:11:DT:C7	2.39	0.52
2:D:29:DA:C2'	2:D:30:DG:H8	2.23	0.52
3:B:148:SER:HB2	3:B:165:VAL:CG1	2.39	0.52
1:C:10:DC:OP1	3:A:136:LYS:NZ	2.42	0.51
3:A:97:LEU:HD22	3:A:101:MET:HG2	1.92	0.51
3:A:120:THR:HG22	3:A:123:GLU:OE1	2.10	0.51
2:D:19:DG:H2"	2:D:20:DA:O5'	2.11	0.51
1:C:7:DC:H2"	1:C:8:DA:H5'	1.91	0.51
3:B:254:SER:HB3	3:B:257:GLU:CB	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:354:GLU:HG3	3:A:355:ILE:H	1.76	0.51
2:F:29:DA:H2"	2:F:30:DG:C8	2.44	0.51
2:F:24:DA:H1'	2:F:25:DA:H5'	1.92	0.50
2:D:29:DA:H2"	2:D:30:DG:H8	1.75	0.50
3:B:110:GLY:HA3	3:B:127:MET:HE3	1.94	0.50
2:F:29:DA:H2'	2:F:30:DG:C8	2.46	0.50
3:A:108:ARG:HG3	3:A:108:ARG:NH1	2.25	0.50
3:B:74:MET:O	3:B:115:PRO:HA	2.11	0.50
3:A:74:MET:O	3:A:115:PRO:HA	2.12	0.50
3:B:349:GLU:O	3:B:350:TRP:HB2	2.11	0.50
3:A:232:GLU:OE1	3:A:239:LYS:HE2	2.12	0.49
3:B:94:GLN:O	3:B:98:ILE:HG13	2.12	0.49
3:A:356:LYS:HB3	3:B:322:ILE:CG2	2.41	0.49
1:C:2:DA:H2"	1:C:3:DA:OP2	2.12	0.49
1:C:1:DC:H2"	1:C:2:DA:C8	2.48	0.49
2:F:33:DG:N7	3:B:287:ARG:NH2	2.60	0.49
1:E:10:DC:H2'	1:E:11:DT:C7	2.42	0.49
2:F:36:DA:H2"	2:F:37:DT:OP2	2.11	0.49
3:A:181:THR:OG1	3:A:184:GLU:HG3	2.12	0.49
3:A:353:ILE:CD1	3:B:330:LEU:HA	2.43	0.48
1:C:3:DA:H1'	1:C:4:DT:H5"	1.93	0.48
3:A:144:ARG:HE	3:A:168:THR:HG23	1.78	0.48
2:F:28:DA:H3'	3:B:93:LYS:HD3	1.95	0.48
3:A:83:TYR:O	3:A:86:ILE:HB	2.13	0.48
3:B:99:ASN:O	3:B:103:LYS:HG2	2.12	0.48
2:F:29:DA:C2'	2:F:30:DG:H8	2.27	0.48
2:D:23:DC:H2"	2:D:24:DA:H8	1.76	0.47
3:A:244:THR:C	3:A:246:LEU:H	2.16	0.47
3:A:276:ARG:HG2	3:A:276:ARG:NH1	2.29	0.47
3:A:144:ARG:HE	3:A:168:THR:CG2	2.28	0.47
3:B:232:GLU:OE1	3:B:239:LYS:HE2	2.15	0.47
3:A:92:ILE:HG23	3:A:96:THR:CB	2.41	0.46
3:B:349:GLU:O	3:B:350:TRP:CB	2.63	0.46
3:B:244:THR:C	3:B:246:LEU:H	2.18	0.46
2:F:14:DT:OP1	3:B:340:SER:HA	2.15	0.46
3:B:276:ARG:HG2	3:B:276:ARG:NH1	2.30	0.46
3:A:153:GLU:HA	3:A:153:GLU:OE1	2.16	0.46
3:B:345:ASP:OD1	3:B:349:GLU:N	2.48	0.46
3:B:83:TYR:O	3:B:86:ILE:HB	2.16	0.46
2:F:29:DA:H2"	2:F:30:DG:H8	1.80	0.46
3:B:115:PRO:O	3:B:116:LEU:O	2.34	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:210:GLY:O	3:B:311:ARG:NH1	2.49	0.45
3:B:144:ARG:HE	3:B:168:THR:HG23	1.80	0.45
3:B:355:ILE:HG22	3:B:355:ILE:O	2.16	0.45
1:E:7:DC:H1'	1:E:8:DA:H5'	1.98	0.45
3:B:153:GLU:OE1	3:B:153:GLU:HA	2.16	0.45
1:E:4:DT:H2"	1:E:5:DG:OP2	2.17	0.44
2:F:39:DG:O3'	3:B:276:ARG:NH2	2.50	0.44
3:A:89:SER:O	3:A:91:GLY:N	2.50	0.44
3:B:328:GLN:HE21	3:B:343:ARG:NE	2.15	0.44
1:C:3:DA:H2'	3:A:288:TYR:OH	2.16	0.44
3:B:328:GLN:HE22	3:B:340:SER:HB3	1.76	0.44
3:A:94:GLN:O	3:A:98:ILE:HG13	2.17	0.44
1:C:3:DA:C2'	1:C:4:DT:H5"	2.48	0.44
3:B:107:ILE:HA	3:B:127:MET:CE	2.48	0.44
3:B:83:TYR:CE1	3:B:150:ALA:HA	2.52	0.44
3:B:103:LYS:O	3:B:107:ILE:HG13	2.16	0.44
3:A:163:ASN:C	3:A:165:VAL:H	2.21	0.43
3:A:210:GLY:O	3:A:311:ARG:NH1	2.51	0.43
3:B:148:SER:HA	3:B:165:VAL:HG21	2.00	0.43
3:B:92:ILE:HG23	3:B:96:THR:CB	2.42	0.43
3:B:338:MET:CE	3:B:343:ARG:HH22	2.31	0.43
1:C:10:DC:H2"	1:C:11:DT:C6	2.53	0.43
2:F:35:DC:H2"	2:F:36:DA:OP2	2.18	0.43
1:C:7:DC:H1'	1:C:8:DA:H5'	2.00	0.43
3:A:115:PRO:O	3:A:116:LEU:O	2.37	0.43
3:B:144:ARG:HE	3:B:168:THR:CG2	2.30	0.43
1:E:10:DC:H2"	1:E:11:DT:C6	2.54	0.43
3:A:247:HIS:HB2	3:A:253:ILE:O	2.19	0.43
3:A:107:ILE:HA	3:A:127:MET:CE	2.49	0.43
3:B:201:LEU:HD21	3:B:261:LYS:HE3	2.01	0.43
1:C:3:DA:H2"	1:C:4:DT:H5"	2.01	0.43
2:F:18:DC:H2"	2:F:19:DG:C8	2.54	0.43
3:A:151:PHE:O	3:A:155:ILE:HG13	2.19	0.43
3:A:179:ARG:HG3	3:A:346:ARG:HB2	2.01	0.43
3:B:179:ARG:HG3	3:B:346:ARG:HB2	2.01	0.42
3:B:270:THR:HG23	3:B:279:PRO:HG3	2.01	0.42
3:A:103:LYS:O	3:A:107:ILE:HG13	2.19	0.42
3:A:328:GLN:HE21	3:A:343:ARG:NE	2.17	0.42
2:F:16:DT:P	3:B:334:LYS:HD2	2.59	0.42
3:A:75:THR:CB	3:A:113:ASP:OD1	2.68	0.42
3:A:216:LEU:HD23	3:A:219:MET:SD	2.59	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:83:TYR:CE2	3:B:87:LEU:HD11	2.54	0.42
3:B:254:SER:O	3:B:257:GLU:HB3	2.20	0.41
3:B:354:GLU:O	3:B:355:ILE:CB	2.67	0.41
3:A:257:GLU:O	3:A:261:LYS:HG3	2.19	0.41
3:B:324:ASP:HB3	3:B:343:ARG:NH1	2.35	0.41
2:F:15:DT:C3'	3:B:334:LYS:HD2	2.50	0.41
3:A:83:TYR:CE1	3:A:150:ALA:HA	2.55	0.41
3:A:303:ASP:HA	3:A:304:PRO:HD3	1.82	0.41
3:A:306:THR:O	3:A:309:GLU:HB2	2.20	0.41
3:B:94:GLN:HA	3:B:94:GLN:NE2	2.34	0.41
3:B:151:PHE:O	3:B:155:ILE:HG13	2.21	0.41
3:B:274:SER:N	3:B:278:GLU:O	2.37	0.41
3:A:152:ARG:HG3	3:A:163:ASN:ND2	2.35	0.41
3:A:216:LEU:HD23	3:A:216:LEU:HA	1.91	0.41
3:A:335:SER:C	3:A:337:THR:N	2.74	0.41
3:B:247:HIS:HB2	3:B:253:ILE:O	2.20	0.41
2:F:28:DA:H1'	2:F:29:DA:H5"	2.03	0.41
2:F:36:DA:H1'	2:F:37:DT:C5'	2.47	0.41
3:A:149:ASP:HA	3:A:152:ARG:NH1	2.36	0.40
2:D:35:DC:H2"	2:D:36:DA:OP2	2.22	0.40
3:A:259:LEU:HD23	3:A:259:LEU:HA	1.91	0.40
3:A:248:ILE:HD11	3:A:318:TYR:CE2	2.57	0.40
3:B:272:ILE:HG22	3:B:280:LEU:HD12	2.04	0.40
3:A:181:THR:HG23	3:A:184:GLU:OE1	2.21	0.40
3:A:324:ASP:OD2	3:A:325:LYS:N	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
3	A	275/283 (97%)	245 (89%)	19 (7%)	11 (4%)	3 14

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
3	B	275/283 (97%)	242 (88%)	22 (8%)	11 (4%)	3 14
All	All	550/566 (97%)	487 (88%)	41 (8%)	22 (4%)	3 14

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	115	PRO
3	A	340	SER
3	A	354	GLU
3	A	355	ILE
3	B	90	ARG
3	B	115	PRO
3	B	340	SER
3	B	354	GLU
3	B	355	ILE
3	A	267	GLY
3	B	89	SER
3	B	267	GLY
3	B	339	ALA
3	A	112	PRO
3	A	164	HIS
3	A	335	SER
3	B	112	PRO
3	A	89	SER
3	A	90	ARG
3	A	322	ILE
3	B	350	TRP
3	B	322	ILE

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
3	A	219/236 (93%)	214 (98%)	5 (2%)	50 78
3	B	208/236 (88%)	201 (97%)	7 (3%)	37 69

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	427/472 (90%)	415 (97%)	12 (3%)	43 74

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	97	LEU
3	A	147	LEU
3	A	168	THR
3	A	179	ARG
3	A	345	ASP
3	B	97	LEU
3	B	109	ARG
3	B	168	THR
3	B	179	ARG
3	B	324	ASP
3	B	345	ASP
3	B	348	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	328	GLN
3	B	94	GLN
3	B	328	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	PTR	B	342	3,1	11,15,17	0.73	0	12,19,24	0.57	0
3	PTR	A	342	3,1	11,15,17	0.64	0	12,19,24	0.53	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PTR	B	342	3,1	-	1/7/10/13	0/1/1/1
3	PTR	A	342	3,1	-	1/7/10/13	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	342	PTR	C-CA-CB-CG
3	B	342	PTR	C-CA-CB-CG

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	C	13/13 (100%)	-0.10	0	100	100	23, 37, 72, 89	0
1	E	13/13 (100%)	-0.05	0	100	100	25, 41, 81, 98	0
2	D	26/26 (100%)	0.15	0	100	100	17, 48, 98, 108	0
2	F	26/26 (100%)	0.31	1 (3%)	40	26	19, 54, 123, 143	0
3	A	279/283 (98%)	0.05	4 (1%)	75	59	20, 49, 69, 79	0
3	B	279/283 (98%)	0.06	8 (2%)	51	35	20, 49, 72, 82	0
All	All	636/644 (98%)	0.06	13 (2%)	65	48	17, 49, 73, 143	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	B	335	SER	3.4
3	B	162	THR	3.3
3	B	164	HIS	2.8
3	A	74	MET	2.6
3	B	169	ARG	2.6
3	A	172	LYS	2.5
2	F	39	DG	2.4
3	B	167	ALA	2.4
3	A	356	LYS	2.4
3	B	302	GLY	2.2
3	A	301	GLU	2.1
3	B	154	ALA	2.1
3	B	161	THR	2.1

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	PTR	B	342	15/17	0.97	0.20	27,34,45,45	0
3	PTR	A	342	15/17	0.98	0.19	30,34,47,47	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.