Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 11:32 PM GMT

PDB ID	:	1XO0
Title	:	High resolution structure of the holliday junction intermediate in cre-loxp site-specific recombination
Authors	:	Ghosh, K.; Lau, C.K.; Guo, F.; Segall, A.M.; Van Duyne, G.D.
Deposited on	:	2004-10-05
Resolution	:	2.00 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

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

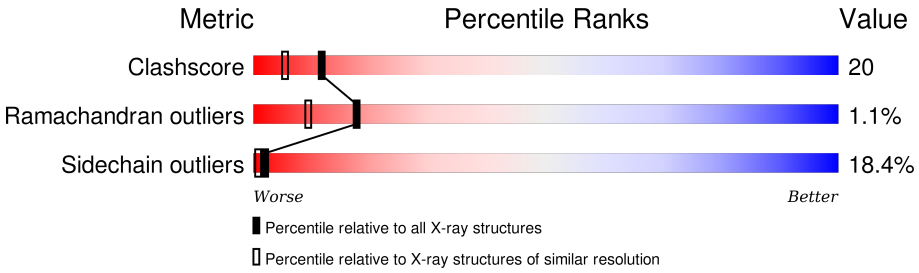
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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 2.00 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))
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

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 on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	C	35	43% 54% .
2	D	35	29% 60% 11%
3	A	324	59% 35% 6% .
3	B	324	63% 28% 6% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7017 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 loxP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	35	Total	C	N	O	P	0	0	0
			715	346	125	210	34			

- Molecule 2 is a DNA chain called loxP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	35	Total	C	N	O	P	0	0	0
			713	345	126	208	34			

- Molecule 3 is a protein called Recombinase CRE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	322	Total	C	N	O	S	0	0	0
			2548	1584	484	465	15			
3	B	322	Total	C	N	O	S	0	0	0
			2548	1584	484	465	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	173	LYS	ARG	ENGINEERED	UNP P06956
B	173	LYS	ARG	ENGINEERED	UNP P06956

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	154	Total	O	0	0
			154	154		
4	B	191	Total	O	0	0
			191	191		
4	C	60	Total	O	0	0
			60	60		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	D	88	Total O 88 88	0	0

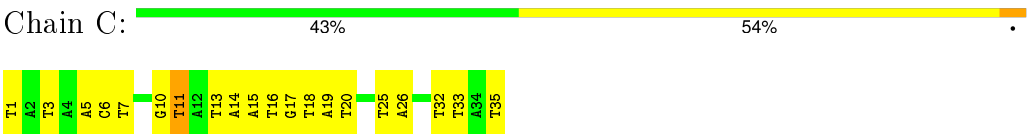
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

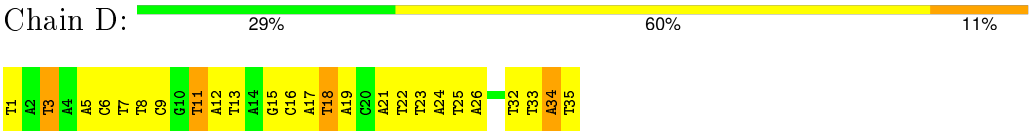
- Molecule 1: loxP

Chain C:



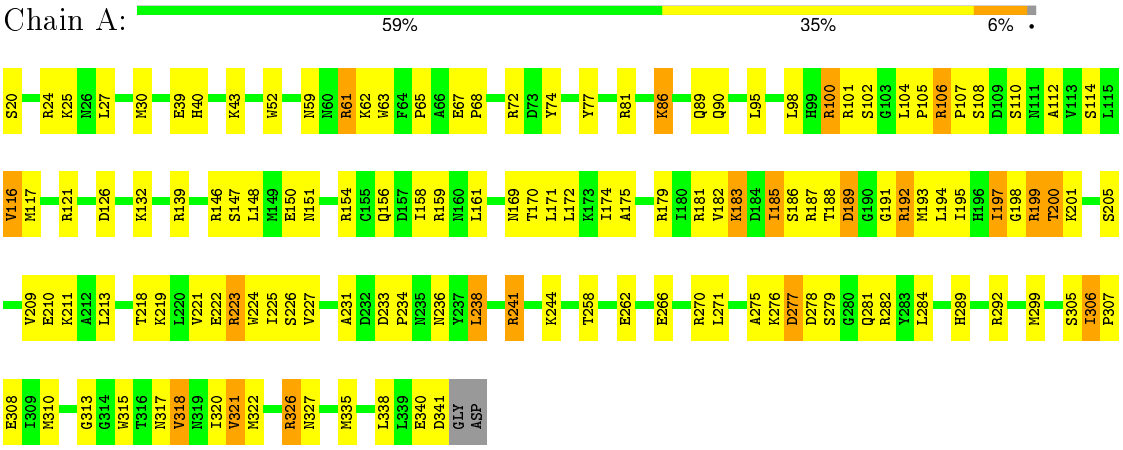
- Molecule 2: loxP

Chain D:



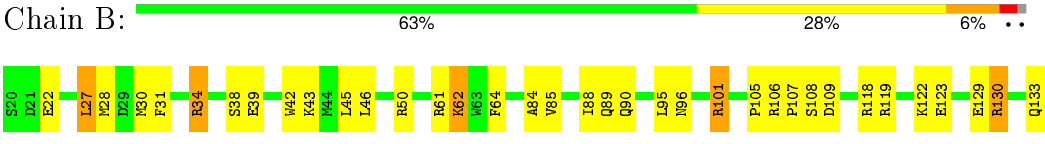
- Molecule 3: Recombinase CRE

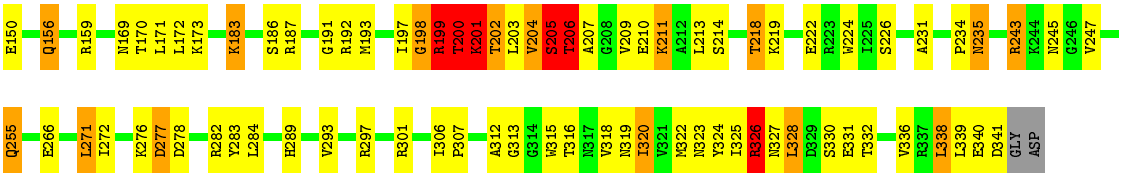
Chain A:



- Molecule 3: Recombinase CRE

Chain B:




E150
Q156
R159
N169
T170
L171
L172
K173
K183
S186
R187
G191
R192
M193
I197
G198
R199
T200
K201
T202
L203
V204
S205
T206
A207
G208
V209
E210
K211
A212
L213
S214
T218
K219
E222
R223
W224
I225
S226
A231
P234
N235
R243
K244
N245
G246
V247
Q255
E266
L271
I272
K276
D277
D278
R282
Y283
L284
H289
V293
R297
R301
I306
P307
A312
G313
G314
W315
T316
N317
V318
N319
I320
V321
M322
N323
Y324
I325
R326
N327
L328
D329
S330
E331
T332
V336
R337
L338
L339
E340
D341
GLY
ASP

4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	107.40Å 123.00Å 180.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	26.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (26.00-2.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.219 , 0.265	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7017	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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	1.39	11/801 (1.4%)	1.68	15/1235 (1.2%)
2	D	1.93	11/799 (1.4%)	1.78	15/1231 (1.2%)
3	A	0.40	0/2589	0.57	0/3490
3	B	1.01	6/2589 (0.2%)	1.13	15/3490 (0.4%)
All	All	1.06	28/6778 (0.4%)	1.17	45/9446 (0.5%)

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
3	B	0	3

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	DT	O3'-P	42.24	2.11	1.61
3	B	199	ARG	CD-NE	-25.77	1.02	1.46
1	C	1	DT	O3'-P	24.51	1.90	1.61
3	B	202	THR	C-N	22.71	1.86	1.34
3	B	206	THR	C-N	-17.07	0.94	1.34
3	B	198	GLY	C-N	-16.93	0.95	1.34
3	B	199	ARG	CG-CD	-15.43	1.13	1.51
3	B	205	SER	C-N	-12.24	1.05	1.34
2	D	13	DT	O3'-P	-11.57	1.47	1.61
2	D	26	DA	O3'-P	-8.53	1.50	1.61
2	D	33	DT	C5-C7	7.51	1.54	1.50
2	D	11	DT	C5-C7	7.50	1.54	1.50
2	D	32	DT	C5-C7	7.39	1.54	1.50
2	D	23	DT	C5-C7	7.38	1.54	1.50
2	D	13	DT	C5-C7	7.36	1.54	1.50
1	C	3	DT	C5-C7	7.34	1.54	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	33	DT	C5-C7	7.31	1.54	1.50
2	D	25	DT	C5-C7	7.30	1.54	1.50
1	C	13	DT	C5-C7	7.30	1.54	1.50
1	C	32	DT	C5-C7	7.27	1.54	1.50
1	C	35	DT	C5-C7	7.23	1.54	1.50
1	C	25	DT	C5-C7	7.20	1.54	1.50
1	C	11	DT	C5-C7	7.19	1.54	1.50
2	D	3	DT	C5-C7	7.16	1.54	1.50
2	D	18	DT	C5-C7	7.11	1.54	1.50
1	C	14	DA	O3'-P	-7.08	1.52	1.61
1	C	3	DT	O3'-P	-6.89	1.52	1.61
1	C	25	DT	O3'-P	-6.24	1.53	1.61

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	1	DT	O3'-P-O5'	35.94	172.28	104.00
2	D	1	DT	OP1-P-O3'	-30.19	38.78	105.20
2	D	1	DT	P-O3'-C3'	25.10	149.82	119.70
3	B	198	GLY	O-C-N	-24.75	83.10	122.70
3	B	201	LYS	C-N-CA	-21.53	67.88	121.70
3	B	201	LYS	CA-C-N	-21.35	70.23	117.20
3	B	202	THR	C-N-CA	-16.00	81.70	121.70
1	C	1	DT	OP1-P-O3'	-15.77	70.50	105.20
2	D	1	DT	OP2-P-O3'	15.67	139.68	105.20
3	B	199	ARG	CA-CB-CG	-15.13	80.12	113.40
3	B	202	THR	O-C-N	-14.45	99.58	122.70
3	B	202	THR	CA-C-N	-13.78	86.88	117.20
3	B	198	GLY	CA-C-N	13.32	146.51	117.20
3	B	204	VAL	O-C-N	-11.65	104.06	122.70
1	C	1	DT	OP2-P-O3'	-11.35	80.23	105.20
2	D	13	DT	P-O3'-C3'	8.90	130.38	119.70
2	D	34	DA	P-O3'-C3'	-8.40	109.62	119.70
3	B	204	VAL	CA-C-N	8.20	135.25	117.20
3	B	199	ARG	NE-CZ-NH2	7.59	124.10	120.30
3	B	326	ARG	NE-CZ-NH2	7.02	123.81	120.30
2	D	18	DT	O4'-C4'-C3'	-6.25	102.00	104.50
1	C	25	DT	P-O3'-C3'	6.11	127.03	119.70
2	D	1	DT	O3'-P-O5'	6.03	115.45	104.00
3	B	322	MET	CG-SD-CE	5.93	109.68	100.20
3	B	201	LYS	O-C-N	5.71	131.84	122.70
1	C	1	DT	P-O3'-C3'	5.68	126.51	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	25	DT	C4-C5-C6	5.50	121.30	118.00
1	C	3	DT	P-O3'-C3'	5.49	126.29	119.70
1	C	13	DT	C4-C5-C6	5.43	121.26	118.00
2	D	33	DT	C4-C5-C6	5.42	121.25	118.00
3	B	200	THR	N-CA-C	5.41	125.59	111.00
1	C	14	DA	P-O3'-C3'	5.40	126.18	119.70
1	C	33	DT	C4-C5-C6	5.37	121.22	118.00
2	D	32	DT	C4-C5-C6	5.34	121.20	118.00
2	D	11	DT	C4-C5-C6	5.28	121.17	118.00
1	C	11	DT	C4-C5-C6	5.25	121.15	118.00
2	D	3	DT	C4-C5-C6	5.24	121.14	118.00
1	C	32	DT	C4-C5-C6	5.23	121.14	118.00
1	C	3	DT	C4-C5-C6	5.21	121.13	118.00
1	C	25	DT	C4-C5-C6	5.19	121.11	118.00
1	C	35	DT	C4-C5-C6	5.15	121.09	118.00
2	D	23	DT	C4-C5-C6	5.09	121.05	118.00
1	C	11	DT	C6-C5-C7	-5.09	119.85	122.90
2	D	13	DT	C4-C5-C6	5.04	121.03	118.00
2	D	18	DT	C4-C5-C6	5.01	121.00	118.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	B	198	GLY	Mainchain
3	B	199	ARG	Sidechain
3	B	201	LYS	Mainchain

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	715	0	401	29	0
2	D	713	0	400	32	0
3	A	2548	0	2571	88	1
3	B	2548	0	2565	142	1
4	A	154	0	0	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	191	0	0	5	0
4	C	60	0	0	1	0
4	D	88	0	0	5	0
All	All	7017	0	5937	253	1

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 20.

All (253) 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)
3:B:202:THR:HG1	3:B:203:LEU:N	0.96	1.43
2:D:24:DA:C5'	3:B:201:LYS:HG2	1.56	1.33
1:C:17:DG:H5"	3:B:202:THR:CB	1.62	1.30
3:B:202:THR:C	3:B:203:LEU:N	1.86	1.28
1:C:17:DG:C5'	3:B:202:THR:HB	1.66	1.26
2:D:24:DA:C5'	3:B:201:LYS:CG	2.22	1.17
3:B:201:LYS:CB	3:B:202:THR:HG23	1.77	1.14
3:B:202:THR:OG1	3:B:203:LEU:N	1.78	1.13
3:B:323:ASN:O	3:B:326:ARG:HB3	1.48	1.12
3:B:202:THR:O	3:B:203:LEU:HA	1.49	1.11
3:B:202:THR:C	3:B:203:LEU:CA	2.19	1.10
1:C:19:DA:H1'	1:C:20:DT:H5"	1.34	1.09
1:C:17:DG:H5"	3:B:202:THR:HB	1.14	1.09
3:B:325:ILE:HG23	3:B:328:LEU:HD23	1.34	1.08
3:B:201:LYS:HB2	3:B:202:THR:HG23	1.13	1.08
1:C:17:DG:C5'	3:B:202:THR:CB	2.25	1.07
1:C:17:DG:H5'	3:B:202:THR:CG2	1.84	1.06
1:C:17:DG:H5'	3:B:202:THR:HG21	1.33	1.04
2:D:24:DA:H5"	3:B:201:LYS:CG	1.88	1.02
2:D:24:DA:H5'	3:B:201:LYS:HG2	1.04	1.01
2:D:24:DA:H5'	3:B:201:LYS:CG	1.86	0.99
3:B:200:THR:HG1	3:B:203:LEU:N	1.60	0.98
2:D:5:DA:H2"	2:D:6:DC:H5"	1.47	0.96
3:B:201:LYS:CB	3:B:202:THR:CG2	2.43	0.96
3:B:199:ARG:HB2	3:B:204:VAL:HA	1.49	0.95
3:B:201:LYS:HB2	3:B:202:THR:CG2	1.97	0.94
2:D:24:DA:H5"	3:B:201:LYS:HG2	1.48	0.91
3:B:202:THR:CA	3:B:203:LEU:N	2.34	0.91
3:B:199:ARG:CA	3:B:203:LEU:O	2.19	0.90
1:C:19:DA:C1'	1:C:20:DT:H5"	2.02	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:24:DA:C4'	3:B:201:LYS:HD2	2.03	0.88
4:D:49:HOH:O	3:A:241:ARG:HD3	1.75	0.87
3:B:133:GLN:OE1	3:B:327:ASN:HB2	1.78	0.84
1:C:17:DG:C5'	3:B:202:THR:CG2	2.52	0.83
2:D:24:DA:H5"	3:B:201:LYS:HG3	1.59	0.83
3:B:320:ILE:O	3:B:323:ASN:HB2	1.79	0.83
2:D:6:DC:H2'	2:D:7:DT:H72	1.62	0.82
3:B:202:THR:C	3:B:203:LEU:HA	1.94	0.81
1:C:19:DA:H1'	1:C:20:DT:C5'	2.08	0.81
2:D:21:DA:H2"	2:D:22:DT:H5"	1.61	0.81
1:C:17:DG:C4'	3:B:202:THR:HB	2.11	0.81
3:A:194:LEU:HG	3:A:210:GLU:OE2	1.82	0.80
3:B:323:ASN:HA	3:B:326:ARG:HB2	1.63	0.80
3:B:306:ILE:HG21	3:B:318:VAL:HG13	1.62	0.79
1:C:5:DA:H2"	1:C:6:DC:H5"	1.66	0.78
3:A:174:ILE:HD12	3:A:258:THR:HB	1.64	0.78
3:A:199:ARG:HD2	3:A:209:VAL:HG21	1.64	0.77
1:C:17:DG:H5"	3:B:202:THR:OG1	1.85	0.77
3:B:199:ARG:CB	3:B:203:LEU:O	2.33	0.76
3:B:200:THR:OG1	3:B:203:LEU:N	2.18	0.76
3:B:199:ARG:HA	3:B:203:LEU:O	1.84	0.75
3:A:317:ASN:ND2	4:A:409:HOH:O	2.18	0.74
3:B:202:THR:CB	3:B:203:LEU:N	2.50	0.74
3:B:325:ILE:HG23	3:B:328:LEU:CD2	2.15	0.74
3:B:323:ASN:O	3:B:326:ARG:CB	2.34	0.74
4:C:73:HOH:O	3:B:156:GLN:HG3	1.86	0.73
3:B:202:THR:O	3:B:203:LEU:CA	2.25	0.73
1:C:17:DG:H2"	1:C:18:DT:H5'	1.71	0.73
3:B:199:ARG:HB2	3:B:203:LEU:O	1.90	0.72
2:D:15:DG:H2"	2:D:16:DC:H5"	1.69	0.71
1:C:17:DG:C5'	3:B:202:THR:HG21	2.13	0.71
3:B:197:ILE:CD1	3:B:211:LYS:HG3	2.22	0.70
3:B:206:THR:HG1	3:B:207:ALA:N	1.90	0.69
2:D:34:DA:H2"	2:D:35:DT:O4'	1.90	0.69
3:B:336:VAL:O	3:B:340:GLU:HG3	1.92	0.69
3:A:193:MET:CE	3:A:218:THR:HG23	2.23	0.69
3:B:146:ARG:O	3:B:150:GLU:HB2	1.92	0.69
1:C:17:DG:H4'	3:B:202:THR:HB	1.74	0.68
2:D:18:DT:H6	2:D:18:DT:H5"	1.59	0.68
1:C:6:DC:H2'	1:C:7:DT:H72	1.74	0.68
3:A:310:MET:SD	3:A:318:VAL:HG13	2.35	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:460:HOH:O	3:B:326:ARG:HD2	1.94	0.66
2:D:19:DA:H5"	4:D:99:HOH:O	1.96	0.65
3:B:101:ARG:HH11	3:B:101:ARG:HG2	1.62	0.65
2:D:16:DC:H2"	2:D:17:DA:C8	2.33	0.64
2:D:6:DC:H2'	2:D:7:DT:C7	2.27	0.64
3:A:210:GLU:O	3:B:331:GLU:HB2	1.97	0.64
3:A:30:MET:HE1	3:A:101:ARG:HB3	1.79	0.64
3:A:193:MET:HE3	3:A:218:THR:HG23	1.79	0.64
3:B:106:ARG:O	3:B:109:ASP:HB2	1.98	0.64
3:B:313:GLY:HA3	3:B:315:TRP:CZ3	2.33	0.63
1:C:19:DA:C2'	1:C:20:DT:H5"	2.28	0.63
3:A:185:ILE:HD11	3:A:238:LEU:HG	1.81	0.63
3:A:318:VAL:HG12	3:A:321:VAL:HG11	1.81	0.63
3:B:319:ASN:O	3:B:323:ASN:ND2	2.32	0.63
3:B:205:SER:HG	3:B:206:THR:N	1.94	0.63
2:D:24:DA:O4'	3:B:201:LYS:HD2	1.98	0.63
3:A:158:ILE:HD13	3:A:227:VAL:HG21	1.81	0.62
3:A:188:THR:HB	3:A:194:LEU:HD22	1.80	0.62
3:B:205:SER:OG	3:B:206:THR:N	2.29	0.61
3:B:214:SER:O	3:B:218:THR:HG23	1.99	0.61
3:B:201:LYS:CA	3:B:202:THR:HG23	2.26	0.61
3:B:197:ILE:HD13	3:B:211:LYS:HG3	1.80	0.61
2:D:24:DA:C5'	3:B:201:LYS:CD	2.79	0.61
4:D:91:HOH:O	3:B:243:ARG:HG2	2.00	0.60
3:B:129:GLU:O	3:B:130:ARG:HD3	2.01	0.60
3:A:262:GLU:O	3:A:266:GLU:HG3	2.02	0.60
2:D:5:DA:C2'	2:D:6:DC:H5"	2.27	0.59
3:B:187:ARG:NH2	3:B:222:GLU:OE2	2.36	0.59
3:B:159:ARG:HB2	3:B:224:TRP:CZ3	2.37	0.59
2:D:24:DA:H4'	3:B:201:LYS:HD2	1.84	0.59
3:B:201:LYS:CB	3:B:202:THR:HG22	2.32	0.59
2:D:24:DA:C4'	3:B:201:LYS:CD	2.79	0.58
1:C:16:DT:OP2	3:B:320:ILE:HG21	2.03	0.58
3:A:308:GLU:OE1	3:B:332:THR:HB	2.03	0.58
3:B:62:LYS:HD3	3:B:64:PHE:O	2.04	0.58
3:A:306:ILE:N	3:A:307:PRO:HD2	2.19	0.58
1:C:6:DC:H2'	1:C:7:DT:C7	2.33	0.57
3:A:197:ILE:HG23	3:A:198:GLY:N	2.17	0.57
3:B:201:LYS:HB3	3:B:202:THR:CG2	2.34	0.57
3:A:81:ARG:NH2	4:A:359:HOH:O	2.30	0.57
3:A:30:MET:CE	3:A:101:ARG:HB3	2.35	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:8:DT:H1'	2:D:9:DC:H5'	1.88	0.56
3:B:42:TRP:O	3:B:46:LEU:HG	2.06	0.56
3:A:102:SER:HB2	3:A:104:LEU:HD12	1.86	0.56
3:B:183:LYS:HD2	4:B:520:HOH:O	2.07	0.55
1:C:15:DA:N3	3:B:201:LYS:NZ	2.55	0.55
3:A:154:ARG:HH21	3:A:156:GLN:HE21	1.54	0.55
3:B:105:PRO:HD2	4:B:436:HOH:O	2.05	0.55
3:A:154:ARG:NH2	3:A:156:GLN:HE21	2.04	0.55
3:A:182:VAL:HG23	3:A:236:ASN:O	2.06	0.55
3:A:100:ARG:NH1	3:A:106:ARG:HD3	2.21	0.55
3:B:85:VAL:O	3:B:89:GLN:HG3	2.07	0.55
3:A:185:ILE:HG22	3:A:193:MET:HG2	1.88	0.55
3:B:209:VAL:HG22	3:B:210:GLU:H	1.70	0.55
3:B:187:ARG:HD3	3:B:191:GLY:HA2	1.89	0.55
3:A:40:HIS:HE1	4:A:410:HOH:O	1.90	0.55
2:D:18:DT:H6	2:D:18:DT:C5'	2.19	0.55
3:B:306:ILE:N	3:B:307:PRO:HD2	2.22	0.54
3:A:175:ALA:O	3:A:179:ARG:HG3	2.07	0.54
3:B:101:ARG:HH11	3:B:101:ARG:CG	2.20	0.54
3:B:133:GLN:HE22	3:B:324:TYR:HA	1.73	0.54
3:A:189:ASP:C	3:A:189:ASP:OD1	2.46	0.54
3:A:200:THR:HG21	3:A:205:SER:HB2	1.90	0.54
2:D:24:DA:H5'	3:B:201:LYS:CD	2.37	0.54
3:A:340:GLU:O	3:A:341:ASP:HB3	2.07	0.54
3:A:223:ARG:HG3	3:A:223:ARG:O	2.07	0.54
3:B:255:GLN:HG3	4:B:503:HOH:O	2.08	0.54
3:B:201:LYS:HB3	3:B:202:THR:HG22	1.91	0.53
3:B:193:MET:CE	3:B:222:GLU:HG3	2.39	0.53
3:B:315:TRP:HZ2	3:B:324:TYR:CE1	2.27	0.53
3:A:139:ARG:NH1	3:B:339:LEU:HA	2.23	0.53
3:B:85:VAL:HG23	3:B:129:GLU:OE2	2.09	0.53
3:A:209:VAL:HG12	3:A:210:GLU:N	2.24	0.53
3:A:59:ASN:O	3:A:61:ARG:HD3	2.08	0.53
3:A:172:LEU:HD21	3:A:197:ILE:HG21	1.91	0.52
3:B:297:ARG:HG2	3:B:328:LEU:HD22	1.90	0.52
3:A:199:ARG:HD2	3:A:209:VAL:CG2	2.37	0.52
3:A:209:VAL:HG12	3:A:210:GLU:H	1.75	0.52
3:A:139:ARG:HH12	3:B:339:LEU:HA	1.74	0.52
3:A:192:ARG:HG3	3:A:192:ARG:NH1	2.25	0.52
3:A:305:SER:C	3:A:307:PRO:HD2	2.30	0.52
3:A:193:MET:HE2	3:A:218:THR:HG23	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:89:GLN:HG3	3:A:117:MET:CE	2.39	0.51
3:B:199:ARG:C	3:B:203:LEU:O	2.49	0.51
4:A:460:HOH:O	3:B:326:ARG:CD	2.57	0.51
4:D:49:HOH:O	3:A:241:ARG:CD	2.44	0.51
3:A:24:ARG:HH11	3:A:24:ARG:HG3	1.75	0.51
2:D:24:DA:C5'	3:B:201:LYS:HD2	2.39	0.51
3:B:202:THR:C	3:B:203:LEU:C	2.69	0.51
3:A:24:ARG:HG3	3:A:24:ARG:NH1	2.26	0.50
3:A:200:THR:O	3:B:130:ARG:NH2	2.44	0.50
3:A:146:ARG:O	3:A:150:GLU:HG3	2.12	0.50
3:A:221:VAL:HG12	3:A:225:ILE:HD11	1.94	0.50
3:A:74:TYR:O	3:A:77:TYR:HB3	2.11	0.50
1:C:19:DA:H2"	1:C:20:DT:C5'	2.40	0.50
3:B:214:SER:O	3:B:218:THR:CG2	2.60	0.49
2:D:24:DA:H4'	3:B:201:LYS:CD	2.42	0.49
3:B:313:GLY:HA3	3:B:315:TRP:CE3	2.47	0.49
3:B:50:ARG:NH2	4:B:529:HOH:O	2.45	0.49
3:B:106:ARG:NH1	3:B:109:ASP:OD2	2.46	0.49
3:A:100:ARG:NH1	3:A:100:ARG:HG3	2.28	0.49
3:B:134:ALA:HA	3:B:283:TYR:CD2	2.48	0.49
3:B:245:ASN:OD1	3:B:247:VAL:HG23	2.12	0.49
3:A:225:ILE:CG2	3:A:231:ALA:HB2	2.42	0.48
3:A:218:THR:O	3:A:222:GLU:HG3	2.14	0.48
3:B:183:LYS:HB3	3:B:234:PRO:HB2	1.95	0.48
3:A:313:GLY:HA3	3:A:315:TRP:CZ3	2.48	0.48
3:B:211:LYS:NZ	3:B:312:ALA:O	2.47	0.48
3:B:272:ILE:N	3:B:272:ILE:HD13	2.28	0.48
3:B:315:TRP:HD1	3:B:320:ILE:HD13	1.78	0.48
3:B:315:TRP:HZ2	3:B:324:TYR:HE1	1.59	0.48
3:B:209:VAL:HG22	3:B:210:GLU:N	2.29	0.47
3:A:89:GLN:HG3	3:A:117:MET:HE2	1.96	0.47
3:A:187:ARG:NH1	3:A:191:GLY:O	2.47	0.47
3:A:68:PRO:HB3	3:A:110:SER:CB	2.45	0.47
3:B:213:LEU:HB2	3:B:218:THR:HG22	1.95	0.47
1:C:19:DA:C2'	1:C:20:DT:C5'	2.92	0.47
3:A:107:PRO:O	3:A:110:SER:HB3	2.15	0.47
3:B:31:PHE:CE2	3:B:34:ARG:NH1	2.83	0.47
3:A:310:MET:HE2	4:A:496:HOH:O	2.13	0.47
3:A:100:ARG:HH11	3:A:100:ARG:HG3	1.79	0.47
3:A:159:ARG:HB2	3:A:224:TRP:CZ3	2.50	0.46
3:A:305:SER:HB2	3:A:307:PRO:HD2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:277:ASP:HB3	3:B:284:LEU:HD13	1.96	0.46
3:B:193:MET:HE3	3:B:222:GLU:HG3	1.97	0.46
2:D:3:DT:O2	3:A:244:LYS:NZ	2.49	0.46
1:C:16:DT:OP2	3:B:320:ILE:CG2	2.64	0.46
3:B:172:LEU:HD11	3:B:197:ILE:HD11	1.97	0.46
1:C:11:DT:OP2	3:B:50:ARG:NH1	2.49	0.45
3:B:119:ARG:O	3:B:123:GLU:HG3	2.16	0.45
3:A:326:ARG:HG2	3:A:327:ASN:N	2.32	0.45
3:A:195:ILE:HD11	3:A:213:LEU:HD11	1.98	0.45
2:D:15:DG:O6	3:A:86:LYS:NZ	2.50	0.45
4:D:38:HOH:O	3:A:156:GLN:HB2	2.16	0.45
3:B:170:THR:O	3:B:171:LEU:HB2	2.17	0.45
3:B:271:LEU:HD13	3:B:272:ILE:HD13	1.98	0.45
3:B:315:TRP:CZ2	3:B:324:TYR:CE1	3.05	0.45
3:B:101:ARG:CG	3:B:101:ARG:NH1	2.77	0.45
3:B:277:ASP:CB	3:B:284:LEU:HD13	2.47	0.45
3:A:183:LYS:HB3	3:A:234:PRO:HB3	1.97	0.45
2:D:8:DT:H2"	2:D:9:DC:H5'	1.97	0.44
1:C:17:DG:H2"	1:C:18:DT:C5'	2.45	0.44
3:B:183:LYS:NZ	3:B:235:ASN:ND2	2.65	0.44
4:A:376:HOH:O	3:B:338:LEU:CD1	2.65	0.44
3:A:340:GLU:O	3:A:341:ASP:CB	2.65	0.44
1:C:6:DC:H2'	1:C:7:DT:C5	2.52	0.44
3:B:277:ASP:CG	3:B:278:ASP:H	2.20	0.44
1:C:26:DA:C8	1:C:26:DA:H5'	2.52	0.44
3:A:65:PRO:HG3	3:A:104:LEU:HD22	1.99	0.43
3:A:192:ARG:CG	3:A:192:ARG:HH11	2.31	0.43
3:B:201:LYS:HB3	3:B:201:LYS:HE2	1.81	0.43
3:B:323:ASN:C	3:B:326:ARG:H	2.21	0.43
3:A:104:LEU:HB3	3:A:105:PRO:HD2	2.00	0.43
3:B:84:ALA:O	3:B:88:ILE:HG13	2.18	0.43
3:B:231:ALA:HB3	4:B:399:HOH:O	2.19	0.42
2:D:11:DT:H2"	2:D:12:DA:OP2	2.19	0.42
3:B:27:LEU:HA	3:B:27:LEU:HD12	1.82	0.42
3:B:193:MET:HB2	3:B:218:THR:HB	2.01	0.42
3:B:315:TRP:CZ2	3:B:324:TYR:HE1	2.36	0.42
3:B:325:ILE:HA	3:B:328:LEU:HB2	2.01	0.42
3:A:72:ARG:HD3	4:A:373:HOH:O	2.19	0.42
3:A:277:ASP:HB2	3:A:284:LEU:HD13	2.02	0.42
3:A:192:ARG:HG3	3:A:192:ARG:HH11	1.82	0.42
3:A:112:ALA:O	3:A:116:VAL:HG22	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:21:DA:C2'	2:D:22:DT:H5"	2.42	0.42
3:A:233:ASP:O	3:A:236:ASN:HB2	2.20	0.42
3:A:170:THR:O	3:A:211:LYS:HE3	2.19	0.42
3:A:100:ARG:HG3	3:A:106:ARG:HD3	2.02	0.42
1:C:10:DG:N7	3:B:43:LYS:NZ	2.49	0.42
3:A:186:SER:OG	3:A:194:LEU:HB3	2.20	0.41
3:B:192:ARG:HG2	3:B:213:LEU:O	2.19	0.41
3:A:182:VAL:HG11	3:A:231:ALA:HA	2.00	0.41
3:B:96:ASN:OD1	3:B:107:PRO:HD2	2.20	0.41
3:A:172:LEU:CD1	3:A:197:ILE:HG13	2.50	0.41
3:A:43:LYS:HD3	3:A:43:LYS:HA	1.94	0.41
3:A:100:ARG:HG3	3:A:106:ARG:CD	2.50	0.41
3:A:25:LYS:NZ	4:A:468:HOH:O	2.51	0.41
3:A:52:TRP:CD1	3:A:63:TRP:HE3	2.38	0.41
3:A:172:LEU:HD11	3:A:197:ILE:HG13	2.03	0.41
3:B:326:ARG:O	3:B:326:ARG:HG2	2.21	0.40
3:B:297:ARG:O	3:B:301:ARG:HG2	2.22	0.40
3:A:227:VAL:HG12	3:A:227:VAL:O	2.20	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:126:ASP:OD2	3:B:199:ARG:NH2[4_566]	2.17	0.03

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	320/324 (99%)	301 (94%)	16 (5%)	3 (1%)	21 13
3	B	320/324 (99%)	302 (94%)	14 (4%)	4 (1%)	15 7

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	640/648 (99%)	603 (94%)	30 (5%)	7 (1%)	17 9

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	281	GLN
3	B	200	THR
3	B	201	LYS
3	B	277	ASP
3	A	275	ALA
3	A	278	ASP
3	B	206	THR

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	A	269/270 (100%)	215 (80%)	54 (20%)	1 0
3	B	269/270 (100%)	224 (83%)	45 (17%)	3 1
All	All	538/540 (100%)	439 (82%)	99 (18%)	2 1

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	20	SER
3	A	27	LEU
3	A	39	GLU
3	A	61	ARG
3	A	62	LYS
3	A	67	GLU
3	A	86	LYS
3	A	90	GLN
3	A	95	LEU
3	A	98	LEU
3	A	100	ARG

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Mol	Chain	Res	Type
3	A	106	ARG
3	A	108	SER
3	A	114	SER
3	A	116	VAL
3	A	121	ARG
3	A	132	LYS
3	A	147	SER
3	A	148	LEU
3	A	151	ASN
3	A	161	LEU
3	A	169	ASN
3	A	171	LEU
3	A	181	ARG
3	A	183	LYS
3	A	185	ILE
3	A	189	ASP
3	A	192	ARG
3	A	197	ILE
3	A	199	ARG
3	A	200	THR
3	A	201	LYS
3	A	219	LYS
3	A	223	ARG
3	A	226	SER
3	A	238	LEU
3	A	241	ARG
3	A	270	ARG
3	A	271	LEU
3	A	276	LYS
3	A	277	ASP
3	A	279	SER
3	A	282	ARG
3	A	289	HIS
3	A	292	ARG
3	A	299	MET
3	A	306	ILE
3	A	318	VAL
3	A	320	ILE
3	A	321	VAL
3	A	322	MET
3	A	326	ARG
3	A	335	MET

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Mol	Chain	Res	Type
3	A	338	LEU
3	B	22	GLU
3	B	27	LEU
3	B	28	MET
3	B	30	MET
3	B	34	ARG
3	B	38	SER
3	B	39	GLU
3	B	45	LEU
3	B	61	ARG
3	B	62	LYS
3	B	90	GLN
3	B	95	LEU
3	B	101	ARG
3	B	108	SER
3	B	118	ARG
3	B	122	LYS
3	B	130	ARG
3	B	156	GLN
3	B	169	ASN
3	B	173	LYS
3	B	183	LYS
3	B	186	SER
3	B	199	ARG
3	B	200	THR
3	B	205	SER
3	B	211	LYS
3	B	218	THR
3	B	219	LYS
3	B	226	SER
3	B	235	ASN
3	B	243	ARG
3	B	255	GLN
3	B	266	GLU
3	B	271	LEU
3	B	276	LYS
3	B	282	ARG
3	B	289	HIS
3	B	293	VAL
3	B	316	THR
3	B	320	ILE
3	B	326	ARG

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Mol	Chain	Res	Type
3	B	328	LEU
3	B	330	SER
3	B	338	LEU
3	B	341	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	89	GLN
3	A	144	GLN
3	B	90	GLN
3	B	133	GLN
3	B	235	ASN
3	B	311	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.