



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 1, 2026 – 10:03 AM UTC

PDB ID : 4G0J / pdb_00004g0j
Title : Crystallographic Analysis of Rotavirus NSP2-RNA Complex Reveals Specific Recognition of 5'-GG Sequence for RTPase activity
Authors : Hu, L.; Prasad, B.V.V.
Deposited on : 2012-07-09
Resolution : 3.40 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

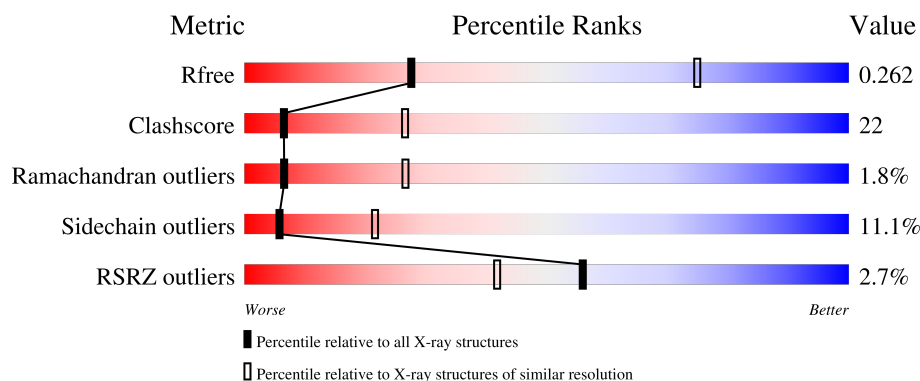
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	294	3% 47% 45% 7%
1	B	294	4% 55% 36% 7%
1	C	294	3% 50% 42% 6%
1	D	294	2% 55% 40%
1	E	294	3% 55% 37% 6%

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Mol	Chain	Length	Quality of chain
1	F	294	2% 54% 39% 5%
1	G	294	2% 47% 46% 5%
1	H	294	2% 61% 32% 6%
1	I	294	3% 52% 42%
1	J	294	4% 54% 40%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 23600 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.

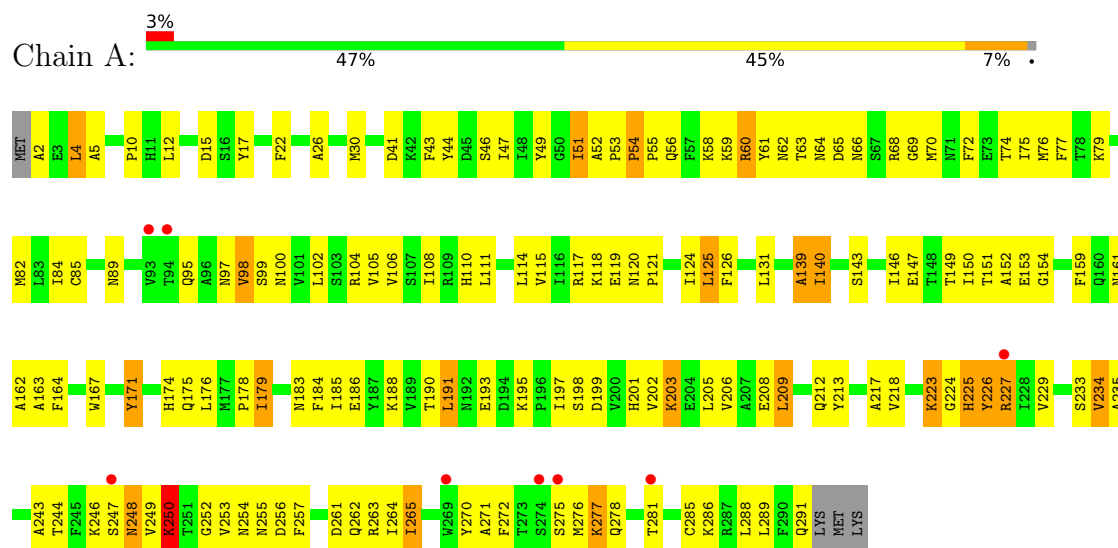
- Molecule 1 is a protein called Non-structural protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	290	Total	C	N	O	S	0	0	0
			2360	1518	400	430	12			
1	B	290	Total	C	N	O	S	0	0	0
			2360	1518	400	430	12			
1	C	290	Total	C	N	O	S	0	0	0
			2360	1518	400	430	12			
1	D	290	Total	C	N	O	S	0	0	0
			2360	1518	400	430	12			
1	E	290	Total	C	N	O	S	0	0	0
			2360	1518	400	430	12			
1	F	290	Total	C	N	O	S	0	0	0
			2360	1518	400	430	12			
1	G	290	Total	C	N	O	S	0	0	0
			2360	1518	400	430	12			
1	H	290	Total	C	N	O	S	0	0	0
			2360	1518	400	430	12			
1	I	290	Total	C	N	O	S	0	0	0
			2360	1518	400	430	12			
1	J	290	Total	C	N	O	S	0	0	0
			2360	1518	400	430	12			

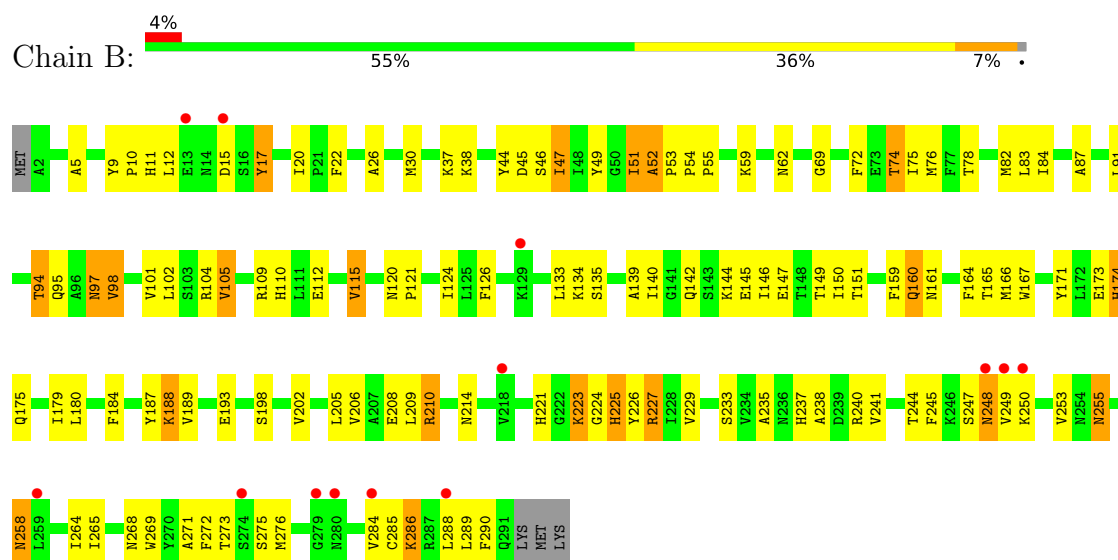
3 Residue-property plots

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: Non-structural protein 2

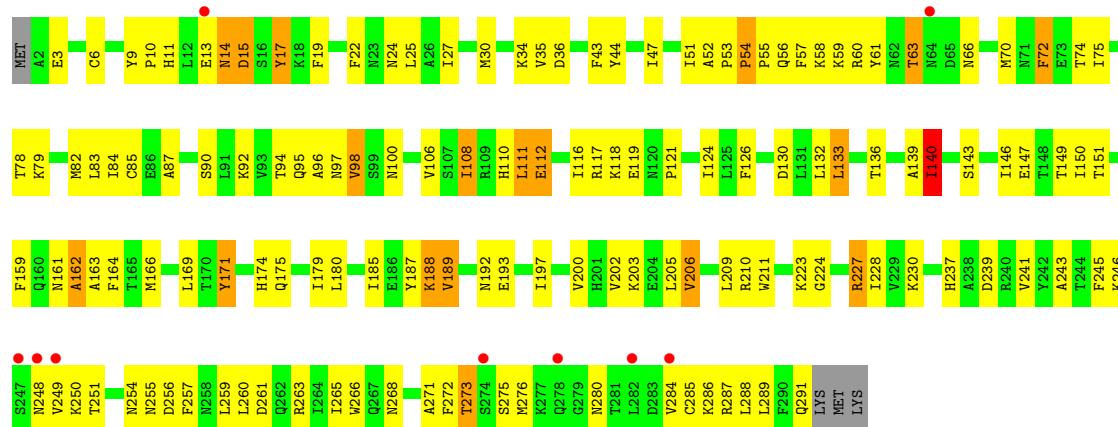


• Molecule 1: Non-structural protein 2

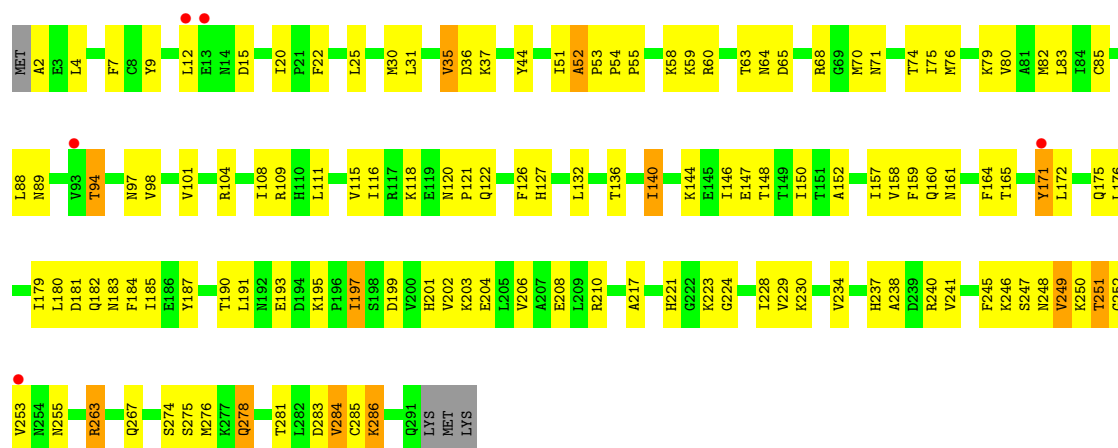


• Molecule 1: Non-structural protein 2

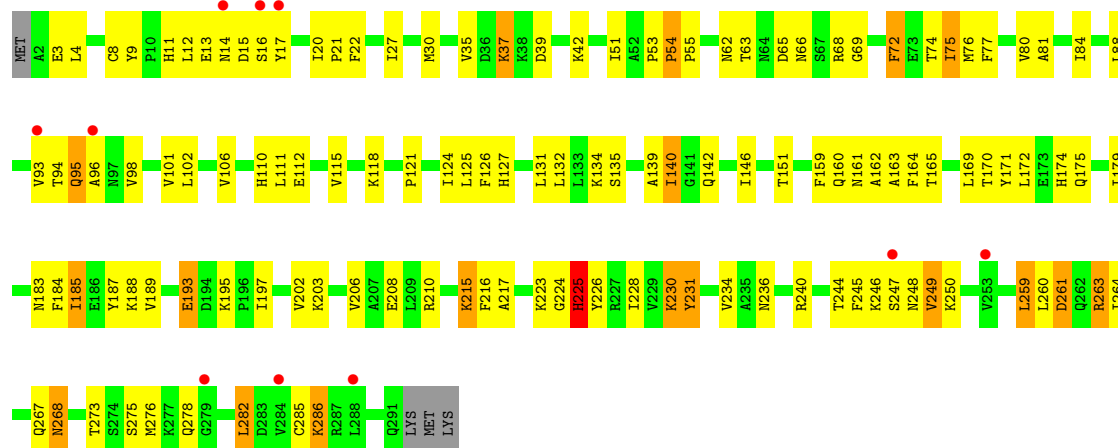




• Molecule 1: Non-structural protein 2



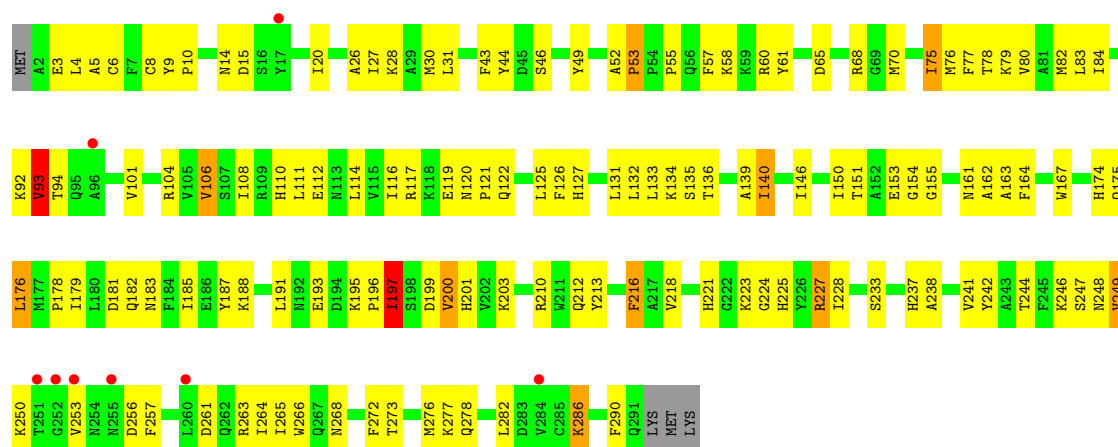
• Molecule 1: Non-structural protein 2



• Molecule 1: Non-structural protein 2

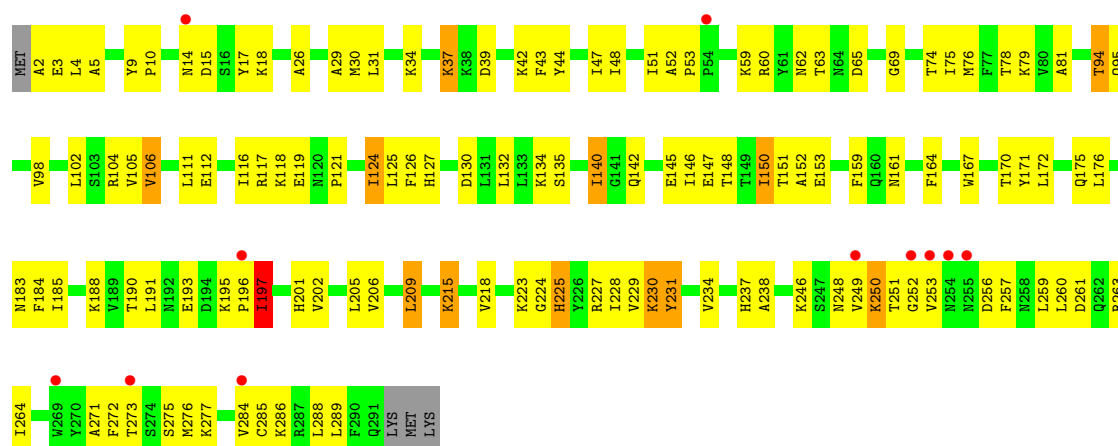
• Molecule 1: Non-structural protein 2

Chain I:  3% 52% 42% ..



• Molecule 1: Non-structural protein 2

Chain J:  4% 54% 40% ..



4 Data and refinement statistics

Property	Value	Source
Space group	P 4	Depositor
Cell constants a, b, c, α , β , γ	122.48Å 122.48Å 301.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	301.31 – 3.40 301.13 – 3.40	Depositor EDS
% Data completeness (in resolution range)	99.3 (301.31-3.40) 99.4 (301.13-3.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 3.44Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.237 , 0.294 0.211 , 0.262	Depositor DCC
R_{free} test set	3536 reflections (5.81%)	wwPDB-VP
Wilson B-factor (Å ²)	67.0	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 28.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.236 for h,-k,-l	Xtriage
Reported twinning fraction	0.299 for H, K, L 0.701 for -H, K, -L	Depositor
Outliers	0 of 60405 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	23600	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 33.80 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.5552e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	A	0.84	0/2411	1.14	10/3261 (0.3%)
1	B	0.86	0/2411	1.15	8/3261 (0.2%)
1	C	0.88	0/2411	1.16	12/3261 (0.4%)
1	D	0.86	0/2411	1.11	9/3261 (0.3%)
1	E	0.90	2/2411 (0.1%)	1.10	7/3261 (0.2%)
1	F	0.88	2/2411 (0.1%)	1.09	7/3261 (0.2%)
1	G	0.86	1/2411 (0.0%)	1.14	9/3261 (0.3%)
1	H	0.87	0/2411	1.16	14/3261 (0.4%)
1	I	0.88	1/2411 (0.0%)	1.14	6/3261 (0.2%)
1	J	0.85	0/2411	1.07	7/3261 (0.2%)
All	All	0.87	6/24110 (0.0%)	1.13	89/32610 (0.3%)

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	95	GLN	CD-NE2	6.43	1.46	1.33
1	F	158	VAL	N-CA	5.50	1.52	1.46
1	E	225	HIS	CG-CD2	5.17	1.41	1.35
1	I	53	PRO	CA-C	5.16	1.54	1.51
1	F	11	HIS	CG-CD2	5.14	1.41	1.35

The worst 5 of 89 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	195	LYS	CA-C-N	10.74	130.86	119.90
1	A	195	LYS	C-N-CA	10.74	130.86	119.90
1	D	252	GLY	N-CA-C	-9.64	103.27	113.58
1	J	252	GLY	N-CA-C	-9.26	102.71	114.37
1	H	252	GLY	N-CA-C	-9.25	103.18	115.21

There are no chirality outliers.

There are no planarity outliers.

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	A	2360	0	2374	135	0
1	B	2360	0	2374	97	0
1	C	2360	0	2374	101	1
1	D	2360	0	2374	96	0
1	E	2360	0	2374	104	0
1	F	2360	0	2374	94	1
1	G	2360	0	2374	141	0
1	H	2360	0	2374	84	0
1	I	2360	0	2374	120	0
1	J	2360	0	2374	112	0
All	All	23600	0	23740	1032	1

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 22.

The worst 5 of 1032 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:H:74:THR:HB	1:H:175:GLN:HE21	1.12	1.14
1:I:65:ASP:HA	1:J:183:ASN:HD21	1.04	1.11
1:H:202:VAL:O	1:H:206:VAL:HG23	1.52	1.09
1:E:74:THR:HB	1:E:175:GLN:HE21	1.10	1.08
1:E:202:VAL:HG13	1:E:264:ILE:HD11	1.35	1.07

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:239:ASP:OD2	1:F:251:THR:OG1[2_654]	2.14	0.06

5.3 Torsion angles

5.3.1 Protein backbone

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	288/294 (98%)	246 (85%)	35 (12%)	7 (2%)	4	22
1	B	288/294 (98%)	246 (85%)	36 (12%)	6 (2%)	5	24
1	C	288/294 (98%)	243 (84%)	41 (14%)	4 (1%)	9	31
1	D	288/294 (98%)	256 (89%)	29 (10%)	3 (1%)	12	40
1	E	288/294 (98%)	260 (90%)	22 (8%)	6 (2%)	5	24
1	F	288/294 (98%)	252 (88%)	27 (9%)	9 (3%)	3	18
1	G	288/294 (98%)	252 (88%)	32 (11%)	4 (1%)	9	31
1	H	288/294 (98%)	254 (88%)	28 (10%)	6 (2%)	5	24
1	I	288/294 (98%)	257 (89%)	28 (10%)	3 (1%)	12	40
1	J	288/294 (98%)	256 (89%)	28 (10%)	4 (1%)	9	31
All	All	2880/2940 (98%)	2522 (88%)	306 (11%)	52 (2%)	6	26

5 of 52 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	249	VAL
1	E	249	VAL
1	F	235	ALA
1	F	250	LYS
1	G	249	VAL

5.3.2 Protein sidechains

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	262/266 (98%)	229 (87%)	33 (13%)	4	18
1	B	262/266 (98%)	227 (87%)	35 (13%)	4	15
1	C	262/266 (98%)	225 (86%)	37 (14%)	3	13
1	D	262/266 (98%)	232 (88%)	30 (12%)	5	20
1	E	262/266 (98%)	237 (90%)	25 (10%)	8	29
1	F	262/266 (98%)	233 (89%)	29 (11%)	6	21
1	G	262/266 (98%)	237 (90%)	25 (10%)	8	29
1	H	262/266 (98%)	233 (89%)	29 (11%)	6	21
1	I	262/266 (98%)	239 (91%)	23 (9%)	9	31
1	J	262/266 (98%)	236 (90%)	26 (10%)	7	26
All	All	2620/2660 (98%)	2328 (89%)	292 (11%)	6	21

5 of 292 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	188	LYS
1	J	215	LYS
1	H	251	THR
1	I	227	ARG
1	C	227	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 90 such sidechains are listed below:

Mol	Chain	Res	Type
1	G	214	ASN
1	I	127	HIS
1	H	110	HIS
1	H	214	ASN
1	I	175	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	290/294 (98%)	0.16	8 (2%) 55 41	37, 73, 144, 188	0
1	B	290/294 (98%)	0.04	13 (4%) 38 28	43, 72, 141, 183	0
1	C	290/294 (98%)	0.02	9 (3%) 51 38	40, 70, 143, 180	0
1	D	290/294 (98%)	-0.07	5 (1%) 69 54	41, 71, 123, 182	0
1	E	290/294 (98%)	-0.04	10 (3%) 48 35	37, 66, 136, 212	0
1	F	290/294 (98%)	0.02	5 (1%) 69 54	43, 73, 124, 206	0
1	G	290/294 (98%)	0.05	5 (1%) 69 54	41, 71, 137, 196	0
1	H	290/294 (98%)	-0.11	5 (1%) 69 54	38, 70, 122, 205	0
1	I	290/294 (98%)	-0.05	8 (2%) 55 41	39, 67, 131, 224	0
1	J	290/294 (98%)	0.03	11 (3%) 44 31	41, 71, 140, 240	0
All	All	2900/2940 (98%)	0.00	79 (2%) 56 42	37, 71, 138, 240	0

The worst 5 of 79 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	17	TYR	6.2
1	J	253	VAL	5.6
1	I	17	TYR	5.6
1	G	238	ALA	4.5
1	E	279	GLY	4.4

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.