



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:03 AM UTC

PDB ID : 3C28 / pdb_00003c28
Title : Crystal structure of the product synapse complex
Authors : Ghosh, K.; Van Duyne, G.D.
Deposited on : 2008-01-24
Resolution : 2.60 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
Xtriage (Phenix)	:	2.0
EDS	:	3.0
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

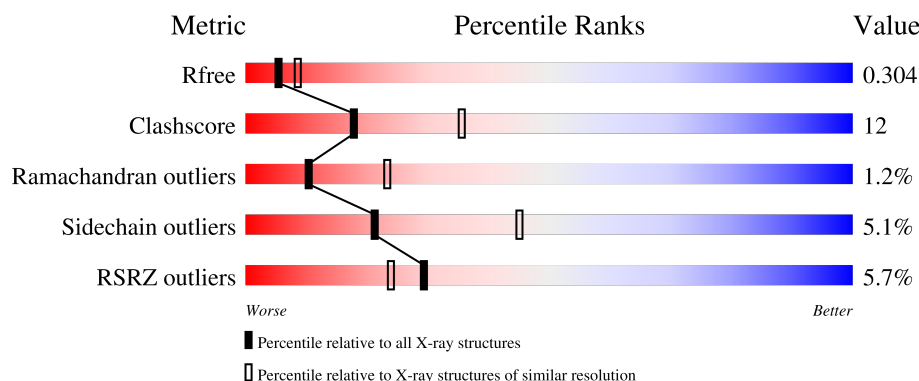
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	C	34	3% 65% 32% .
2	D	34	85% 15%
3	A	322	8% 66% 24% . 7%
3	B	322	4% 76% 20% ...

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6602 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 DNA, chain C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	34	Total	C	N	O	P	0	0	0
			698	336	123	205	34			

- Molecule 2 is a DNA chain called LoxP DNA, chain D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	34	Total	C	N	O	P	0	0	0
			696	335	124	203	34			

- Molecule 3 is a protein called Recombinase cre.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	299	Total	C	N	O	S	0	0	0
			2385	1487	452	431	15			
3	B	315	Total	C	N	O	S	0	0	0
			2499	1555	478	451	15			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	42	Total	O	0	0
			42	42		
4	D	38	Total	O	0	0
			38	38		
4	A	93	Total	O	0	0
			93	93		
4	B	151	Total	O	0	0
			151	151		

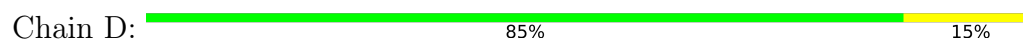
3 Residue-property plots

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.

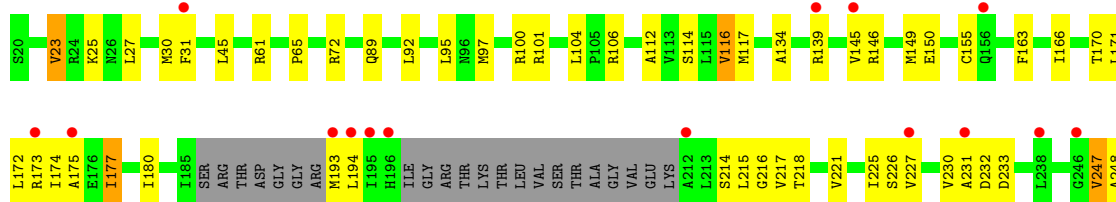
- Molecule 1: LoxP DNA, chain C



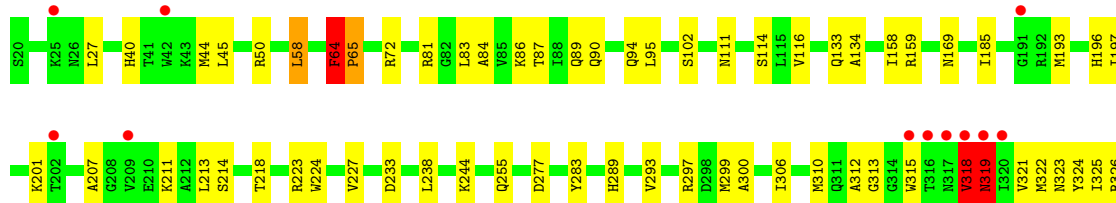
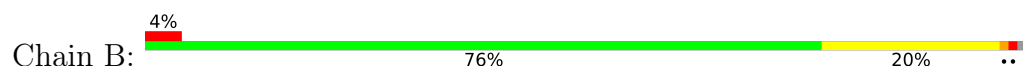
- Molecule 2: LoxP DNA, chain D



- Molecule 3: Recombinase cre



- Molecule 3: Recombinase cre



ASN	
LEU	
ASP	
SER	
GLU	
THR	
G333	
ALA	
M336	
V336	
R337	
L338	
I339	
E340	
D341	

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	108.25Å 122.45Å 178.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.10 – 2.60 40.10 – 2.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (40.10-2.60) 99.9 (40.10-2.60)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.53 (at 2.61Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.225 , 0.292 (Not available) , 0.304	Depositor DCC
R_{free} test set	1871 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	56.8	Xtriage
Anisotropy	0.112	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 38.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6602	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% 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](#)

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	0.42	0/782	1.19	1/1205 (0.1%)
2	D	0.42	0/780	1.18	0/1201
3	A	0.56	0/2423	0.79	0/3264
3	B	0.62	0/2538	0.84	3/3417 (0.1%)
All	All	0.56	0/6523	0.93	4/9087 (0.0%)

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	1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	64	PHE	CA-C-N	-9.58	107.86	119.84
3	B	64	PHE	C-N-CA	-9.58	107.86	119.84
1	C	18	DT	P-O3'-C3'	7.13	130.90	120.20
3	B	319	ASN	N-CA-C	5.39	122.28	110.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	B	64	PHE	Peptide

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	C	698	0	388	18	0
2	D	696	0	387	4	0
3	A	2385	0	2393	78	0
3	B	2499	0	2525	72	0
4	A	93	0	0	20	0
4	B	151	0	0	10	0
4	C	42	0	0	3	0
4	D	38	0	0	1	0
All	All	6602	0	5693	146	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 12.

All (146) 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:318:VAL:HG23	4:B:483:HOH:O	1.49	1.11
3:A:171:LEU:HD11	3:B:335:MET:HE3	1.41	1.02
1:C:18:DT:C7	3:B:319:ASN:HD21	1.75	0.99
1:C:18:DT:H71	3:B:319:ASN:HD21	1.26	0.99
1:C:18:DT:H71	3:B:319:ASN:ND2	1.83	0.92
1:C:18:DT:C7	3:B:319:ASN:ND2	2.35	0.86
3:A:72:ARG:HG3	3:A:116:VAL:HG21	1.56	0.86
3:A:72:ARG:HD2	4:A:424:HOH:O	1.77	0.84
3:A:171:LEU:CD1	3:B:335:MET:HE3	2.08	0.83
3:A:171:LEU:HD13	4:A:430:HOH:O	1.79	0.83
3:B:64:PHE:O	3:B:65:PRO:C	2.20	0.81
3:B:27:LEU:HD13	3:B:102:SER:OG	1.84	0.78
3:A:247:VAL:HG23	4:A:428:HOH:O	1.91	0.70
3:B:169:ASN:HD21	3:B:214:SER:H	1.38	0.70
3:B:306:ILE:HG22	3:B:310:MET:CE	2.23	0.69
3:A:318:VAL:HG12	3:A:322:MET:HG2	1.75	0.68
3:A:112:ALA:C	4:A:424:HOH:O	2.37	0.67
3:A:215:LEU:HD23	3:A:216:GLY:H	1.59	0.67
3:A:233:ASP:O	4:A:427:HOH:O	2.13	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:215:LEU:HD22	3:B:340:GLU:HG2	1.77	0.65
3:A:171:LEU:HD11	3:B:335:MET:CE	2.20	0.65
1:C:13:DT:H2'	3:B:87:THR:HG22	1.79	0.64
3:B:233:ASP:HB3	4:B:477:HOH:O	1.97	0.64
3:A:45:LEU:HD22	3:A:97:MET:HE1	1.79	0.64
3:A:299:MET:HE3	3:A:309:ILE:HG23	1.79	0.64
3:A:45:LEU:HD22	3:A:97:MET:CE	2.28	0.64
3:B:300:ALA:CB	3:B:325:ILE:HD12	2.28	0.63
3:A:277:ASP:HB2	3:A:284:LEU:HD13	1.81	0.63
1:C:13:DT:H6	3:B:87:THR:HG22	1.63	0.63
3:B:185:ILE:HD11	3:B:238:LEU:HD22	1.81	0.62
3:A:231:ALA:HB2	4:A:414:HOH:O	1.98	0.62
3:B:64:PHE:O	3:B:65:PRO:O	2.16	0.62
3:A:248:ALA:HB2	4:A:367:HOH:O	2.00	0.60
3:A:310:MET:SD	3:A:318:VAL:HG22	2.41	0.60
3:B:90:GLN:HE22	3:B:94:GLN:HE21	1.49	0.59
3:A:112:ALA:HB1	4:A:424:HOH:O	2.03	0.59
3:B:90:GLN:NE2	3:B:94:GLN:HE21	2.01	0.58
3:A:193:MET:HE2	3:A:218:THR:HG23	1.85	0.58
3:B:193:MET:HG3	3:B:218:THR:HG23	1.85	0.58
1:C:27:DC:P	4:C:76:HOH:O	2.61	0.58
3:A:193:MET:HE1	3:A:221:VAL:HB	1.84	0.58
1:C:13:DT:C6	3:B:87:THR:HG22	2.40	0.57
3:B:223:ARG:O	3:B:227:VAL:HG23	2.03	0.57
1:C:13:DT:OP2	3:B:87:THR:HG21	2.05	0.56
3:B:196:HIS:C	3:B:197:ILE:HD12	2.31	0.56
3:A:45:LEU:HD13	3:A:97:MET:CE	2.36	0.56
3:B:306:ILE:HG22	3:B:310:MET:HE3	1.87	0.55
3:A:296:ALA:HA	3:A:299:MET:HE2	1.88	0.55
3:A:145:VAL:CG1	3:A:149:MET:HE2	2.36	0.55
3:B:211:LYS:NZ	3:B:312:ALA:O	2.40	0.53
1:C:13:DT:O4	3:B:44:MET:HE2	2.09	0.53
3:B:64:PHE:CD2	3:B:65:PRO:HD3	2.44	0.52
3:B:321:VAL:HG23	4:B:425:HOH:O	2.10	0.52
3:A:89:GLN:HG2	3:A:117:MET:HE1	1.92	0.52
3:A:225:ILE:HG23	4:A:414:HOH:O	2.10	0.51
3:A:100:ARG:HH11	3:A:106:ARG:HD2	1.75	0.51
3:B:193:MET:HB2	3:B:213:LEU:HD12	1.93	0.51
1:C:24:DA:H2'	1:C:25:DT:C6	2.46	0.51
3:A:312:ALA:HB1	4:A:430:HOH:O	2.10	0.51
3:A:116:VAL:HG23	4:A:424:HOH:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:25:DT:H71	3:B:40:HIS:CD2	2.46	0.51
3:B:306:ILE:CG2	3:B:310:MET:CE	2.88	0.51
1:C:13:DT:H2'	3:B:87:THR:CG2	2.39	0.51
3:A:101:ARG:HH11	3:B:111:ASN:ND2	2.09	0.50
3:A:248:ALA:N	4:A:367:HOH:O	2.43	0.50
3:B:326:ARG:O	4:B:480:HOH:O	2.18	0.50
3:A:146:ARG:O	3:A:150:GLU:HB2	2.12	0.50
3:B:72:ARG:HG3	3:B:116:VAL:HG11	1.94	0.50
1:C:19:DA:H2''	1:C:20:DT:OP2	2.12	0.49
3:A:139:ARG:HH21	3:B:338:LEU:CD1	2.26	0.49
3:B:159:ARG:HB2	3:B:224:TRP:CZ3	2.48	0.49
3:B:325:ILE:HG22	3:B:325:ILE:O	2.13	0.49
3:B:233:ASP:HB3	4:B:482:HOH:O	2.13	0.49
3:B:300:ALA:HB2	3:B:325:ILE:HD12	1.94	0.49
3:A:89:GLN:CG	3:A:117:MET:HE1	2.44	0.48
3:A:72:ARG:HD3	4:A:351:HOH:O	2.13	0.48
3:B:27:LEU:CD1	3:B:102:SER:OG	2.60	0.48
3:A:100:ARG:HD2	3:A:106:ARG:HD2	1.96	0.48
3:B:300:ALA:HB3	3:B:325:ILE:HD12	1.95	0.48
3:A:293:VAL:HG22	3:A:324:TYR:CD1	2.49	0.48
3:A:23:VAL:HG12	4:A:420:HOH:O	2.14	0.48
3:A:301:ARG:NH2	3:A:328:LEU:HD21	2.29	0.48
3:A:317:ASN:N	3:A:317:ASN:OD1	2.45	0.47
4:D:56:HOH:O	3:A:320:ILE:HD11	2.13	0.47
3:B:313:GLY:HA3	3:B:315:TRP:CZ3	2.49	0.47
1:C:21:DG:H5''	4:C:72:HOH:O	2.15	0.47
3:A:163:PHE:CE1	3:A:261:LEU:HD22	2.49	0.47
3:B:81:ARG:NH1	3:B:83:LEU:HD11	2.30	0.47
3:B:333:GLY:O	3:B:335:MET:N	2.48	0.47
3:A:116:VAL:CG2	4:A:424:HOH:O	2.62	0.46
3:A:291:ALA:HB3	4:A:358:HOH:O	2.15	0.46
3:A:45:LEU:HD13	3:A:97:MET:HE2	1.97	0.46
3:A:139:ARG:HH21	3:B:338:LEU:HD13	1.79	0.46
3:A:341:ASP:C	3:A:341:ASP:OD1	2.58	0.46
3:B:255:GLN:NE2	4:B:383:HOH:O	2.48	0.46
2:D:16:DC:H2''	2:D:17:DA:C8	2.51	0.46
3:A:72:ARG:CG	3:A:116:VAL:HG21	2.37	0.46
3:A:306:ILE:HD13	3:A:309:ILE:HD12	1.96	0.46
3:A:65:PRO:HB3	3:A:104:LEU:HD13	1.98	0.45
3:A:166:ILE:O	3:A:170:THR:HG23	2.16	0.45
3:B:133:GLN:NE2	3:B:323:ASN:HB3	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:134:ALA:HA	3:B:283:TYR:CD1	2.52	0.45
3:A:134:ALA:HA	3:A:283:TYR:CD1	2.51	0.45
3:A:173:ARG:O	3:A:177:ILE:HG13	2.18	0.44
3:A:214:SER:OG	3:A:215:LEU:N	2.51	0.44
3:A:248:ALA:CB	4:A:367:HOH:O	2.62	0.44
3:A:271:LEU:C	3:A:271:LEU:HD13	2.42	0.44
3:B:84:ALA:HB3	3:B:87:THR:HG23	2.00	0.44
3:A:311:GLN:CG	3:B:322:MET:HE1	2.48	0.43
3:B:72:ARG:HG3	3:B:116:VAL:CG1	2.49	0.43
3:A:230:VAL:HG23	4:A:414:HOH:O	2.19	0.43
3:A:325:ILE:HG22	3:A:331:GLU:HG3	2.00	0.43
3:B:299:MET:HE1	3:B:312:ALA:CB	2.49	0.43
1:C:18:DT:H72	3:B:319:ASN:ND2	2.31	0.43
3:B:244:LYS:NZ	4:B:486:HOH:O	2.51	0.43
2:D:23:DT:O2	3:B:201:LYS:HE3	2.18	0.43
3:A:174:ILE:HD12	3:A:258:THR:HB	2.00	0.43
3:A:320:ILE:HG23	4:A:395:HOH:O	2.19	0.43
3:B:158:ILE:HD13	3:B:227:VAL:HG21	2.00	0.42
3:A:45:LEU:HD22	3:A:97:MET:HE3	2.00	0.42
3:A:310:MET:HE1	3:A:317:ASN:N	2.34	0.42
3:B:293:VAL:HG22	3:B:324:TYR:CE1	2.54	0.42
3:B:333:GLY:C	3:B:336:VAL:HG22	2.44	0.42
2:D:24:DA:H2'	2:D:25:DT:C6	2.54	0.42
3:A:311:GLN:HE21	3:B:322:MET:HE3	1.84	0.42
3:A:30:MET:HE2	3:A:30:MET:HB3	1.94	0.42
3:A:217:VAL:HG23	4:A:364:HOH:O	2.19	0.42
3:A:27:LEU:O	3:A:31:PHE:CD2	2.73	0.42
3:A:193:MET:HG3	3:A:218:THR:HG23	2.01	0.42
3:B:86:LYS:HA	3:B:89:GLN:HE21	1.84	0.42
3:B:207:ALA:HB3	4:B:435:HOH:O	2.19	0.42
3:A:194:LEU:N	3:A:194:LEU:HD12	2.35	0.42
3:B:58:LEU:HD13	4:B:437:HOH:O	2.20	0.42
3:A:177:ILE:HG23	3:A:180:ILE:HD12	2.02	0.42
3:A:226:SER:OG	3:A:227:VAL:N	2.53	0.42
3:A:89:GLN:HG2	3:A:117:MET:CE	2.49	0.41
3:B:297:ARG:HA	3:B:325:ILE:HD13	2.02	0.41
3:A:304:VAL:HG22	3:A:308:GLU:OE1	2.20	0.41
3:A:311:GLN:HG3	3:B:322:MET:HE1	2.01	0.41
1:C:28:DG:OP1	3:A:276:LYS:NZ	2.48	0.40
3:B:197:ILE:HD12	3:B:197:ILE:N	2.37	0.40
1:C:7:DT:H2'	1:C:8:DT:H71	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:11:DT:OP2	3:B:50:ARG:NH1	2.54	0.40
3:B:319:ASN:CG	4:B:400:HOH:O	2.64	0.40
4:C:47:HOH:O	3:A:175:ALA:HB3	2.22	0.40
3:A:101:ARG:HH11	3:B:111:ASN:HD22	1.70	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	291/322 (90%)	272 (94%)	17 (6%)	2 (1%)	18	38
3	B	310/322 (96%)	301 (97%)	4 (1%)	5 (2%)	7	16
All	All	601/644 (93%)	573 (95%)	21 (4%)	7 (1%)	10	23

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	B	64	PHE
3	A	232	ASP
3	B	277	ASP
3	B	65	PRO
3	A	319	ASN
3	B	319	ASN
3	B	318	VAL

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	251/269 (93%)	232 (92%)	19 (8%)	12	27
3	B	263/269 (98%)	256 (97%)	7 (3%)	39	67
All	All	514/538 (96%)	488 (95%)	26 (5%)	21	45

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	23	VAL
3	A	25	LYS
3	A	61	ARG
3	A	92	LEU
3	A	95	LEU
3	A	114	SER
3	A	116	VAL
3	A	155	CYS
3	A	172	LEU
3	A	177	ILE
3	A	247	VAL
3	A	271	LEU
3	A	277	ASP
3	A	304	VAL
3	A	316	THR
3	A	317	ASN
3	A	318	VAL
3	A	323	ASN
3	A	341	ASP
3	B	45	LEU
3	B	58	LEU
3	B	95	LEU
3	B	114	SER
3	B	289	HIS
3	B	318	VAL
3	B	338	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	89	GLN

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Mol	Chain	Res	Type
3	A	144	GLN
3	A	156	GLN
3	A	269	HIS
3	A	281	GLN
3	A	311	GLN
3	A	319	ASN
3	B	40	HIS
3	B	89	GLN
3	B	94	GLN
3	B	111	ASN
3	B	169	ASN
3	B	255	GLN
3	B	269	HIS
3	B	319	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 oligosaccharides 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

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	34/34 (100%)	0.09	1 (2%) 53 48	36, 53, 87, 89	0
2	D	34/34 (100%)	0.01	0 100 100	34, 52, 71, 77	0
3	A	299/322 (92%)	0.80	25 (8%) 17 13	43, 69, 95, 101	0
3	B	315/322 (97%)	0.37	13 (4%) 41 36	35, 52, 78, 104	0
All	All	682/712 (95%)	0.53	39 (5%) 29 24	34, 59, 89, 104	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	B	319	ASN	4.8
3	B	333	GLY	4.4
3	A	156	GLN	4.3
3	A	314	GLY	4.2
3	A	212	ALA	3.5
3	A	315	TRP	3.1
3	A	139	ARG	3.1
3	A	311	GLN	3.0
3	B	318	VAL	3.0
3	A	325	ILE	3.0
3	A	280	GLY	2.8
3	A	194	LEU	2.7
3	A	282	ARG	2.7
3	A	231	ALA	2.7
3	A	31	PHE	2.7
3	B	209	VAL	2.6
3	A	145	VAL	2.4
3	B	320	ILE	2.4
3	A	319	ASN	2.4
3	B	339	LEU	2.3
3	A	193	MET	2.3

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Mol	Chain	Res	Type	RSRZ
3	A	173	ARG	2.3
3	A	246	GLY	2.3
3	A	175	ALA	2.3
3	A	296	ALA	2.3
3	B	316	THR	2.2
1	C	17	DG	2.2
3	A	227	VAL	2.2
3	A	196	HIS	2.2
3	B	191	GLY	2.2
3	A	272	ILE	2.2
3	B	317	ASN	2.2
3	B	25	LYS	2.2
3	A	195	ILE	2.1
3	B	315	TRP	2.1
3	B	202	THR	2.0
3	A	238	LEU	2.0
3	A	317	ASN	2.0
3	B	42	TRP	2.0

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

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.