



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:55 PM GMT

PDB ID : 1VTO
Title : 1.9 Å RESOLUTION REFINED STRUCTURE OF TBP RECOGNIZING
THE MINOR GROOVE OF TATAAAAG
Authors : Kim, J.L.; Burley, S.K.
Deposited on : 1996-09-06
Resolution : 1.90 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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

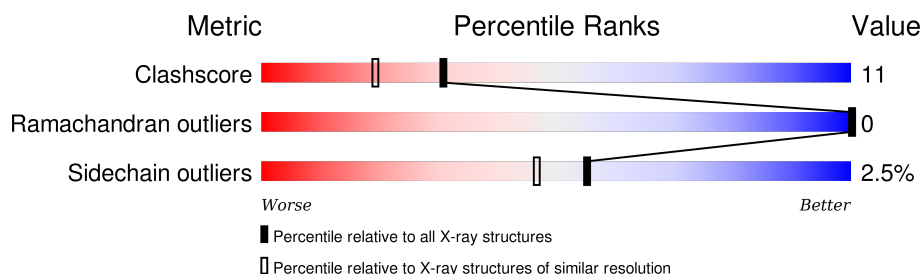
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)

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 $\leq 5\%$. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	C	14	
1	E	14	
2	D	14	
2	F	14	
3	A	190	
3	B	190	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6447 atoms, of which 1830 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 DNA (5'-D(*GP*CP*TP*AP*TP*AP*AP*AP*AP*GP*GP*GP*CP*A)-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	C	14	Total	C	H	N	O	P	0	0	0
			318	137	31	60	77	13			
1	E	14	Total	C	H	N	O	P	0	0	0
			321	138	32	60	78	13			

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

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	D	14	Total	C	H	N	O	P	0	0	0
			305	136	26	44	86	13			
2	F	14	Total	C	H	N	O	P	0	0	0
			305	136	26	44	86	13			

- Molecule 3 is a protein called TATA BINDING PROTEIN.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	A	187	Total	C	H	N	O	S	0	0	0
			1806	960	333	250	255	8			
3	B	188	Total	C	H	N	O	S	0	1	0
			1831	970	342	255	256	8			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	172	Total	H	O	0	0
			516	344	172		
4	B	185	Total	H	O	0	0
			553	368	185		
4	C	38	Total	H	O	0	0
			114	76	38		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	39	Total	H	O	0	0
			117	78	39		
4	E	37	Total	H	O	0	0
			111	74	37		
4	F	50	Total	H	O	0	0
			150	100	50		

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.

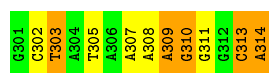
- Molecule 1: DNA (5'-D(*GP*CP*TP*AP*TP*AP*AP*AP*AP*GP*GP*GP*CP*A)-3')

Chain C: 



- Molecule 1: DNA (5'-D(*GP*CP*TP*AP*TP*AP*AP*AP*AP*GP*GP*GP*CP*A)-3')

Chain E: 



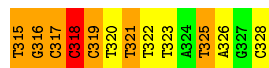
- Molecule 2: DNA (5'-D(*TP*GP*CP*CP*CP*TP*TP*TP*TP*AP*TP*AP*GP*C)-3')

Chain D: 




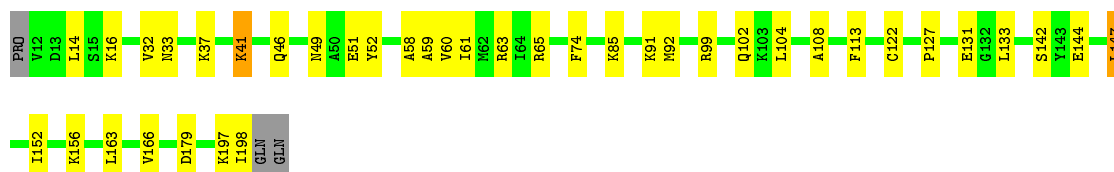
- Molecule 2: DNA (5'-D(*TP*GP*CP*CP*CP*TP*TP*TP*TP*AP*TP*AP*GP*C)-3')

Chain F: 

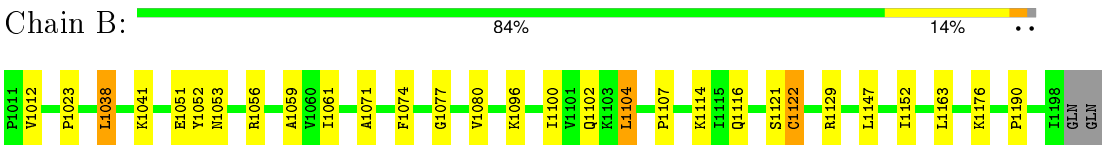


- Molecule 3: TATA BINDING PROTEIN

Chain A: 



● Molecule 3: TATA BINDING PROTEIN



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	42.00 Å 147.00 Å 57.00 Å 90.00° 90.50° 90.00°	Depositor
Resolution (Å)	6.00 – 1.90	Depositor
% Data completeness (in resolution range)	94.4 (6.00-1.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	5.20	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.194 , 0.264	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6447	wwPDB-VP
Average B, all atoms (Å ²)	20.0	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	1.37	0/324	1.68	7/499 (1.4%)
1	E	1.43	0/326	1.82	13/502 (2.6%)
2	D	1.46	2/310 (0.6%)	2.02	15/476 (3.2%)
2	F	1.53	5/310 (1.6%)	2.42	22/476 (4.6%)
3	A	0.65	1/1503 (0.1%)	0.77	0/2024
3	B	0.68	1/1525 (0.1%)	0.79	0/2052
All	All	0.96	9/4298 (0.2%)	1.31	57/6029 (0.9%)

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
2	D	0	1
2	F	0	1
All	All	0	3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1122	CYS	CB-SG	-8.31	1.68	1.82
2	D	215	DT	C5-C7	6.84	1.54	1.50
2	F	322	DT	C5-C7	6.69	1.54	1.50
2	F	315	DT	C5-C7	5.93	1.53	1.50
2	D	221	DT	C5-C7	5.75	1.53	1.50
2	F	325	DT	C5-C7	5.34	1.53	1.50
2	F	328	DC	C4'-O4'	-5.24	1.39	1.45
3	A	122	CYS	CB-SG	-5.11	1.73	1.81
2	F	318	DC	P-O5'	5.00	1.64	1.59

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	317	DC	O4'-C4'-C3'	-13.58	97.85	106.00
2	D	227	DG	O4'-C1'-N9	13.48	117.44	108.00
2	F	320	DT	C6-C5-C7	-10.79	116.43	122.90
2	F	317	DC	O4'-C1'-N1	10.71	115.49	108.00
2	F	316	DG	O4'-C1'-N9	8.91	114.24	108.00
2	F	319	DC	O4'-C4'-C3'	-8.83	100.70	106.00
2	F	321	DT	O4'-C4'-C3'	-8.66	100.81	106.00
2	F	328	DC	C1'-O4'-C4'	-8.24	101.86	110.10
2	F	323	DT	C6-C5-C7	-8.21	117.97	122.90
1	C	202	DC	O4'-C1'-N1	8.21	113.75	108.00
2	F	328	DC	O4'-C4'-C3'	-8.19	101.09	106.00
2	F	321	DT	C6-C5-C7	-7.78	118.23	122.90
2	F	322	DT	C6-C5-C7	-7.66	118.31	122.90
2	F	315	DT	P-O3'-C3'	7.60	128.81	119.70
1	E	310	DG	O4'-C1'-N9	7.59	113.31	108.00
1	E	313	DC	C1'-O4'-C4'	-7.54	102.56	110.10
2	D	223	DT	O4'-C4'-C3'	-7.44	101.53	104.50
2	F	317	DC	C1'-O4'-C4'	-7.17	102.93	110.10
1	E	314	DA	O4'-C1'-N9	6.88	112.82	108.00
2	D	223	DT	C6-C5-C7	-6.65	118.91	122.90
2	D	222	DT	C6-C5-C7	-6.64	118.92	122.90
2	D	216	DG	O4'-C1'-C2'	-6.61	100.62	105.90
2	F	325	DT	C6-C5-C7	-6.59	118.94	122.90
2	F	318	DC	O4'-C4'-C3'	-6.53	101.89	104.50
1	C	201	DG	O4'-C1'-N9	6.49	112.55	108.00
2	F	315	DT	C6-C5-C7	-6.49	119.00	122.90
1	E	313	DC	O4'-C1'-N1	6.42	112.49	108.00
1	E	313	DC	C6-N1-C2	6.25	122.80	120.30
1	C	207	DA	O4'-C1'-N9	6.25	112.37	108.00
1	E	303	DT	C6-C5-C7	-6.22	119.17	122.90
2	D	219	DC	C1'-O4'-C4'	-6.16	103.94	110.10
1	E	313	DC	O4'-C1'-C2'	-6.15	100.98	105.90
2	F	320	DT	C4-C5-C7	6.06	122.64	119.00
1	C	203	DT	C4-C5-C6	6.06	121.64	118.00
2	D	227	DG	C4'-C3'-C2'	-6.00	97.70	103.10
2	F	319	DC	C1'-O4'-C4'	-5.97	104.13	110.10
1	C	203	DT	C6-C5-C7	-5.84	119.40	122.90
1	E	302	DC	O4'-C1'-N1	5.83	112.08	108.00
2	D	220	DT	C6-C5-C7	-5.81	119.41	122.90
2	D	222	DT	C4-C5-C6	5.72	121.44	118.00
1	C	205	DT	C6-C5-C7	-5.72	119.47	122.90
2	D	225	DT	C4-C5-C6	5.67	121.40	118.00
1	E	305	DT	C6-C5-C7	-5.66	119.51	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	225	DT	C6-C5-C7	-5.66	119.51	122.90
1	C	213	DC	O4'-C1'-C2'	-5.64	101.38	105.90
2	F	315	DT	O4'-C1'-C2'	-5.57	101.44	105.90
2	D	221	DT	C6-C5-C7	-5.54	119.57	122.90
1	E	310	DG	N3-C4-N9	-5.50	122.70	126.00
1	E	310	DG	N9-C4-C5	5.48	107.59	105.40
2	F	326	DA	O4'-C1'-N9	5.47	111.83	108.00
2	D	220	DT	C4-C5-C6	5.42	121.25	118.00
2	F	325	DT	O4'-C4'-C3'	-5.41	102.34	104.50
2	D	223	DT	C4-C5-C6	5.29	121.18	118.00
1	E	309	DA	O4'-C1'-N9	5.22	111.65	108.00
1	E	310	DG	C8-N9-C1'	5.18	133.73	127.00
2	F	315	DT	C4-C5-C6	5.17	121.10	118.00
2	D	221	DT	C4-C5-C6	5.07	121.04	118.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	207	DA	Sidechain
2	D	227	DG	Sidechain
2	F	318	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	287	31	154	4	0
1	E	289	32	158	5	5
2	D	279	26	162	8	0
2	F	279	26	162	13	0
3	A	1473	333	1549	31	5
3	B	1489	342	1575	26	0
4	A	172	344	0	8	5
4	B	185	368	0	8	4
4	C	38	76	0	1	0
4	D	39	78	0	1	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	E	37	74	0	2	1
4	F	50	100	0	0	1
All	All	4617	1830	3760	81	11

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 11.

All (81) 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:A:156:LYS:HA	3:A:156:LYS:HE2	1.43	0.97
3:B:1041:LYS:HD3	3:B:1052:TYR:HE2	1.33	0.89
2:F:317:DC:H2'	2:F:318:DC:C6	2.09	0.87
4:C:924:HOH:O	3:A:58:ALA:HB3	1.75	0.87
3:A:198:ILE:HD11	3:B:1190:PRO:HG2	1.56	0.86
2:F:315:DT:H2''	2:F:316:DG:H2'	1.64	0.78
3:B:1041:LYS:HD3	3:B:1052:TYR:CE2	2.17	0.78
3:B:1041:LYS:HG2	4:B:759:HOH:O	1.83	0.78
3:A:131:GLU:HB2	4:A:819:HOH:O	1.85	0.77
2:D:219:DC:H2'	2:D:220:DT:H71	1.68	0.76
2:D:217:DC:H5''	4:D:844:HOH:O	1.88	0.73
3:B:1096:LYS:HG3	4:B:919:HOH:O	1.89	0.72
3:A:41:LYS:HE3	3:A:41:LYS:O	1.95	0.67
3:A:102:GLN:HG3	3:A:108:ALA:HB3	1.74	0.67
2:F:315:DT:H2''	2:F:316:DG:C8	2.30	0.65
3:A:152:ILE:HD12	3:A:163:LEU:HD22	1.79	0.65
2:F:319:DC:OP2	2:F:319:DC:H6	1.79	0.64
1:C:202:DC:H2''	1:C:203:DT:H71	1.80	0.64
2:F:315:DT:H1'	2:F:316:DG:H5'	1.81	0.63
2:D:218:DC:H2''	2:D:219:DC:OP2	1.98	0.63
2:F:318:DC:H2'	2:F:319:DC:O4'	2.01	0.61
3:A:92:MET:HG3	4:A:774:HOH:O	2.00	0.61
3:A:156:LYS:CE	3:A:156:LYS:HA	2.23	0.60
3:A:16:LYS:O	3:A:16:LYS:HD3	2.02	0.60
3:B:1121:SER:HB2	4:B:880:HOH:O	2.01	0.60
3:A:51:GLU:HG3	3:A:61:ILE:HB	1.84	0.59
2:F:317:DC:H2'	2:F:318:DC:H6	1.66	0.58
3:B:1041:LYS:CD	3:B:1052:TYR:HE2	2.13	0.55
3:A:198:ILE:HD11	3:B:1190:PRO:CG	2.31	0.55
3:A:63:ARG:HD2	4:A:555:HOH:O	2.07	0.54
2:D:218:DC:H2''	2:D:219:DC:C6	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:52:TYR:HB2	3:A:60:VAL:HG22	1.91	0.52
2:F:325:DT:H5'	4:B:880:HOH:O	2.09	0.52
3:A:59:ALA:HB2	3:A:74:PHE:CE2	2.45	0.52
3:B:1116:GLN:O	3:B:1176:LYS:HE2	2.09	0.52
2:F:315:DT:H4'	2:F:316:DG:OP1	2.09	0.51
3:A:152:ILE:CD1	3:A:163:LEU:HD22	2.41	0.50
3:B:1053:ASN:HB2	3:B:1056:ARG:HB3	1.94	0.50
3:A:197:LYS:NZ	4:A:917:HOH:O	2.45	0.48
3:A:85:LYS:HE2	4:A:458:HOH:O	2.12	0.48
3:A:144:GLU:O	3:A:147:LEU:HB2	2.13	0.48
2:D:227:DG:H2''	2:D:228:DC:C5	2.49	0.47
3:A:63:ARG:CD	4:A:555:HOH:O	2.61	0.47
1:E:303:DT:H2''	3:B:1147:LEU:O	2.15	0.47
1:E:307:DA:OP2	4:E:631:HOH:O	2.20	0.47
3:A:16:LYS:HD3	4:A:911:HOH:O	2.15	0.46
2:F:315:DT:H2''	2:F:316:DG:H8	1.80	0.46
3:A:51:GLU:CG	3:A:61:ILE:HB	2.45	0.46
3:A:37:LYS:HB3	4:A:677:HOH:O	2.16	0.45
1:E:313:DC:H2''	1:E:314:DA:C8	2.51	0.45
3:B:1152:ILE:HD12	3:B:1163:LEU:HD22	1.99	0.45
1:E:310:DG:H2''	4:E:662:HOH:O	2.16	0.45
3:B:1051:GLU:HB2	3:B:1061:ILE:HB	1.99	0.44
3:B:1038:LEU:HD22	3:B:1077:GLY:HA2	1.99	0.44
2:F:315:DT:C1'	2:F:316:DG:H5'	2.45	0.44
2:D:218:DC:H2''	2:D:219:DC:H6	1.83	0.44
3:B:1012:VAL:HG22	4:B:618:HOH:O	2.17	0.44
3:B:1114:LYS:HG2	3:B:1116:GLN:NE2	2.33	0.44
3:B:1107:PRO:HD3	4:B:609:HOH:O	2.17	0.43
3:B:1053:ASN:O	3:B:1056:ARG:HB3	2.17	0.43
3:B:1023:PRO:HA	3:B:1122:CYS:HB3	2.01	0.43
3:A:99:ARG:HD3	3:A:102:GLN:OE1	2.18	0.43
2:F:321:DT:H5'	3:B:1061:ILE:HD13	2.00	0.43
3:B:1129:ARG:HD3	4:B:816:HOH:O	2.18	0.43
3:B:1059:ALA:HB2	3:B:1074:PHE:CE2	2.54	0.42
3:A:127:PRO:HA	3:A:166:VAL:O	2.19	0.42
3:A:37:LYS:HA	3:A:37:LYS:HD3	1.77	0.42
2:F:315:DT:C2'	2:F:316:DG:C8	3.01	0.42
1:C:202:DC:H2''	1:C:203:DT:C7	2.46	0.42
3:A:91:LYS:HA	3:A:113:PHE:CE1	2.55	0.42
1:C:202:DC:N3	2:D:227:DG:O6	2.52	0.42
3:B:1100:ILE:O	3:B:1104:LEU:HD22	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:1056:ARG:HD3	4:B:596:HOH:O	2.20	0.41
3:A:152:ILE:HD12	3:A:163:LEU:CD2	2.49	0.41
3:A:32:VAL:HG22	3:A:33:ASN:N	2.35	0.41
3:B:1071:ALA:HA	3:B:1080:VAL:O	2.21	0.41
3:A:142:SER:HB2	3:A:152:ILE:HB	2.02	0.40
1:C:201:DG:N3	1:C:201:DG:H2'	2.36	0.40
3:A:104:LEU:HD12	3:A:104:LEU:HA	1.94	0.40
2:D:228:DC:H6	2:D:228:DC:O5'	2.05	0.40
1:E:309:DA:H1'	3:B:1074:PHE:CE1	2.56	0.40

All (11) 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)
4:A:693:HOH:H1	4:B:613:HOH:H1[2_647]	1.18	0.42
4:A:693:HOH:O	4:B:613:HOH:H1[2_647]	1.19	0.41
1:E:311:DG:N7	3:A:65:ARG:HH22[2_657]	1.21	0.39
3:A:46:GLN:HE21	4:B:885:HOH:O[2_647]	1.31	0.29
4:A:693:HOH:O	4:B:613:HOH:O[2_647]	2.03	0.17
1:E:311:DG:O6	3:A:65:ARG:HH12[2_657]	1.49	0.11
1:E:308:DA:OP1	3:A:49:ASN:ND2[2_657]	2.09	0.11
4:E:662:HOH:H1	4:A:601:HOH:O[2_657]	1.50	0.10
1:E:311:DG:N7	3:A:65:ARG:NH2[2_657]	2.13	0.07
1:E:309:DA:OP2	4:A:824:HOH:O[2_657]	2.14	0.06
4:D:622:HOH:O	4:F:674:HOH:H2[2_546]	1.55	0.05

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	185/190 (97%)	177 (96%)	8 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	B	187/190 (98%)	179 (96%)	8 (4%)	0	100	100
All	All	372/380 (98%)	356 (96%)	16 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	159/163 (98%)	154 (97%)	5 (3%)	47	37
3	B	162/163 (99%)	159 (98%)	3 (2%)	65	59
All	All	321/326 (98%)	313 (98%)	8 (2%)	55	47

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	14	LEU
3	A	41	LYS
3	A	133	LEU
3	A	147	LEU
3	A	179	ASP
3	B	1038	LEU
3	B	1102	GLN
3	B	1104	LEU

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

Mol	Chain	Res	Type
3	A	49	ASN
3	B	1053	ASN
3	B	1137	HIS

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 carbohydrates 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 ⓘ

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.