



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 9, 2023 – 01:13 PM EDT

PDB ID : 8SSQ
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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.35.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35.1

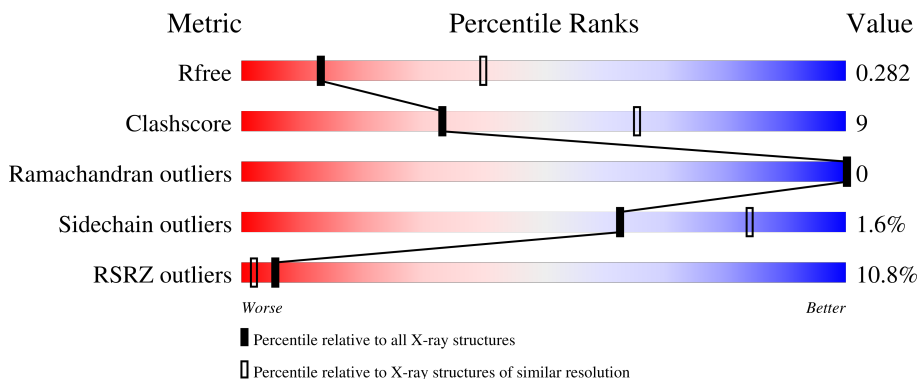
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	130704	1292 (3.14-3.10)
Clashscore	141614	1389 (3.14-3.10)
Ramachandran outliers	138981	1337 (3.14-3.10)
Sidechain outliers	138945	1337 (3.14-3.10)
RSRZ outliers	127900	1260 (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	<div style="display: flex; align-items: center;"> <div style="width: 7%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 68%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 22%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: grey;"></div> </div>
1	D	288	<div style="display: flex; align-items: center;"> <div style="width: 14%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 64%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 16%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 20%; height: 10px; background-color: grey;"></div> </div>
2	B	35	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 60%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 40%; height: 10px; background-color: yellow;"></div> </div>
2	E	35	<div style="display: flex; align-items: center;"> <div style="width: 14%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 60%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 40%; height: 10px; background-color: yellow;"></div> </div>
3	C	35	<div style="display: flex; align-items: center;"> <div style="width: 60%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 40%; height: 10px; background-color: yellow;"></div> </div>

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Mol	Chain	Length	Quality of chain
3	F	35	 <p>6% 51% 49%</p>

2 Entry composition [i](#)

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
			Total	C	N	O	S			
1	A	260	Total 2152	C 1330	N 425	O 368	S 29	0	0	0
1	D	230	Total 1899	C 1175	N 372	O 326	S 26	0	0	0

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

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

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

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

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

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

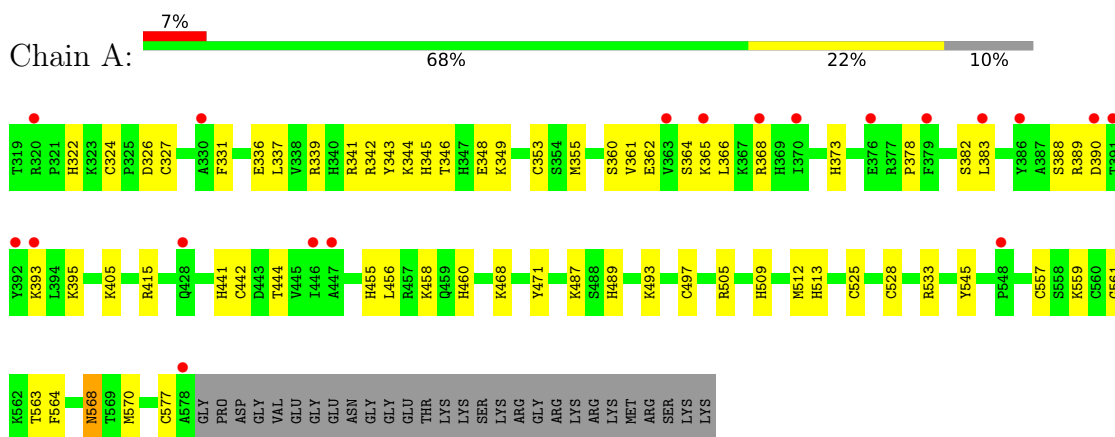
- Molecule 6 is water.

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

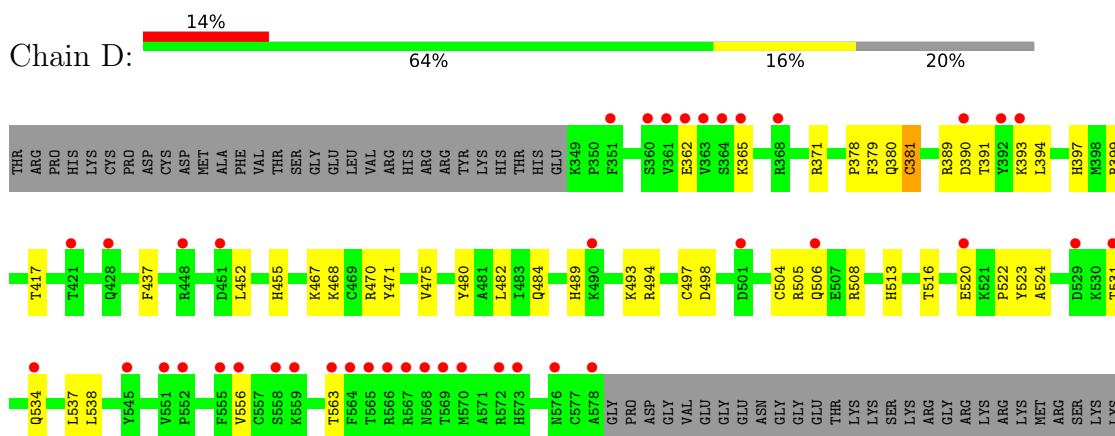
3 Residue-property plots [i](#)

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 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.

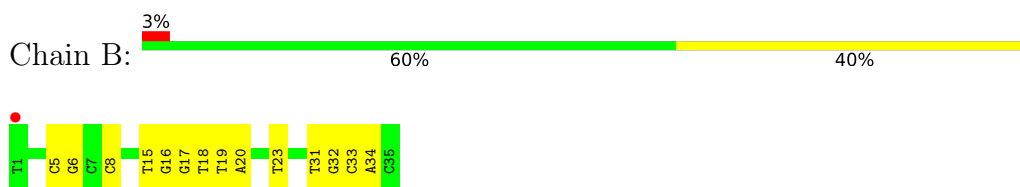
- Molecule 1: Transcriptional repressor CTCF



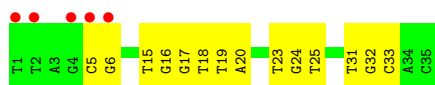
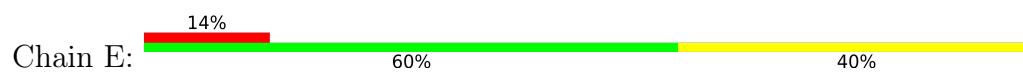
- Molecule 1: Transcriptional repressor CTCF



- 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 i

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 44.83 – 3.12	Depositor EDS
% Data completeness (in resolution range)	97.8 (41.86-3.12) 90.8 (44.83-3.12)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.62 (at 3.12Å)	Xtrriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.240 , 0.271 0.247 , 0.282	Depositor DCC
R_{free} test set	1278 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	88.9	Xtrriage
Anisotropy	0.210	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.23 , 79.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	0.033 for -h,-k,l	Xtrriage
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

Xtrriage'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	0.24	0/2209	0.49	0/2954
1	D	0.23	0/1948	0.47	0/2604
2	B	0.48	0/795	0.92	0/1225
2	E	0.47	0/795	0.92	0/1225
3	C	0.51	0/806	0.85	0/1241
3	F	0.47	0/806	0.81	0/1241
All	All	0.36	0/7359	0.70	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	46	0
1	D	1899	0	1834	29	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	13	0
4	A	9	0	0	0	0

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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	107	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 (107) 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
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
2:E:5:DC:H2''	2:E:6:DG:C8	2.41	0.56
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:509:HIS:HA	1:A:512:MET:HE2	1.91	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:458:LYS:NZ	6:A:801:HOH:O	2.33	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: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:D:468:LYS:HA	1:D:475:VAL:HA	1.95	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:D:513:HIS:O	1:D:516:THR:OG1	2.22	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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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
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
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%)	53	78
1	D	208/259 (80%)	206 (99%)	2 (1%)	76	89
All	All	444/518 (86%)	437 (98%)	7 (2%)	62	84

All (7) 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

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Mol	Chain	Res	Type
1	D	381	CYS
1	D	494	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

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%)	0.33	19 (7%) 15 6	91, 156, 256, 308	0
1	D	230/288 (79%)	0.81	41 (17%) 1 0	88, 211, 282, 320	0
2	B	35/35 (100%)	0.26	1 (2%) 51 29	100, 145, 249, 286	0
2	E	35/35 (100%)	0.37	5 (14%) 2 1	140, 206, 375, 448	0
3	C	35/35 (100%)	0.42	0 100 100	88, 142, 234, 260	0
3	F	35/35 (100%)	0.51	2 (5%) 23 11	132, 207, 385, 439	0
All	All	630/716 (87%)	0.52	68 (10%) 5 2	88, 176, 284, 448	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	392	TYR	13.0
1	D	569	THR	7.6
1	D	363	VAL	6.8
1	D	501	ASP	6.8
1	D	563	THR	6.4
1	A	392	TYR	6.2
1	D	565	THR	5.2
2	E	1	DT	5.2
1	D	361	VAL	4.7
3	F	35	DA	4.3
1	D	556	VAL	4.3
1	A	578	ALA	4.2
1	D	390	ASP	4.0
1	D	368	ARG	4.0
1	A	548	PRO	3.8
1	D	564	PHE	3.8
1	D	365	LYS	3.6
1	D	567	ARG	3.5
1	A	376	GLU	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	390	ASP	3.4
1	D	506	GLN	3.4
2	E	2	DT	3.4
2	E	5	DC	3.4
1	D	566	ARG	3.2
1	D	534	GLN	3.2
1	D	551	VAL	3.0
1	D	393	LYS	3.0
1	D	568	ASN	3.0
1	A	370	ILE	2.9
1	D	520	GLU	2.9
1	D	364	SER	2.9
1	D	428	GLN	2.9
1	D	552	PRO	2.9
1	D	545	TYR	2.8
1	D	558	SER	2.8
1	A	393	LYS	2.8
1	D	572	ARG	2.7
2	E	6	DG	2.7
1	D	360	SER	2.7
1	A	330	ALA	2.6
1	A	365	LYS	2.6
3	F	1	DG	2.6
1	D	451	ASP	2.6
1	D	578	ALA	2.5
1	A	391	THR	2.5
1	A	363	VAL	2.5
1	A	379	PHE	2.5
1	D	351	PHE	2.4
2	E	4	DG	2.4
1	D	448	ARG	2.3
1	D	555	PHE	2.3
2	B	1	DT	2.2
1	A	320	ARG	2.2
1	D	570	MET	2.2
1	D	573	HIS	2.1
1	A	368	ARG	2.1
1	D	362	GLU	2.1
1	D	421	THR	2.1
1	D	559	LYS	2.1
1	A	428	GLN	2.1
1	A	447	ALA	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	531	THR	2.0
1	A	386	TYR	2.0
1	A	446	ILE	2.0
1	D	576	ASN	2.0
1	D	529	ASP	2.0
1	D	490	LYS	2.0
1	A	383	LEU	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 monosaccharides 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
4	ZN	D	708	1/1	0.42	0.07	307,307,307,307	0
4	ZN	D	707	1/1	0.65	0.12	275,275,275,275	0
4	ZN	A	709	1/1	0.71	0.08	290,290,290,290	0
4	ZN	D	701	1/1	0.80	0.06	255,255,255,255	0
4	ZN	A	708	1/1	0.84	0.26	204,204,204,204	0
5	NA	E	101	1/1	0.86	0.33	139,139,139,139	0
5	NA	A	710	1/1	0.90	0.24	122,122,122,122	0
4	ZN	A	701	1/1	0.91	0.13	216,216,216,216	0
4	ZN	A	704	1/1	0.94	0.28	121,121,121,121	0
4	ZN	D	702	1/1	0.94	0.21	179,179,179,179	0
4	ZN	D	703	1/1	0.94	0.23	134,134,134,134	0
4	ZN	D	706	1/1	0.94	0.14	266,266,266,266	0
4	ZN	A	702	1/1	0.95	0.13	242,242,242,242	0
4	ZN	A	706	1/1	0.95	0.29	109,109,109,109	0
4	ZN	A	707	1/1	0.97	0.20	144,144,144,144	0
4	ZN	D	705	1/1	0.98	0.33	145,145,145,145	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	ZN	D	704	1/1	0.98	0.27	109,109,109,109	0
4	ZN	A	703	1/1	0.99	0.25	117,117,117,117	0
4	ZN	A	705	1/1	0.99	0.28	140,140,140,140	0

6.5 Other polymers [i](#)

There are no such residues in this entry.