



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 27, 2026 – 11:05 PM UTC

PDB ID : 8SSQ / pdb_00008ssq
Title : ZnFs 3-11 of CCCTC-binding factor (CTCF) Complexed with 35mer DNA 35-4
Authors : Horton, J.R.; Yang, J.; Cheng, X.
Deposited on : 2023-05-08
Resolution : 3.12 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

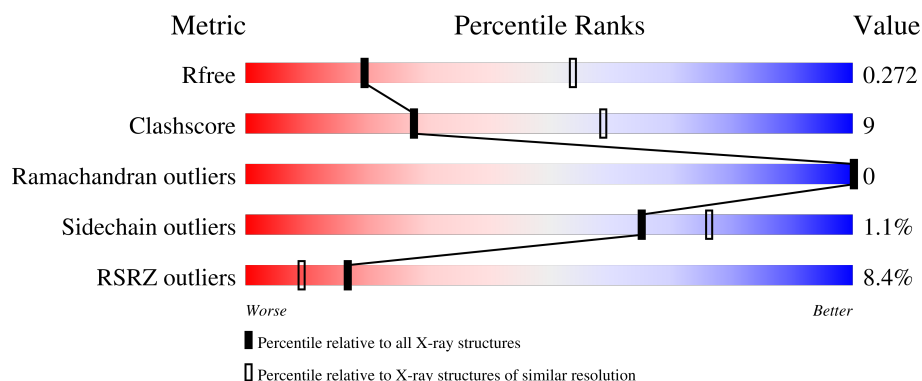
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.12 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.



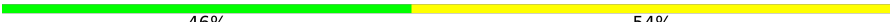
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1816 (3.14-3.10)
Clashscore	190562	1906 (3.14-3.10)
Ramachandran outliers	187476	1802 (3.14-3.10)
Sidechain outliers	187428	1802 (3.14-3.10)
RSRZ outliers	180081	1816 (3.14-3.10)

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	A	288	9% 67% 23% 10%
1	D	288	9% 63% 17% 20%
2	B	35	60% 40%
2	E	35	60% 40%
3	C	35	60% 40%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	F	35	 A horizontal bar chart showing the quality of chain F. The bar is divided into two segments: a green segment on the left representing 46% and a yellow segment on the right representing 54%. The percentages are labeled below the respective segments.

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6941 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 protein called Transcriptional repressor CTCF.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	260	Total	C	N	O	S	0	0	0
			2152	1330	425	368	29			
1	D	230	Total	C	N	O	S	0	0	0
			1899	1175	372	326	26			

- Molecule 2 is a DNA chain called DNA (35-MER) Strand I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	35	Total	C	N	O	P	0	0	0
			711	340	125	212	34			
2	E	35	Total	C	N	O	P	0	0	0
			711	340	125	212	34			

- Molecule 3 is a DNA chain called DNA (35-MER) Strand 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	35	Total	C	N	O	P	0	0	0
			717	341	139	203	34			
3	F	35	Total	C	N	O	P	0	0	0
			717	341	139	203	34			

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	9	Total	Zn	0	0
			9	9		
4	D	8	Total	Zn	0	0
			8	8		

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Na	0	0
			1	1		
5	E	1	Total	Na	0	0
			1	1		

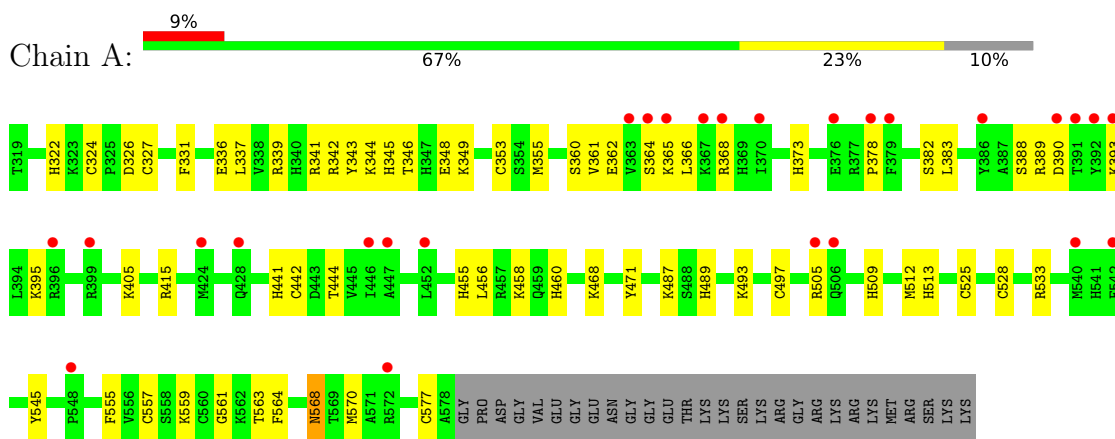
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	2	Total	O	0	0
			2	2		
6	B	4	Total	O	0	0
			4	4		
6	C	2	Total	O	0	0
			2	2		
6	D	6	Total	O	0	0
			6	6		
6	E	1	Total	O	0	0
			1	1		

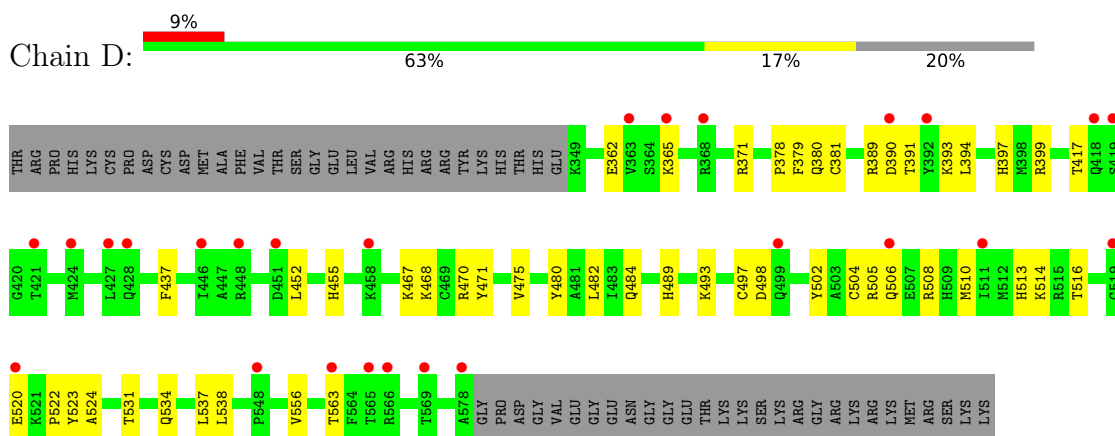
3 Residue-property plots [i](#)

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: Transcriptional repressor CTCF



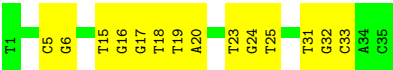
• Molecule 1: Transcriptional repressor CTCF



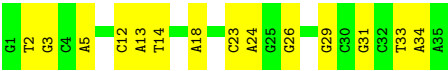
• Molecule 2: DNA (35-MER) Strand I



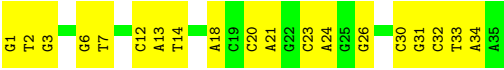
• Molecule 2: DNA (35-MER) Strand I



● Molecule 3: DNA (35-MER) Strand 2



● Molecule 3: DNA (35-MER) Strand 2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	352.07Å 67.77Å 61.19Å 90.00° 92.59° 90.00°	Depositor
Resolution (Å)	41.86 – 3.12 41.86 – 3.12	Depositor EDS
% Data completeness (in resolution range)	97.8 (41.86-3.12) 90.8 (41.86-3.12)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.62 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.240 , 0.271 0.243 , 0.272	Depositor DCC
R_{free} test set	1278 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	88.9	Xtriage
Anisotropy	0.210	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 178.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.033 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6941	wwPDB-VP
Average B, all atoms (Å ²)	185.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.26% 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](#)

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

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	A	0.09	0/2209	0.24	0/2954
1	D	0.07	0/1948	0.20	0/2604
2	B	0.17	0/795	0.36	0/1225
2	E	0.17	0/795	0.36	0/1225
3	C	0.19	0/806	0.36	0/1241
3	F	0.16	0/806	0.31	0/1241
All	All	0.13	0/7359	0.29	0/10490

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	A	2152	0	2078	47	0
1	D	1899	0	1834	32	0
2	B	711	0	397	11	0
2	E	711	0	397	11	0
3	C	717	0	393	13	0
3	F	717	0	393	15	0
4	A	9	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	8	0	0	0	0
5	A	1	0	0	0	0
5	E	1	0	0	0	0
6	A	2	0	0	1	0
6	B	4	0	0	0	0
6	C	2	0	0	0	0
6	D	6	0	0	0	0
6	E	1	0	0	0	0
All	All	6941	0	5492	112	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 9.

All (112) 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)
1:A:344:LYS:HE2	3:C:29:DG:H5''	1.65	0.78
1:A:353:CYS:SG	1:A:373:HIS:CD2	2.85	0.70
1:A:390:ASP:HB2	1:A:393:LYS:HG2	1.73	0.70
1:D:520:GLU:HG2	1:D:522:PRO:HD3	1.78	0.65
1:A:455:HIS:ND1	3:C:18:DA:OP1	2.27	0.64
1:A:362:GLU:HB2	1:A:365:LYS:HG2	1.81	0.63
1:A:487:LYS:NZ	2:B:23:DT:OP1	2.32	0.62
1:D:390:ASP:HB3	1:D:393:LYS:HG2	1.80	0.62
2:B:15:DT:H2''	2:B:16:DG:C8	2.36	0.60
1:D:380:GLN:NE2	1:D:381:CYS:O	2.34	0.60
3:F:13:DA:H2''	3:F:14:DT:H5''	1.84	0.59
1:D:397:HIS:ND1	3:F:24:DA:OP1	2.36	0.59
1:A:533:ARG:NH1	3:C:5:DA:OP1	2.34	0.59
1:D:523:TYR:HB3	1:D:538:LEU:HD22	1.83	0.58
2:E:31:DT:H2''	2:E:32:DG:C8	2.39	0.58
1:A:388:SER:HB2	1:A:393:LYS:HB2	1.86	0.58
1:D:389:ARG:NH2	3:F:26:DG:OP2	2.36	0.58
1:A:456:LEU:HD23	1:A:460:HIS:HD2	1.68	0.57
1:D:399:ARG:NH1	1:D:417:THR:O	2.36	0.57
1:D:510:MET:HE3	1:D:514:LYS:HE3	1.85	0.57
2:B:16:DG:H2''	2:B:17:DG:C8	2.39	0.57
1:A:389:ARG:NH2	3:C:26:DG:OP2	2.38	0.56
3:C:13:DA:H2''	3:C:14:DT:H5''	1.88	0.56
1:D:513:HIS:O	1:D:516:THR:OG1	2.22	0.56
2:E:5:DC:H2''	2:E:6:DG:C8	2.41	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:PHE:HD2	1:A:336:GLU:HB3	1.72	0.55
1:A:360:SER:HB3	1:A:365:LYS:HB2	1.89	0.54
1:D:437:PHE:HB3	1:D:452:LEU:HD22	1.89	0.54
3:F:33:DT:H2'	3:F:34:DA:C8	2.42	0.54
2:E:15:DT:H2''	2:E:16:DG:C8	2.42	0.54
1:A:395:LYS:NZ	2:B:8:DC:OP2	2.36	0.53
1:D:556:VAL:HG22	1:D:563:THR:HG22	1.90	0.53
1:A:458:LYS:NZ	6:A:801:HOH:O	2.33	0.53
2:B:17:DG:H1'	2:B:18:DT:H5'	1.90	0.53
1:A:471:TYR:HB2	1:A:489:HIS:HD2	1.74	0.52
1:D:493:LYS:HD3	1:D:504:CYS:HA	1.90	0.52
1:D:455:HIS:ND1	3:F:18:DA:OP1	2.39	0.52
1:D:493:LYS:HG2	1:D:505:ARG:HG3	1.92	0.51
1:A:442:CYS:SG	1:A:444:THR:OG1	2.62	0.51
1:A:568:ASN:ND2	2:B:31:DT:OP2	2.34	0.50
1:A:441:HIS:HB2	1:A:460:HIS:CD2	2.46	0.50
1:A:468:LYS:HD3	1:D:468:LYS:H	1.76	0.50
3:C:33:DT:H2''	3:C:34:DA:O5'	2.11	0.50
2:E:17:DG:H1'	2:E:18:DT:H5'	1.93	0.50
1:D:502:TYR:OH	3:F:7:DT:OP2	2.27	0.50
1:A:324:CYS:HB2	1:A:337:LEU:HD11	1.94	0.49
2:B:33:DC:H2''	2:B:34:DA:C8	2.47	0.49
3:F:2:DT:H2''	3:F:3:DG:C8	2.47	0.49
3:C:33:DT:H4'	3:C:34:DA:OP1	2.12	0.49
2:B:19:DT:H2''	2:B:20:DA:C8	2.47	0.49
1:D:470:ARG:NH2	2:E:25:DT:OP1	2.46	0.49
1:A:509:HIS:HA	1:A:512:MET:HE2	1.93	0.48
1:D:468:LYS:HA	1:D:475:VAL:HA	1.95	0.48
1:A:555:PHE:HB3	1:A:570:MET:HE2	1.96	0.48
3:F:23:DC:H2'	3:F:24:DA:C8	2.49	0.48
2:B:31:DT:H2''	2:B:32:DG:C8	2.49	0.47
1:A:525:CYS:HB3	1:A:528:CYS:SG	2.55	0.47
1:A:327:CYS:HB3	1:A:345:HIS:CE1	2.50	0.47
1:A:471:TYR:HB2	1:A:489:HIS:CD2	2.49	0.47
1:D:379:PHE:HB3	1:D:394:LEU:HD22	1.97	0.47
2:E:16:DG:H2''	2:E:17:DG:C8	2.49	0.47
3:F:30:DC:H2''	3:F:31:DG:C8	2.49	0.47
2:E:19:DT:H2''	2:E:20:DA:C8	2.49	0.47
1:A:564:PHE:HB2	1:A:570:MET:HB2	1.96	0.47
1:D:506:GLN:OE1	1:D:508:ARG:NE	2.48	0.46
2:B:16:DG:H2''	2:B:17:DG:H8	1.79	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:557:CYS:HB3	1:A:561:GLY:H	1.79	0.46
3:F:31:DG:H2''	3:F:32:DC:H5''	1.96	0.46
1:A:326:ASP:OD2	1:A:341:ARG:NH1	2.49	0.46
3:F:20:DC:H2''	3:F:21:DA:C8	2.51	0.46
3:C:23:DC:H2'	3:C:24:DA:C8	2.51	0.46
1:D:524:ALA:HA	1:D:531:THR:HG22	1.97	0.46
1:D:471:TYR:HB2	1:D:489:HIS:CD2	2.51	0.46
1:A:393:LYS:NZ	3:C:26:DG:O6	2.49	0.45
1:A:468:LYS:HB3	1:D:468:LYS:HB2	1.98	0.45
1:A:382:SER:HB2	1:A:383:LEU:HD12	1.97	0.45
1:D:378:PRO:HD2	1:D:389:ARG:HA	1.97	0.45
1:D:497:CYS:HA	1:D:510:MET:HE1	1.98	0.45
1:A:355:MET:HB3	1:A:373:HIS:NE2	2.32	0.45
1:A:441:HIS:HB2	1:A:460:HIS:NE2	2.32	0.45
1:A:360:SER:HB2	1:A:366:LEU:HB2	1.99	0.44
2:B:5:DC:H2''	2:B:6:DG:H8	1.82	0.44
1:A:497:CYS:SG	1:A:513:HIS:HD2	2.40	0.44
1:A:559:LYS:HD2	1:A:577:CYS:HB3	1.99	0.44
3:C:2:DT:H2''	3:C:3:DG:C8	2.53	0.44
1:A:405:LYS:O	1:A:415:ARG:NH2	2.51	0.44
2:E:31:DT:H2''	2:E:32:DG:H8	1.80	0.43
1:A:493:LYS:HG2	1:A:505:ARG:HG3	2.01	0.43
3:F:6:DG:H2'	3:F:7:DT:H71	2.00	0.43
1:A:342:ARG:HA	1:A:346:THR:OG1	2.18	0.43
3:F:33:DT:H2''	3:F:34:DA:OP1	2.18	0.43
2:E:23:DT:H2''	2:E:24:DG:C8	2.53	0.42
2:E:32:DG:H1'	2:E:33:DC:H5'	2.01	0.42
3:F:12:DC:H2''	3:F:13:DA:C8	2.53	0.42
1:D:467:LYS:HB3	1:D:482:LEU:HD22	2.02	0.42
1:A:322:HIS:HB3	1:A:337:LEU:HD22	2.02	0.42
1:D:362:GLU:HB2	1:D:365:LYS:HG2	2.01	0.42
2:E:16:DG:H2''	2:E:17:DG:H8	1.85	0.42
1:A:339:ARG:NH1	3:C:31:DG:N7	2.64	0.41
3:C:12:DC:H2''	3:C:13:DA:C8	2.55	0.41
3:F:1:DG:C8	3:F:2:DT:H72	2.55	0.41
1:A:349:LYS:HG2	1:A:361:VAL:HG23	2.02	0.41
1:A:326:ASP:HB2	1:A:345:HIS:CD2	2.55	0.41
1:D:534:GLN:HB2	1:D:537:LEU:HG	2.02	0.41
1:A:378:PRO:HD2	1:A:389:ARG:HA	2.03	0.41
1:D:480:TYR:O	1:D:484:GLN:HG2	2.21	0.41
1:D:497:CYS:SG	1:D:498:ASP:N	2.94	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:365:LYS:O	1:A:368:ARG:HB2	2.20	0.41
1:A:545:TYR:OH	3:C:3:DG:OP1	2.25	0.41
1:D:390:ASP:OD1	1:D:391:THR:N	2.54	0.41
1:D:371:ARG:HB3	1:D:389:ARG:HE	1.86	0.40
1:A:331:PHE:CD2	1:A:336:GLU:HB3	2.56	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	
1	A	258/288 (90%)	249 (96%)	9 (4%)	0	100	100
1	D	228/288 (79%)	226 (99%)	2 (1%)	0	100	100
All	All	486/576 (84%)	475 (98%)	11 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	
1	A	236/259 (91%)	231 (98%)	5 (2%)	47	69
1	D	208/259 (80%)	208 (100%)	0	100	100
All	All	444/518 (86%)	439 (99%)	5 (1%)	65	76

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	343	TYR
1	A	348	GLU
1	A	364	SER
1	A	563	THR
1	A	568	ASN

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

Mol	Chain	Res	Type
1	A	459	GLN
1	A	486	GLN
1	D	433	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](#)

Of 19 ligands modelled in this entry, 19 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	A	260/288 (90%)	0.50	27 (10%) 11 6	91, 156, 256, 308	0
1	D	230/288 (79%)	0.68	26 (11%) 10 5	88, 211, 282, 320	0
2	B	35/35 (100%)	-0.14	0 100 100	100, 145, 249, 286	0
2	E	35/35 (100%)	-0.11	0 100 100	140, 206, 375, 448	0
3	C	35/35 (100%)	0.02	0 100 100	88, 142, 234, 260	0
3	F	35/35 (100%)	0.01	0 100 100	132, 207, 385, 439	0
All	All	630/716 (87%)	0.44	53 (8%) 17 9	88, 176, 284, 448	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	392	TYR	10.8
1	A	392	TYR	5.6
1	A	390	ASP	5.5
1	D	428	GLN	5.4
1	D	563	THR	5.3
1	D	424	MET	5.2
1	A	379	PHE	5.2
1	D	363	VAL	5.1
1	D	390	ASP	4.9
1	A	391	THR	4.6
1	A	447	ALA	4.3
1	A	368	ARG	4.2
1	A	376	GLU	3.9
1	A	572	ARG	3.9
1	D	565	THR	3.8
1	A	505	ARG	3.7
1	A	393	LYS	3.5
1	A	364	SER	3.4
1	D	446	ILE	3.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	368	ARG	3.0
1	A	548	PRO	3.0
1	A	386	TYR	3.0
1	A	396	ARG	3.0
1	A	365	LYS	2.9
1	D	506	GLN	2.9
1	D	520	GLU	2.9
1	D	451	ASP	2.9
1	D	548	PRO	2.8
1	D	499	GLN	2.8
1	A	424	MET	2.8
1	A	363	VAL	2.8
1	A	367	LYS	2.7
1	A	446	ILE	2.7
1	A	540	MET	2.7
1	A	399	ARG	2.7
1	D	418	GLN	2.6
1	D	566	ARG	2.6
1	D	578	ALA	2.5
1	A	370	ILE	2.5
1	D	448	ARG	2.5
1	D	427	LEU	2.4
1	D	511	ILE	2.4
1	D	419	SER	2.3
1	D	458	LYS	2.3
1	D	569	THR	2.3
1	A	542	PHE	2.3
1	A	452	LEU	2.2
1	A	428	GLN	2.2
1	D	519	GLY	2.1
1	A	506	GLN	2.1
1	A	378	PRO	2.1
1	D	421	THR	2.0
1	D	365	LYS	2.0

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

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](#)

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
5	NA	A	710	1/1	0.73	0.26	122,122,122,122	0
5	NA	E	101	1/1	0.75	0.09	139,139,139,139	0
4	ZN	A	709	1/1	0.92	0.08	290,290,290,290	0
4	ZN	D	707	1/1	0.92	0.07	275,275,275,275	0
4	ZN	D	706	1/1	0.93	0.07	266,266,266,266	0
4	ZN	D	708	1/1	0.94	0.08	307,307,307,307	0
4	ZN	D	701	1/1	0.95	0.04	255,255,255,255	0
4	ZN	D	703	1/1	0.96	0.07	134,134,134,134	0
4	ZN	A	701	1/1	0.97	0.07	216,216,216,216	0
4	ZN	A	702	1/1	0.97	0.06	242,242,242,242	0
4	ZN	A	704	1/1	0.97	0.08	121,121,121,121	0
4	ZN	A	708	1/1	0.97	0.10	204,204,204,204	0
4	ZN	A	707	1/1	0.98	0.05	144,144,144,144	0
4	ZN	A	706	1/1	0.98	0.11	109,109,109,109	0
4	ZN	A	703	1/1	0.99	0.04	117,117,117,117	0
4	ZN	D	702	1/1	0.99	0.04	179,179,179,179	0
4	ZN	A	705	1/1	0.99	0.12	140,140,140,140	0
4	ZN	D	704	1/1	0.99	0.06	109,109,109,109	0
4	ZN	D	705	1/1	0.99	0.10	145,145,145,145	0

6.5 Other polymers [i](#)

There are no such residues in this entry.