Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 07:55 AM EDT

PDB ID	:	1EOO
Title	:	ECORV BOUND TO COGNATE DNA
Authors	:	Horton, N.C.; Perona, J.J.
Deposited on	:	2000-03-23
Resolution	:	2.16 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:

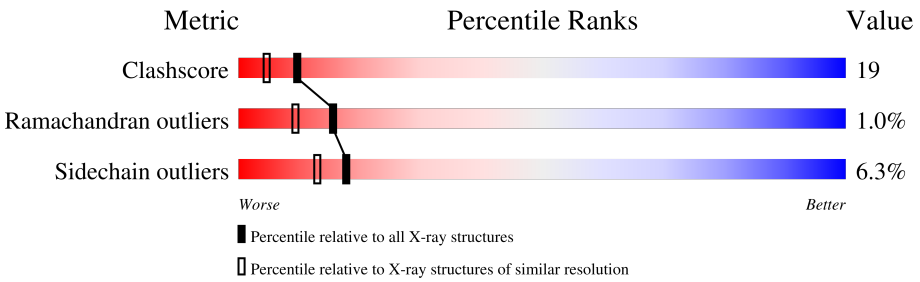
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 2.16 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	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)

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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	C	12	67% 33%
1	D	12	8% 50% 33% 8%
2	A	245	64% 32% •
2	B	245	62% 35% •

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4667 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 DNA (5'-D(*GP*AP*AP*GP*AP*TP*AP*TP*CP*TP*TP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	C	12	243	118	44	70	11	0	0	0
1	D	12	243	118	44	70	11	0	0	0

- Molecule 2 is a protein called TYPE II RESTRICTION ENZYME ECORV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	A	244	1947	1250	328	368	1	0	0	0
2	B	244	1941	1249	325	366	1	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	C	28	28	28	0	0
3	D	40	40	40	0	0
3	A	113	113	113	0	0
3	B	112	112	112	0	0

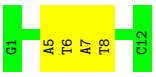
3 Residue-property plots (i)

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. 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. 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.

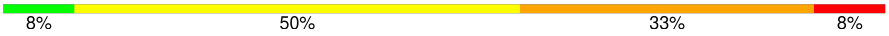
• Molecule 1: DNA (5'-D(*GP*AP*AP*GP*AP*TP*AP*TP*CP*TP*TP*C)-3')

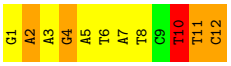
Chain C:



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

Chain D:



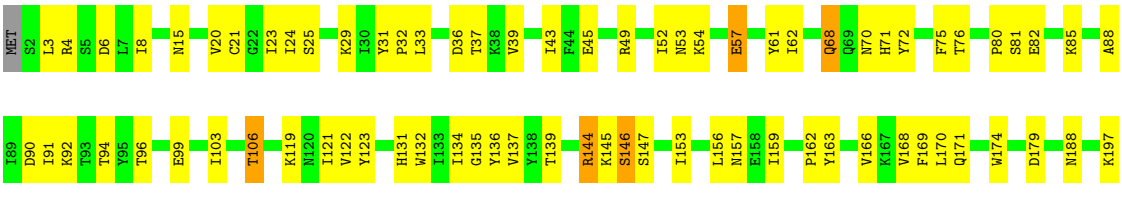


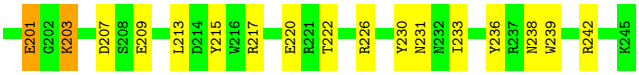
• Molecule 2: TYPE II RESTRICTION ENZYME ECORV

Chain A:

• Molecule 2: TYPE II RESTRICTION ENZYME ECORV

Chain B:




E201
G202
K203
D207
S208
E209
L213
D214
Y215
W216
R217
E220
R221
T222
R226
Y230
N231
N232
I233
Y236
R237
N238
W239
R242
K245

4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	72.60Å 120.40Å 182.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	4.80 – 2.16	Depositor
% Data completeness (in resolution range)	80.3 (4.80-2.16)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.231 , 0.316	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4667	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

5 Model quality (i)

5.1 Standard geometry (i)

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	0.98	0/272	1.08	0/418
1	D	1.10	0/272	1.43	3/418 (0.7%)
2	A	0.61	0/1995	0.80	1/2710 (0.0%)
2	B	0.63	0/1989	0.84	1/2705 (0.0%)
All	All	0.68	0/4528	0.89	5/6251 (0.1%)

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	D	0	5

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	4	DG	O4'-C4'-C3'	-6.99	101.70	104.50
1	D	4	DG	O4'-C1'-N9	6.45	112.51	108.00
1	D	10	DT	O4'-C1'-C2'	6.20	110.86	105.90
2	A	106	THR	N-CA-C	-5.36	96.53	111.00
2	B	106	THR	N-CA-C	-5.01	97.48	111.00

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	1	DG	Sidechain
1	D	10	DT	Sidechain
1	D	11	DT	Sidechain
1	D	12	DC	Sidechain

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Mol	Chain	Res	Type	Group
1	D	2	DA	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	243	0	138	6	0
1	D	243	0	138	15	0
2	A	1947	0	1813	81	0
2	B	1941	0	1817	79	0
3	A	113	0	0	8	0
3	B	112	0	0	6	0
3	C	28	0	0	0	0
3	D	40	0	0	0	0
All	All	4667	0	3906	161	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 (161) 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:B:68:GLN:HG2	3:B:339:HOH:O	1.69	0.91
2:A:49:ARG:HG2	2:A:75:PHE:HZ	1.42	0.85
2:A:201:GLU:HB3	2:A:203:LYS:NZ	1.93	0.84
2:A:201:GLU:HB3	2:A:203:LYS:HZ3	1.52	0.74
2:B:103:ILE:HD12	2:B:137:VAL:HG21	1.69	0.74
2:A:14:GLU:HG2	2:A:51:ILE:HD11	1.68	0.73
1:D:3:DA:H2"	1:D:4:DG:O5'	1.91	0.71
2:A:8:ILE:CG2	2:A:170:LEU:HB3	2.23	0.68
1:D:11:DT:H2"	1:D:12:DC:O5'	1.95	0.67
2:B:15:ASN:OD1	2:B:136:TYR:HE1	1.80	0.65
2:B:8:ILE:HG12	2:B:170:LEU:HB3	1.79	0.65
2:B:49:ARG:HG2	2:B:75:PHE:HZ	1.61	0.65
2:B:68:GLN:HB2	2:B:71:HIS:CD2	2.33	0.64
2:A:193:HIS:HE1	3:A:295:HOH:O	1.81	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:206:PHE:CE1	2:A:233:ILE:HD13	2.33	0.64
2:B:231:ASN:ND2	3:B:279:HOH:O	2.31	0.63
2:A:153:ILE:HG23	2:B:153:ILE:HD13	1.79	0.63
1:D:4:DG:H2'	1:D:5:DA:C8	2.34	0.63
2:B:197:LYS:O	2:B:201:GLU:HB2	2.00	0.62
2:B:238:ASN:O	2:B:242:ARG:HG3	1.99	0.62
2:A:57:GLU:HG2	2:A:58:LYS:N	2.14	0.62
2:B:29:LYS:HD3	2:B:31:TYR:OH	1.99	0.61
2:B:4:ARG:O	2:B:8:ILE:HG13	1.99	0.61
1:D:10:DT:H2"	1:D:11:DT:O5'	2.00	0.60
2:B:52:ILE:HD11	2:B:134:ILE:HD12	1.84	0.60
1:C:7:DA:H62	2:A:186:THR:HG21	1.67	0.60
1:D:4:DG:H2'	1:D:5:DA:H8	1.68	0.59
2:A:86:LYS:NZ	2:A:127:GLN:OE1	2.34	0.59
2:B:39:VAL:O	2:B:43:ILE:HG12	2.03	0.59
1:D:6:DT:O4'	2:B:70:ASN:HB3	2.03	0.58
2:B:94:THR:HG22	2:B:137:VAL:HG22	1.84	0.58
2:A:49:ARG:HB2	2:A:50:PRO:HD3	1.86	0.58
2:B:137:VAL:O	2:B:166:VAL:HA	2.04	0.58
2:B:201:GLU:HB3	2:B:203:LYS:HG2	1.86	0.58
2:A:23:ILE:HG12	2:A:31:TYR:HB2	1.85	0.57
2:A:153:ILE:HD13	2:B:153:ILE:CG2	2.35	0.57
2:B:213:LEU:O	2:B:217:ARG:HG3	2.04	0.57
2:A:85:LYS:HA	2:A:129:ILE:HG23	1.86	0.57
2:A:91:ILE:O	2:A:92:LYS:NZ	2.37	0.57
2:A:43:ILE:HD12	2:B:23:ILE:HG21	1.86	0.57
2:A:96:THR:HG22	2:A:139:THR:HA	1.87	0.57
2:A:193:HIS:CE1	3:A:295:HOH:O	2.58	0.56
2:A:201:GLU:HB3	2:A:203:LYS:HZ2	1.67	0.55
2:B:157:ASN:H	2:B:157:ASN:ND2	2.04	0.55
2:A:103:ILE:HG21	2:A:137:VAL:HG11	1.86	0.55
2:A:102:LYS:NZ	3:A:324:HOH:O	2.39	0.55
2:B:4:ARG:HB2	2:B:132:TRP:CH2	2.41	0.55
2:A:2:SER:O	2:A:6:ASP:HB2	2.06	0.55
2:A:69:GLN:HE22	2:B:37:THR:H	1.52	0.55
2:B:20:VAL:HG11	2:B:43:ILE:CG2	2.37	0.54
2:A:3:LEU:HD22	2:A:61:TYR:CE1	2.42	0.54
1:C:7:DA:H3'	2:A:92:LYS:HE3	1.90	0.54
2:A:18:TYR:CZ	2:B:29:LYS:HD2	2.43	0.54
2:A:116:ASN:HB3	2:A:119:LYS:HB2	1.90	0.54
1:D:2:DA:H5"	2:A:223:SER:HB3	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:5:DA:H2'	1:C:6:DT:H72	1.91	0.53
2:A:167:LYS:HG2	3:A:311:HOH:O	2.07	0.53
1:C:7:DA:N6	2:A:186:THR:HG21	2.24	0.53
1:D:4:DG:H2"	1:D:5:DA:O5'	2.09	0.53
2:A:59:HIS:O	2:A:79:LYS:NZ	2.42	0.53
2:A:82:GLU:HB3	2:A:85:LYS:HB2	1.92	0.53
2:B:21:CYS:SG	2:B:156:LEU:HD22	2.49	0.53
2:B:20:VAL:HG11	2:B:43:ILE:HG23	1.91	0.52
2:A:49:ARG:HG2	2:A:75:PHE:CZ	2.33	0.52
1:D:7:DA:H2"	1:D:8:DT:O5'	2.09	0.52
2:A:213:LEU:O	2:A:217:ARG:HG3	2.11	0.51
2:B:222:THR:O	2:B:226:ARG:HG3	2.10	0.51
1:C:5:DA:H2'	1:C:6:DT:C7	2.40	0.51
2:A:7:LEU:O	2:A:11:LEU:HG	2.11	0.51
2:A:203:LYS:HD2	2:A:203:LYS:N	2.25	0.51
2:B:68:GLN:NE2	2:B:71:HIS:HD2	2.08	0.51
2:B:236:TYR:O	2:B:239:TRP:HB3	2.11	0.51
2:A:69:GLN:NE2	2:B:37:THR:H	2.08	0.51
2:B:68:GLN:HB3	3:B:301:HOH:O	2.11	0.50
2:A:5:SER:HA	2:A:8:ILE:HD11	1.94	0.50
2:A:116:ASN:HB3	2:A:119:LYS:CB	2.41	0.50
2:B:96:THR:O	2:B:139:THR:HA	2.12	0.50
2:A:69:GLN:HE21	2:B:36:ASP:HB2	1.77	0.50
2:A:214:ASP:HB3	2:A:236:TYR:OH	2.12	0.49
2:B:94:THR:CG2	2:B:137:VAL:HG22	2.42	0.49
2:A:32:PRO:HG2	2:A:147:SER:O	2.12	0.49
2:A:53:ASN:O	2:A:57:GLU:HB3	2.13	0.48
2:B:122:VAL:HG21	3:B:306:HOH:O	2.12	0.48
1:D:2:DA:H2"	1:D:3:DA:C5'	2.43	0.48
2:B:144:ARG:C	2:B:146:SER:H	2.15	0.48
1:D:4:DG:C2'	1:D:5:DA:C8	2.96	0.48
2:B:76:THR:HG21	2:B:123:TYR:CD2	2.49	0.48
2:B:220:GLU:HB2	2:B:226:ARG:HG2	1.96	0.48
2:A:121:ILE:HG12	2:A:123:TYR:O	2.13	0.47
2:A:153:ILE:HD13	2:B:153:ILE:HG21	1.95	0.47
2:A:20:VAL:HA	2:B:25:SER:HA	1.96	0.47
2:A:8:ILE:HG22	2:A:170:LEU:HB3	1.96	0.47
2:A:104:LYS:HG2	3:A:294:HOH:O	2.14	0.47
1:D:4:DG:C2'	1:D:5:DA:H8	2.28	0.47
2:A:58:LYS:HG3	2:A:59:HIS:HD2	1.80	0.47
2:A:72:TYR:O	2:A:121:ILE:HA	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:179:ASP:OD1	2:B:179:ASP:N	2.48	0.46
2:A:69:GLN:NE2	2:B:36:ASP:HB2	2.30	0.46
2:B:174:TRP:HB3	2:B:209:GLU:OE1	2.16	0.46
2:B:54:LYS:HE2	2:B:54:LYS:HB3	1.72	0.45
2:A:183:SER:OG	2:A:188:ASN:HB2	2.16	0.45
2:B:163:TYR:CD1	2:B:166:VAL:HG23	2.51	0.45
2:B:226:ARG:HE	2:B:231:ASN:ND2	2.15	0.45
2:B:32:PRO:HB3	2:B:159:ILE:HG23	1.99	0.45
2:A:192:ILE:HG13	2:A:199:PHE:CZ	2.52	0.45
2:A:4:ARG:HB2	2:A:132:TRP:CH2	2.52	0.45
2:B:20:VAL:HG12	2:B:33:LEU:HD11	1.99	0.45
2:B:68:GLN:CB	2:B:71:HIS:CD2	3.00	0.45
2:B:99:GLU:HA	2:B:139:THR:OG1	2.16	0.45
2:A:232:ASN:OD1	2:A:235:GLU:HG3	2.17	0.45
2:B:220:GLU:HB2	2:B:226:ARG:CG	2.47	0.44
2:B:8:ILE:HG12	2:B:170:LEU:CB	2.46	0.44
2:B:220:GLU:O	2:B:226:ARG:HD3	2.18	0.44
1:D:10:DT:H2"	1:D:11:DT:C5'	2.46	0.44
2:B:157:ASN:H	2:B:157:ASN:HD22	1.65	0.44
2:A:58:LYS:HG3	2:A:59:HIS:CD2	2.52	0.44
2:A:153:ILE:HG21	2:B:153:ILE:HG21	1.99	0.44
2:B:15:ASN:ND2	3:B:310:HOH:O	2.51	0.44
2:B:121:ILE:HG12	2:B:123:TYR:O	2.17	0.44
2:B:45:GLU:OE1	2:B:91:ILE:HD12	2.18	0.43
2:B:54:LYS:O	2:B:57:GLU:HG3	2.18	0.43
2:A:87:ILE:HG23	2:A:130:ALA:HB3	1.99	0.43
2:A:193:HIS:HB3	3:A:269:HOH:O	2.18	0.43
2:B:88:ALA:HB3	2:B:131:HIS:ND1	2.33	0.43
2:B:135:GLY:O	2:B:168:VAL:HA	2.18	0.43
2:B:170:LEU:HD23	2:B:171:GLN:N	2.33	0.43
2:A:69:GLN:NE2	3:A:306:HOH:O	2.50	0.43
2:A:99:GLU:O	2:A:100:ASN:HB2	2.18	0.43
2:A:91:ILE:C	2:A:92:LYS:HG2	2.39	0.43
2:B:169:PHE:C	2:B:169:PHE:CD1	2.91	0.43
2:A:23:ILE:HG22	2:B:23:ILE:HG22	2.00	0.43
2:B:99:GLU:HA	2:B:139:THR:HG1	1.84	0.43
2:A:20:VAL:HG21	2:A:43:ILE:HG21	1.99	0.43
2:B:3:LEU:HD22	2:B:61:TYR:CE2	2.54	0.42
2:B:62:ILE:HD11	2:B:80:PRO:HG3	2.00	0.42
2:A:24:ILE:HD12	2:B:24:ILE:CD1	2.48	0.42
2:A:48:SER:HA	2:A:51:ILE:HD12	1.99	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:86:LYS:HB2	2:A:127:GLN:O	2.19	0.42
2:A:236:TYR:O	2:A:239:TRP:HB3	2.19	0.42
2:B:32:PRO:HG2	2:B:147:SER:O	2.18	0.42
1:D:6:DT:C4	1:D:7:DA:N7	2.87	0.42
2:A:206:PHE:HE1	2:A:233:ILE:HD13	1.84	0.42
1:C:8:DT:OP2	2:A:92:LYS:HE3	2.19	0.42
2:B:31:TYR:CD1	2:B:31:TYR:N	2.87	0.42
2:B:53:ASN:O	2:B:57:GLU:HG3	2.20	0.42
2:B:72:TYR:O	2:B:121:ILE:HA	2.20	0.42
2:B:215:TYR:O	2:B:230:TYR:OH	2.34	0.42
1:D:2:DA:H2"	1:D:3:DA:H5'	2.01	0.41
2:A:214:ASP:OD2	2:A:244:ARG:NH1	2.53	0.41
2:B:82:GLU:CD	2:B:85:LYS:HD3	2.41	0.41
2:B:207:ASP:HB2	3:B:321:HOH:O	2.19	0.41
2:A:24:ILE:HD12	2:B:24:ILE:HD13	2.01	0.41
2:A:2:SER:N	2:A:5:SER:HG	2.19	0.41
2:A:120:ASN:ND2	3:A:258:HOH:O	2.48	0.41
2:A:85:LYS:HA	2:A:129:ILE:CG2	2.50	0.41
2:A:103:ILE:HD13	2:A:137:VAL:HG21	2.02	0.40
2:A:189:ILE:HG23	2:A:216:TRP:CE2	2.56	0.40
2:B:92:LYS:HD3	2:B:106:THR:O	2.20	0.40
2:A:31:TYR:CD1	2:A:31:TYR:N	2.90	0.40
2:A:137:VAL:O	2:A:166:VAL:HA	2.21	0.40
2:A:209:GLU:O	2:A:209:GLU:HG3	2.21	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
2	A	242/245 (99%)	226 (93%)	15 (6%)	1 (0%)	34 29

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
2	B	242/245 (99%)	225 (93%)	13 (5%)	4 (2%)	9 3
All	All	484/490 (99%)	451 (93%)	28 (6%)	5 (1%)	15 9

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	144	ARG
2	B	146	SER
2	B	145	LYS
2	B	162	PRO
2	A	100	ASN

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
2	A	197/221 (89%)	182 (92%)	15 (8%)	13 8
2	B	197/221 (89%)	187 (95%)	10 (5%)	24 20
All	All	394/442 (89%)	369 (94%)	25 (6%)	18 13

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	6	ASP
2	A	8	ILE
2	A	57	GLU
2	A	71	HIS
2	A	78	TYR
2	A	90	ASP
2	A	92	LYS
2	A	96	THR
2	A	154	ASN
2	A	188	ASN
2	A	203	LYS

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Mol	Chain	Res	Type
2	A	209	GLU
2	A	228	ASP
2	A	231	ASN
2	A	234	SER
2	B	6	ASP
2	B	57	GLU
2	B	68	GLN
2	B	81	SER
2	B	90	ASP
2	B	119	LYS
2	B	188	ASN
2	B	201	GLU
2	B	203	LYS
2	B	233	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	69	GLN
2	A	193	HIS
2	A	238	ASN
2	B	15	ASN
2	B	70	ASN
2	B	71	HIS
2	B	157	ASN
2	B	231	ASN

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 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)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.