



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 20, 2024 – 03:56 pm GMT

PDB ID : 7QFC
Title : Crystal structure of cytotoxin 13 from *Naja naja*, orthorhombic form
Authors : Samygina, V.R.; Dubova, K.M.; Bourenkov, G.; Utkin, Y.N.; Dubovskii, P.V.
Deposited on : 2021-12-05
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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

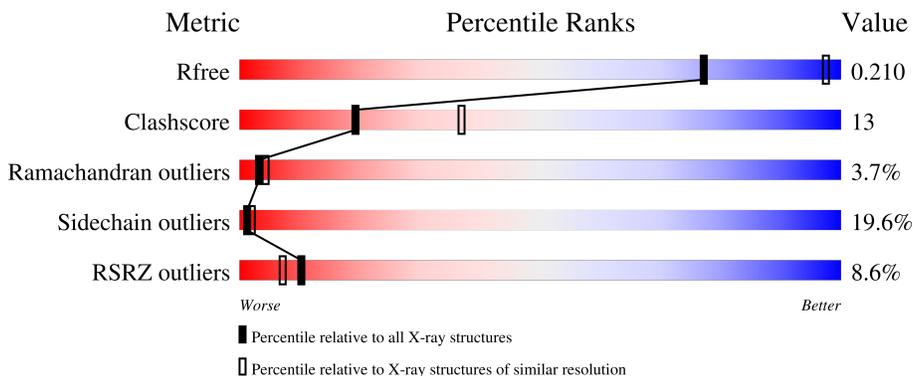
1 Overall quality at a glance

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}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (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	AAA	60	
1	BBB	60	
1	CCC	60	
1	DDD	60	
1	EEE	60	

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Mol	Chain	Length	Quality of chain
1	FFF	60	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into four segments: a red segment (22%), a green segment (38%), a yellow segment (50%), and an orange segment (12%). The percentages are labeled above and below the segments.</p>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 2801 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 Cytotoxin 13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	AAA	60	Total 464	C 294	N 81	O 79	S 10	0	0	0
1	BBB	60	Total 464	C 294	N 81	O 79	S 10	0	0	0
1	CCC	60	Total 464	C 294	N 81	O 79	S 10	0	0	0
1	DDD	60	Total 464	C 294	N 81	O 79	S 10	0	0	0
1	EEE	60	Total 464	C 294	N 81	O 79	S 10	0	0	0
1	FFF	60	Total 464	C 294	N 81	O 79	S 10	0	0	0

- Molecule 2 is water.

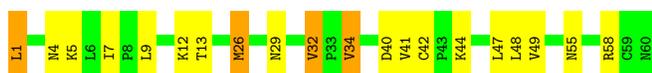
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	AAA	7	Total 7	O 7	0	0
2	BBB	3	Total 3	O 3	0	0
2	CCC	4	Total 4	O 4	0	0
2	DDD	1	Total 1	O 1	0	0
2	EEE	1	Total 1	O 1	0	0
2	FFF	1	Total 1	O 1	0	0

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: Cytotoxin 13

Chain AAA: 



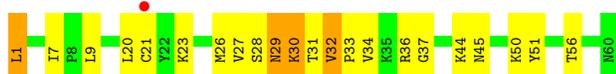
- Molecule 1: Cytotoxin 13

Chain BBB: 



- Molecule 1: Cytotoxin 13

Chain CCC: 



- Molecule 1: Cytotoxin 13

Chain DDD: 



- Molecule 1: Cytotoxin 13

Chain EEE: 



- Molecule 1: Cytotoxin 13



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	88.76Å 97.17Å 131.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.89 – 2.60 14.89 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.2 (14.89-2.60) 99.8 (14.89-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.20	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 2.61Å)	Xtrriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.205 , 0.255 0.210 , 0.210	Depositor DCC
R_{free} test set	913 reflections (5.14%)	wwPDB-VP
Wilson B-factor (Å ²)	79.7	Xtrriage
Anisotropy	0.252	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 63.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2801	wwPDB-VP
Average B, all atoms (Å ²)	100.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.95% 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	AAA	0.89	1/471 (0.2%)	1.20	1/632 (0.2%)
1	BBB	0.78	0/471	1.19	0/632
1	CCC	0.83	1/471 (0.2%)	1.19	0/632
1	DDD	0.79	0/471	1.11	1/632 (0.2%)
1	EEE	0.71	0/471	1.06	0/632
1	FFF	0.78	0/471	1.08	0/632
All	All	0.80	2/2826 (0.1%)	1.14	2/3792 (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	CCC	0	1
1	DDD	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AAA	55	ASN	C-O	-5.23	1.13	1.23
1	CCC	1	LEU	N-CA	5.04	1.56	1.46

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	DDD	42	CYS	CA-CB-SG	-5.57	103.97	114.00
1	AAA	34	VAL	N-CA-CB	-5.08	100.33	111.50

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	CCC	20	LEU	Peptide
1	DDD	39	ILE	Peptide

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	AAA	464	0	492	7	0
1	BBB	464	0	492	16	0
1	CCC	464	0	492	14	0
1	DDD	464	0	492	10	0
1	EEE	464	0	492	9	0
1	FFF	464	0	492	23	0
2	AAA	7	0	0	0	0
2	BBB	3	0	0	0	0
2	CCC	4	0	0	0	0
2	DDD	1	0	0	0	0
2	EEE	1	0	0	0	0
2	FFF	1	0	0	0	0
All	All	2801	0	2952	74	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 13.

All (74) 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:CCC:45:ASN:ND2	1:CCC:51:TYR:H	1.81	0.79
1:FFF:5:LYS:HG2	1:FFF:11:TYR:HA	1.65	0.76
1:BBB:4:ASN:ND2	1:BBB:58:ARG:HE	1.86	0.73
1:CCC:31:THR:HG21	1:DDD:9:LEU:HD11	1.73	0.71
1:BBB:4:ASN:HD21	1:BBB:58:ARG:HE	1.38	0.70
1:CCC:45:ASN:HD21	1:CCC:51:TYR:H	1.41	0.69
1:CCC:29:ASN:HD22	1:CCC:30:LYS:N	1.91	0.68
1:FFF:25:PHE:CD1	1:FFF:29:ASN:HA	2.30	0.67
1:FFF:21:CYS:HB2	1:FFF:54:CYS:SG	2.36	0.66
1:AAA:4:ASN:HD21	1:AAA:58:ARG:HE	1.42	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:FFF:23:LYS:HB2	1:FFF:25:PHE:CE2	2.32	0.65
1:CCC:45:ASN:HD22	1:CCC:50:LYS:HA	1.59	0.65
1:DDD:29:ASN:C	1:DDD:29:ASN:HD22	2.04	0.61
1:DDD:43:PRO:HB2	1:DDD:51:TYR:CE2	2.36	0.61
1:AAA:40:ASP:OD2	1:AAA:41:VAL:N	2.34	0.60
1:DDD:43:PRO:HB2	1:DDD:51:TYR:CD2	2.37	0.60
1:AAA:26:MET:HE3	1:AAA:48:LEU:O	2.03	0.59
1:CCC:45:ASN:ND2	1:CCC:50:LYS:HA	2.18	0.59
1:EEE:29:ASN:ND2	1:EEE:32:VAL:O	2.35	0.59
1:CCC:29:ASN:HD22	1:CCC:29:ASN:C	2.06	0.58
1:FFF:25:PHE:CG	1:FFF:29:ASN:HA	2.40	0.57
1:DDD:45:ASN:ND2	1:DDD:51:TYR:H	2.03	0.56
1:EEE:40:ASP:OD2	1:EEE:40:ASP:N	2.23	0.53
1:FFF:57:ASP:OD1	1:FFF:57:ASP:N	2.41	0.53
1:FFF:32:VAL:HG23	1:FFF:33:PRO:HD2	1.90	0.53
1:EEE:2:LYS:HG3	1:EEE:13:THR:HG22	1.91	0.53
1:BBB:29:ASN:HD21	1:BBB:30:LYS:HE2	1.74	0.52
1:FFF:1:LEU:HD21	1:FFF:56:THR:O	2.10	0.52
1:FFF:26:MET:O	1:FFF:27:VAL:HB	2.11	0.51
1:FFF:26:MET:HA	1:FFF:26:MET:HE2	1.93	0.51
1:FFF:1:LEU:HD23	1:FFF:2:LYS:N	2.26	0.50
1:EEE:1:LEU:HD11	1:EEE:55:ASN:HA	1.93	0.50
1:FFF:2:LYS:HA	1:FFF:13:THR:HG22	1.92	0.50
1:BBB:46:SER:HB2	1:FFF:35:LYS:HE3	1.93	0.50
1:DDD:3:CYS:SG	1:DDD:14:CYS:SG	3.10	0.49
1:EEE:29:ASN:C	1:EEE:31:THR:H	2.16	0.49
1:BBB:46:SER:OG	1:BBB:47:LEU:N	2.47	0.48
1:CCC:7:ILE:O	1:CCC:9:LEU:N	2.46	0.48
1:AAA:26:MET:HE3	1:AAA:49:VAL:HG22	1.95	0.47
1:FFF:23:LYS:HB2	1:FFF:25:PHE:HE2	1.78	0.47
1:AAA:1:LEU:HD23	1:AAA:1:LEU:HA	1.82	0.47
1:DDD:4:ASN:HD21	1:DDD:58:ARG:NE	2.12	0.47
1:CCC:23:LYS:O	1:CCC:51:TYR:HA	2.15	0.46
1:EEE:2:LYS:CG	1:EEE:13:THR:HG22	2.45	0.46
1:FFF:1:LEU:HD12	1:FFF:19:ASN:HA	1.98	0.46
1:BBB:25:PHE:CE1	1:BBB:33:PRO:HG3	2.51	0.46
1:FFF:2:LYS:HG2	1:FFF:13:THR:CG2	2.46	0.46
1:BBB:26:MET:CE	1:BBB:26:MET:HA	2.46	0.44
1:FFF:9:LEU:HD23	1:FFF:9:LEU:HA	1.90	0.44
1:FFF:20:LEU:HD12	1:FFF:41:VAL:HA	1.99	0.44
1:BBB:29:ASN:C	1:BBB:29:ASN:HD22	2.21	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BBB:26:MET:CE	1:BBB:26:MET:CA	2.95	0.44
1:BBB:4:ASN:HD21	1:BBB:58:ARG:NE	2.10	0.43
1:CCC:29:ASN:C	1:CCC:29:ASN:ND2	2.70	0.42
1:BBB:56:THR:HG22	1:BBB:59:CYS:HB3	2.01	0.42
1:CCC:32:VAL:HA	1:CCC:33:PRO:HD3	1.75	0.42
1:CCC:27:VAL:HG12	1:CCC:28:SER:N	2.33	0.42
1:BBB:19:ASN:O	1:BBB:20:LEU:HD23	2.20	0.42
1:DDD:43:PRO:HG2	1:DDD:51:TYR:CE1	2.55	0.42
1:AAA:32:VAL:HG21	1:BBB:32:VAL:HG13	2.01	0.41
1:FFF:4:ASN:OD1	1:FFF:60:ASN:ND2	2.52	0.41
1:FFF:2:LYS:HG2	1:FFF:13:THR:HG22	2.03	0.41
1:FFF:5:LYS:HA	1:FFF:36:ARG:HB2	2.01	0.41
1:AAA:32:VAL:CG2	1:BBB:32:VAL:HG13	2.50	0.41
1:FFF:11:TYR:O	1:FFF:12:LYS:HB3	2.20	0.41
1:BBB:30:LYS:HB3	1:CCC:27:VAL:HG21	2.03	0.41
1:EEE:26:MET:N	1:EEE:26:MET:SD	2.93	0.41
1:BBB:45:ASN:ND2	1:BBB:51:TYR:H	2.19	0.40
1:EEE:11:TYR:CG	1:EEE:58:ARG:NH1	2.89	0.40
1:FFF:30:LYS:O	1:FFF:31:THR:O	2.39	0.40
1:CCC:21:CYS:HA	1:CCC:37:GLY:O	2.21	0.40
1:DDD:4:ASN:ND2	1:DDD:58:ARG:HG2	2.37	0.40
1:DDD:39:ILE:HG12	1:DDD:41:VAL:O	2.21	0.40
1:EEE:46:SER:OG	1:EEE:47:LEU:N	2.55	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	AAA	58/60 (97%)	55 (95%)	3 (5%)	0	100 100
1	BBB	58/60 (97%)	50 (86%)	6 (10%)	2 (3%)	3 5

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	CCC	58/60 (97%)	54 (93%)	4 (7%)	0	100	100
1	DDD	58/60 (97%)	52 (90%)	3 (5%)	3 (5%)	2	2
1	EEE	58/60 (97%)	48 (83%)	7 (12%)	3 (5%)	2	2
1	FFF	58/60 (97%)	47 (81%)	6 (10%)	5 (9%)	1	0
All	All	348/360 (97%)	306 (88%)	29 (8%)	13 (4%)	3	4

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	DDD	40	ASP
1	BBB	46	SER
1	DDD	56	THR
1	FFF	9	LEU
1	FFF	40	ASP
1	BBB	40	ASP
1	FFF	15	PRO
1	FFF	31	THR
1	DDD	27	VAL
1	EEE	9	LEU
1	EEE	28	SER
1	EEE	29	ASN
1	FFF	58	ARG

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	AAA	56/56 (100%)	43 (77%)	13 (23%)	1	1
1	BBB	56/56 (100%)	48 (86%)	8 (14%)	3	5
1	CCC	56/56 (100%)	47 (84%)	9 (16%)	2	4
1	DDD	56/56 (100%)	45 (80%)	11 (20%)	1	2
1	EEE	56/56 (100%)	43 (77%)	13 (23%)	1	1

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	FFF	56/56 (100%)	44 (79%)	12 (21%)	1	1
All	All	336/336 (100%)	270 (80%)	66 (20%)	1	2

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	1	LEU
1	AAA	5	LYS
1	AAA	7	ILE
1	AAA	9	LEU
1	AAA	12	LYS
1	AAA	13	THR
1	AAA	26	MET
1	AAA	29	ASN
1	AAA	32	VAL
1	AAA	34	VAL
1	AAA	42	CYS
1	AAA	44	LYS
1	AAA	47	LEU
1	BBB	1	LEU
1	BBB	26	MET
1	BBB	28	SER
1	BBB	29	ASN
1	BBB	34	VAL
1	BBB	40	ASP
1	BBB	41	VAL
1	BBB	56	THR
1	CCC	1	LEU
1	CCC	26	MET
1	CCC	29	ASN
1	CCC	30	LYS
1	CCC	32	VAL
1	CCC	34	VAL
1	CCC	36	ARG
1	CCC	44	LYS
1	CCC	56	THR
1	DDD	13	THR
1	DDD	29	ASN
1	DDD	34	VAL
1	DDD	39	ILE
1	DDD	40	ASP
1	DDD	41	VAL

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Mol	Chain	Res	Type
1	DDD	44	LYS
1	DDD	47	LEU
1	DDD	52	VAL
1	DDD	56	THR
1	DDD	57	ASP
1	EEE	24	MET
1	EEE	26	MET
1	EEE	28	SER
1	EEE	30	LYS
1	EEE	31	THR
1	EEE	32	VAL
1	EEE	34	VAL
1	EEE	40	ASP
1	EEE	42	CYS
1	EEE	45	ASN
1	EEE	46	SER
1	EEE	47	LEU
1	EEE	55	ASN
1	FFF	4	ASN
1	FFF	12	LYS
1	FFF	24	MET
1	FFF	30	LYS
1	FFF	32	VAL
1	FFF	34	VAL
1	FFF	42	CYS
1	FFF	44	LYS
1	FFF	47	LEU
1	FFF	52	VAL
1	FFF	57	ASP
1	FFF	59	CYS

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](#)

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	AAA	60/60 (100%)	-0.47	0 100 100	58, 72, 93, 100	0
1	BBB	60/60 (100%)	-0.41	0 100 100	59, 78, 109, 119	0
1	CCC	60/60 (100%)	-0.31	1 (1%) 70 66	68, 77, 108, 127	0
1	DDD	60/60 (100%)	0.31	6 (10%) 7 4	72, 95, 132, 141	0
1	EEE	60/60 (100%)	0.77	11 (18%) 1 0	77, 123, 158, 171	0
1	FFF	60/60 (100%)	1.02	13 (21%) 0 0	72, 149, 178, 194	0
All	All	360/360 (100%)	0.15	31 (8%) 10 7	58, 90, 164, 194	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	FFF	17	GLY	6.5
1	FFF	2	LYS	6.0
1	DDD	2	LYS	5.7
1	DDD	3	CYS	5.5
1	FFF	18	LYS	4.8
1	FFF	58	ARG	4.1
1	EEE	21	CYS	4.0
1	FFF	12	LYS	4.0
1	EEE	55	ASN	3.9
1	DDD	21	CYS	3.8
1	EEE	3	CYS	3.7
1	FFF	16	ALA	3.4
1	FFF	57	ASP	3.3
1	FFF	55	ASN	3.3
1	EEE	20	LEU	3.1
1	FFF	21	CYS	3.0
1	EEE	19	ASN	2.9
1	EEE	30	LYS	2.9
1	EEE	60	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	EEE	16	ALA	2.8
1	FFF	40	ASP	2.7
1	DDD	1	LEU	2.7
1	FFF	56	THR	2.5
1	FFF	44	LYS	2.4
1	DDD	58	ARG	2.4
1	DDD	56	THR	2.3
1	EEE	23	LYS	2.2
1	CCC	21	CYS	2.2
1	EEE	15	PRO	2.1
1	EEE	2	LYS	2.1
1	FFF	19	ASN	2.1

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.