



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 06:42 PM UTC

PDB ID : 8EAW / pdb_00008eaw
Title : An asymmetric disk assembly formed by tandem dimers of the tobacco mosaic viral capsid protein (TMV)
Authors : Dai, J.; Pereira, J.H.; Adams, P.D.; Francis, M.B.
Deposited on : 2022-08-29
Resolution : 2.80 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

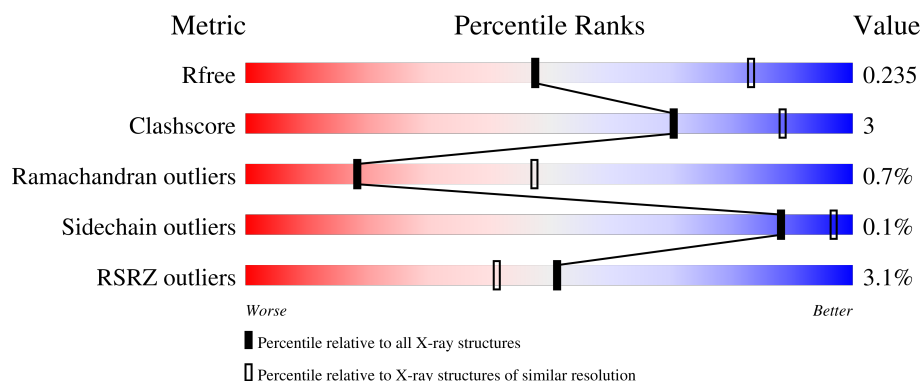
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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.



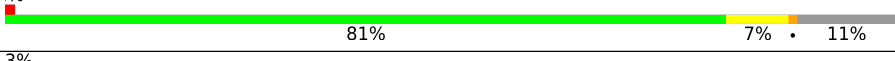
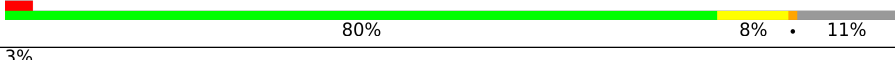
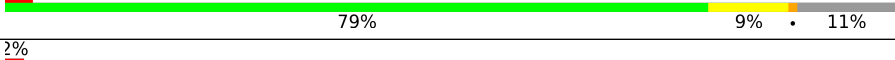
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	323	2% 80% 9% 11%
1	b	323	2% 79% 10% 11%
1	c	323	3% 81% 7% 11%
1	d	323	3% 79% 9% 11%
1	e	323	4% 81% 7% 11%

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Mol	Chain	Length	Quality of chain
1	f	323	
1	g	323	
1	h	323	
1	i	323	
1	j	323	
1	k	323	
1	l	323	
1	m	323	
1	n	323	
1	o	323	
1	p	323	
1	q	323	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 38318 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 Capsid protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	a	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			
1	b	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			
1	c	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			
1	d	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			
1	e	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			
1	f	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			
1	g	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			
1	h	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			
1	i	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			
1	j	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			
1	k	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			
1	l	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			
1	m	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			
1	n	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			
1	o	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			
1	p	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	q	287	Total	C	N	O	S	0	0	0
			2254	1422	392	436	4			

There are 153 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
a	123	CYS	SER	engineered mutation	UNP P69687
a	159	GLY	-	linker	UNP P69687
a	160	GLY	-	linker	UNP P69687
a	161	GLY	-	linker	UNP P69687
a	162	GLU	-	linker	UNP P69687
a	163	GLY	-	linker	UNP P69687
a	164	GLY	-	linker	UNP P69687
a	165	GLY	-	linker	UNP P69687
a	288	CYS	SER	engineered mutation	UNP P69687
b	123	CYS	SER	engineered mutation	UNP P69687
b	159	GLY	-	linker	UNP P69687
b	160	GLY	-	linker	UNP P69687
b	161	GLY	-	linker	UNP P69687
b	162	GLU	-	linker	UNP P69687
b	163	GLY	-	linker	UNP P69687
b	164	GLY	-	linker	UNP P69687
b	165	GLY	-	linker	UNP P69687
b	288	CYS	SER	engineered mutation	UNP P69687
c	123	CYS	SER	engineered mutation	UNP P69687
c	159	GLY	-	linker	UNP P69687
c	160	GLY	-	linker	UNP P69687
c	161	GLY	-	linker	UNP P69687
c	162	GLU	-	linker	UNP P69687
c	163	GLY	-	linker	UNP P69687
c	164	GLY	-	linker	UNP P69687
c	165	GLY	-	linker	UNP P69687
c	288	CYS	SER	engineered mutation	UNP P69687
d	123	CYS	SER	engineered mutation	UNP P69687
d	159	GLY	-	linker	UNP P69687
d	160	GLY	-	linker	UNP P69687
d	161	GLY	-	linker	UNP P69687
d	162	GLU	-	linker	UNP P69687
d	163	GLY	-	linker	UNP P69687
d	164	GLY	-	linker	UNP P69687
d	165	GLY	-	linker	UNP P69687
d	288	CYS	SER	engineered mutation	UNP P69687

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Chain	Residue	Modelled	Actual	Comment	Reference
e	123	CYS	SER	engineered mutation	UNP P69687
e	159	GLY	-	linker	UNP P69687
e	160	GLY	-	linker	UNP P69687
e	161	GLY	-	linker	UNP P69687
e	162	GLU	-	linker	UNP P69687
e	163	GLY	-	linker	UNP P69687
e	164	GLY	-	linker	UNP P69687
e	165	GLY	-	linker	UNP P69687
e	288	CYS	SER	engineered mutation	UNP P69687
f	123	CYS	SER	engineered mutation	UNP P69687
f	159	GLY	-	linker	UNP P69687
f	160	GLY	-	linker	UNP P69687
f	161	GLY	-	linker	UNP P69687
f	162	GLU	-	linker	UNP P69687
f	163	GLY	-	linker	UNP P69687
f	164	GLY	-	linker	UNP P69687
f	165	GLY	-	linker	UNP P69687
f	288	CYS	SER	engineered mutation	UNP P69687
g	123	CYS	SER	engineered mutation	UNP P69687
g	159	GLY	-	linker	UNP P69687
g	160	GLY	-	linker	UNP P69687
g	161	GLY	-	linker	UNP P69687
g	162	GLU	-	linker	UNP P69687
g	163	GLY	-	linker	UNP P69687
g	164	GLY	-	linker	UNP P69687
g	165	GLY	-	linker	UNP P69687
g	288	CYS	SER	engineered mutation	UNP P69687
h	123	CYS	SER	engineered mutation	UNP P69687
h	159	GLY	-	linker	UNP P69687
h	160	GLY	-	linker	UNP P69687
h	161	GLY	-	linker	UNP P69687
h	162	GLU	-	linker	UNP P69687
h	163	GLY	-	linker	UNP P69687
h	164	GLY	-	linker	UNP P69687
h	165	GLY	-	linker	UNP P69687
h	288	CYS	SER	engineered mutation	UNP P69687
i	123	CYS	SER	engineered mutation	UNP P69687
i	159	GLY	-	linker	UNP P69687
i	160	GLY	-	linker	UNP P69687
i	161	GLY	-	linker	UNP P69687
i	162	GLU	-	linker	UNP P69687
i	163	GLY	-	linker	UNP P69687

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Chain	Residue	Modelled	Actual	Comment	Reference
i	164	GLY	-	linker	UNP P69687
i	165	GLY	-	linker	UNP P69687
i	288	CYS	SER	engineered mutation	UNP P69687
j	123	CYS	SER	engineered mutation	UNP P69687
j	159	GLY	-	linker	UNP P69687
j	160	GLY	-	linker	UNP P69687
j	161	GLY	-	linker	UNP P69687
j	162	GLU	-	linker	UNP P69687
j	163	GLY	-	linker	UNP P69687
j	164	GLY	-	linker	UNP P69687
j	165	GLY	-	linker	UNP P69687
j	288	CYS	SER	engineered mutation	UNP P69687
k	123	CYS	SER	engineered mutation	UNP P69687
k	159	GLY	-	linker	UNP P69687
k	160	GLY	-	linker	UNP P69687
k	161	GLY	-	linker	UNP P69687
k	162	GLU	-	linker	UNP P69687
k	163	GLY	-	linker	UNP P69687
k	164	GLY	-	linker	UNP P69687
k	165	GLY	-	linker	UNP P69687
k	288	CYS	SER	engineered mutation	UNP P69687
l	123	CYS	SER	engineered mutation	UNP P69687
l	159	GLY	-	linker	UNP P69687
l	160	GLY	-	linker	UNP P69687
l	161	GLY	-	linker	UNP P69687
l	162	GLU	-	linker	UNP P69687
l	163	GLY	-	linker	UNP P69687
l	164	GLY	-	linker	UNP P69687
l	165	GLY	-	linker	UNP P69687
l	288	CYS	SER	engineered mutation	UNP P69687
m	123	CYS	SER	engineered mutation	UNP P69687
m	159	GLY	-	linker	UNP P69687
m	160	GLY	-	linker	UNP P69687
m	161	GLY	-	linker	UNP P69687
m	162	GLU	-	linker	UNP P69687
m	163	GLY	-	linker	UNP P69687
m	164	GLY	-	linker	UNP P69687
m	165	GLY	-	linker	UNP P69687
m	288	CYS	SER	engineered mutation	UNP P69687
n	123	CYS	SER	engineered mutation	UNP P69687
n	159	GLY	-	linker	UNP P69687
n	160	GLY	-	linker	UNP P69687

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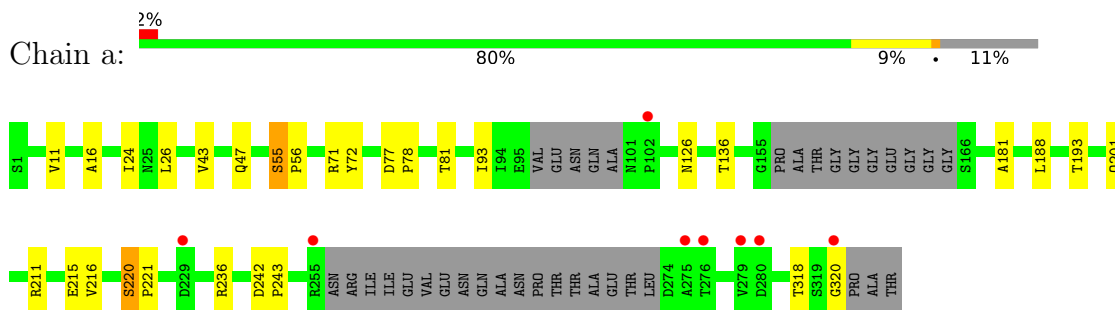
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Chain	Residue	Modelled	Actual	Comment	Reference
n	161	GLY	-	linker	UNP P69687
n	162	GLU	-	linker	UNP P69687
n	163	GLY	-	linker	UNP P69687
n	164	GLY	-	linker	UNP P69687
n	165	GLY	-	linker	UNP P69687
n	288	CYS	SER	engineered mutation	UNP P69687
o	123	CYS	SER	engineered mutation	UNP P69687
o	159	GLY	-	linker	UNP P69687
o	160	GLY	-	linker	UNP P69687
o	161	GLY	-	linker	UNP P69687
o	162	GLU	-	linker	UNP P69687
o	163	GLY	-	linker	UNP P69687
o	164	GLY	-	linker	UNP P69687
o	165	GLY	-	linker	UNP P69687
o	288	CYS	SER	engineered mutation	UNP P69687
p	123	CYS	SER	engineered mutation	UNP P69687
p	159	GLY	-	linker	UNP P69687
p	160	GLY	-	linker	UNP P69687
p	161	GLY	-	linker	UNP P69687
p	162	GLU	-	linker	UNP P69687
p	163	GLY	-	linker	UNP P69687
p	164	GLY	-	linker	UNP P69687
p	165	GLY	-	linker	UNP P69687
p	288	CYS	SER	engineered mutation	UNP P69687
q	123	CYS	SER	engineered mutation	UNP P69687
q	159	GLY	-	linker	UNP P69687
q	160	GLY	-	linker	UNP P69687
q	161	GLY	-	linker	UNP P69687
q	162	GLU	-	linker	UNP P69687
q	163	GLY	-	linker	UNP P69687
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q	288	CYS	SER	engineered mutation	UNP P69687

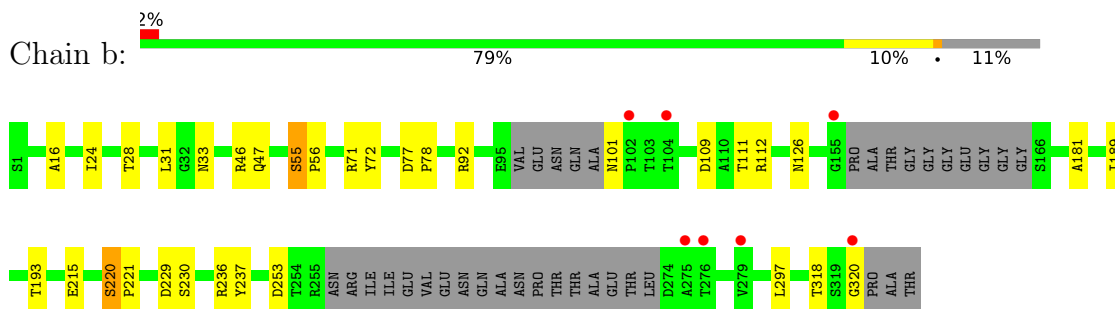
3 Residue-property plots [i](#)

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.

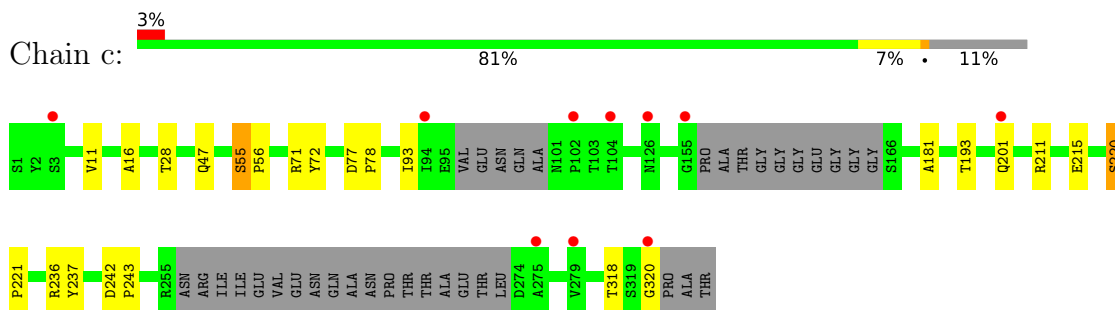
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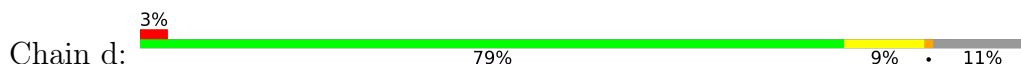
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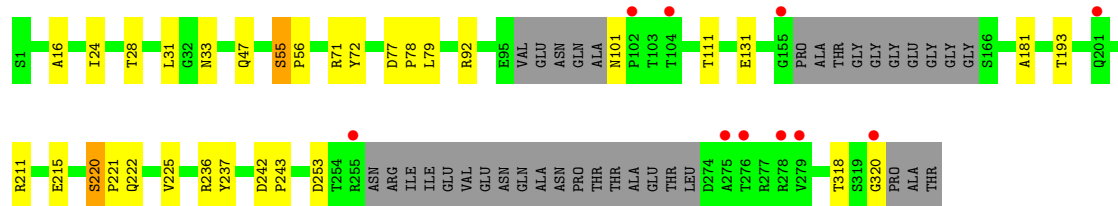


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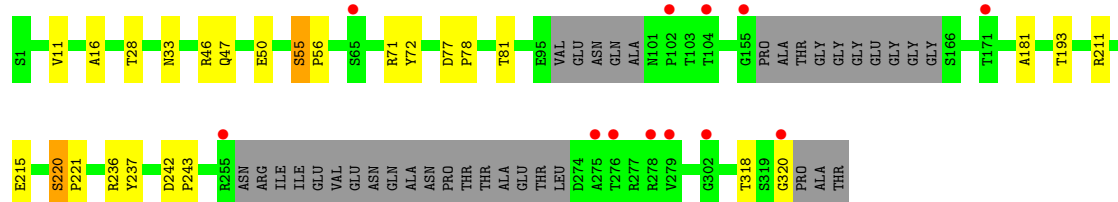
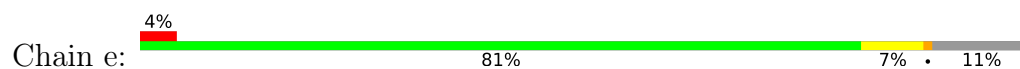


- Molecule 1: Capsid protein

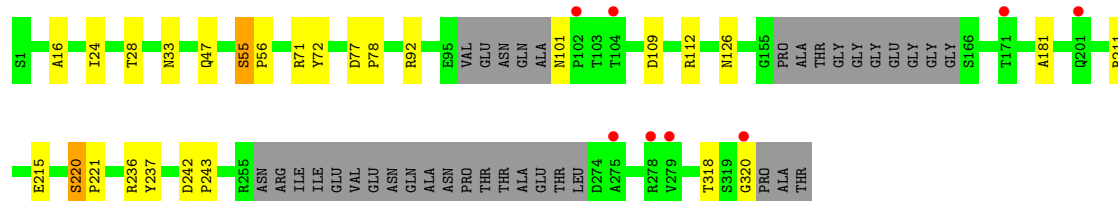
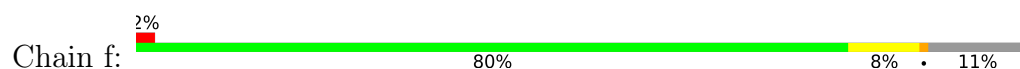




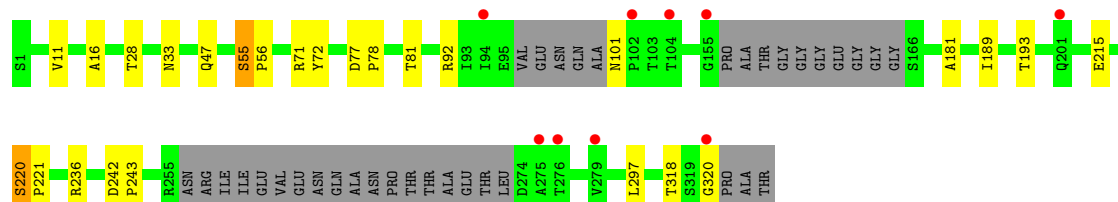
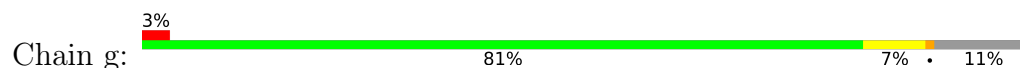
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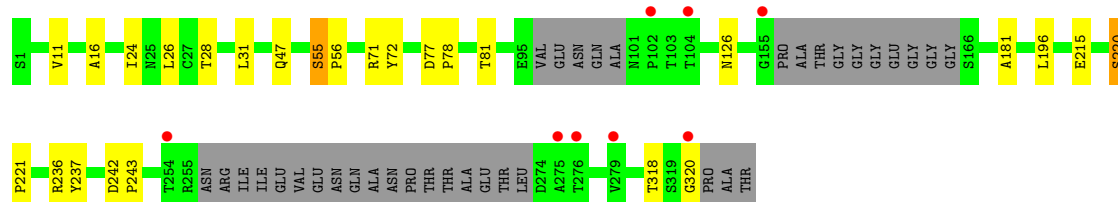
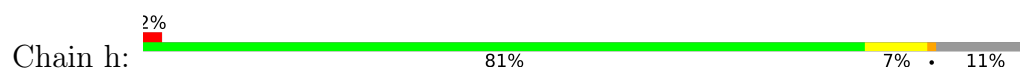
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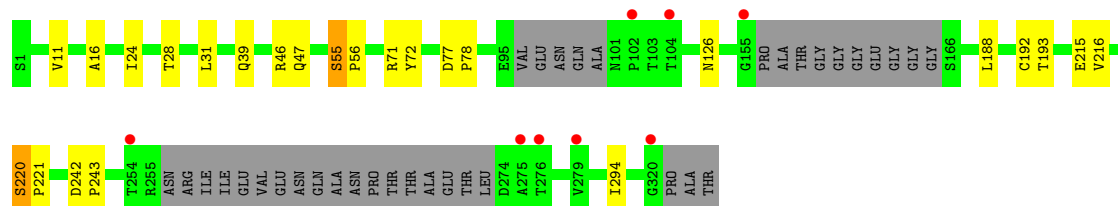
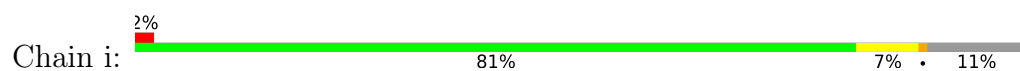
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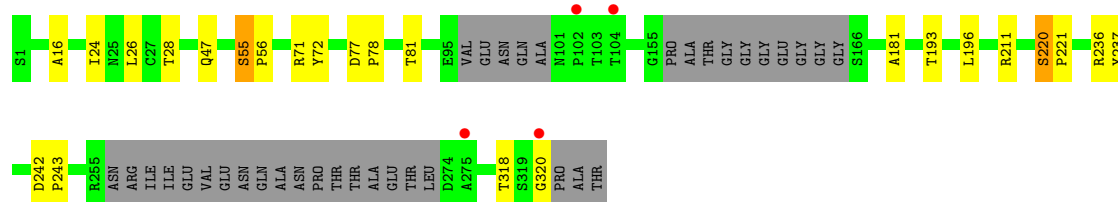
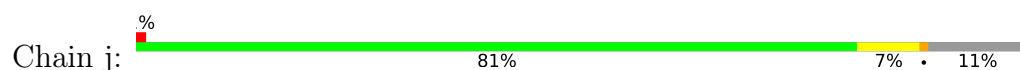
• Molecule 1: Capsid protein



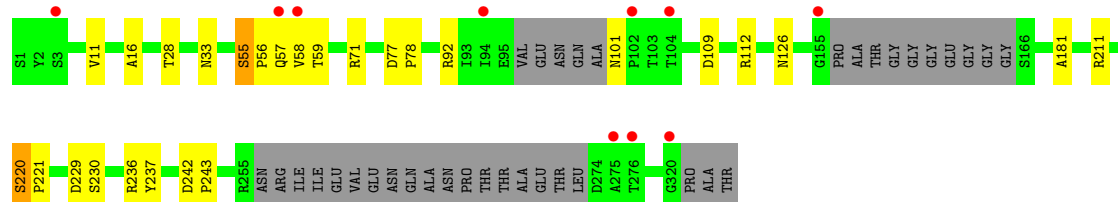
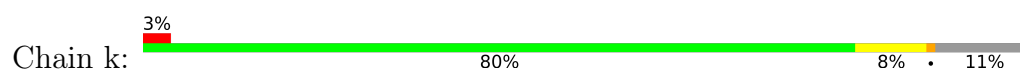
- Molecule 1: Capsid protein



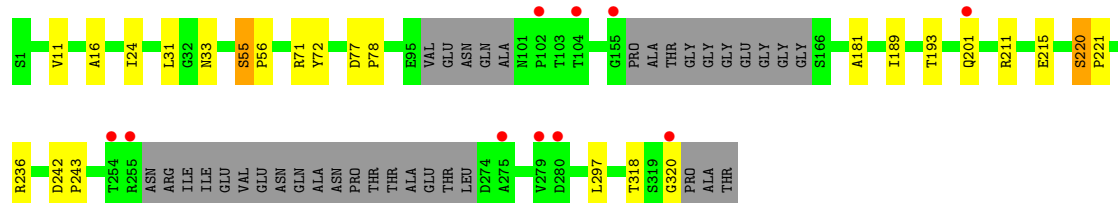
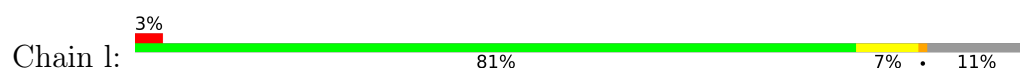
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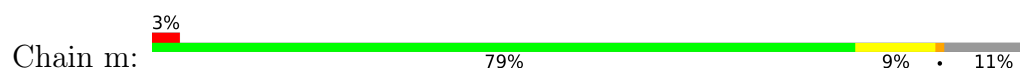
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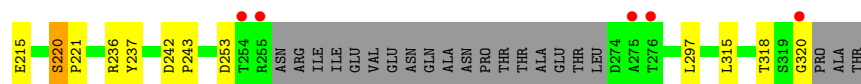


- Molecule 1: Capsid protein

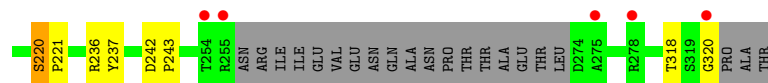
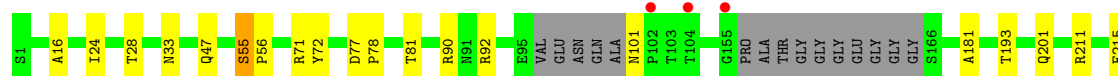
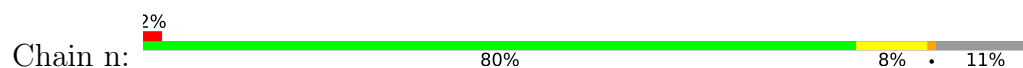


- Molecule 1: Capsid protein

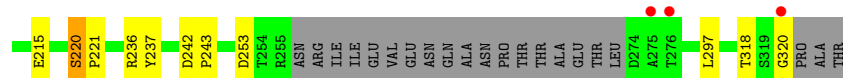
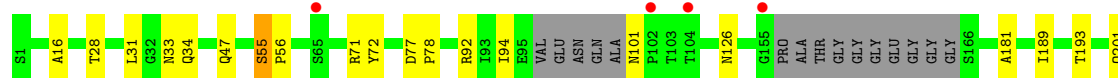
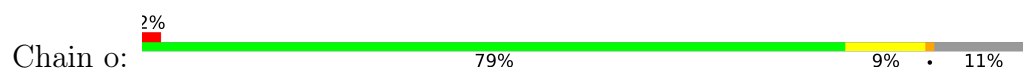




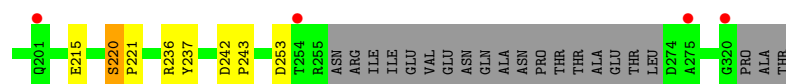
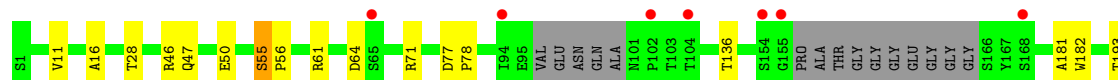
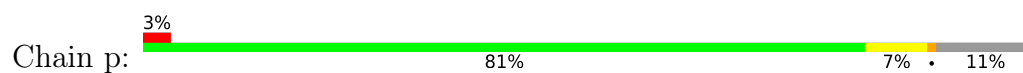
• Molecule 1: Capsid protein



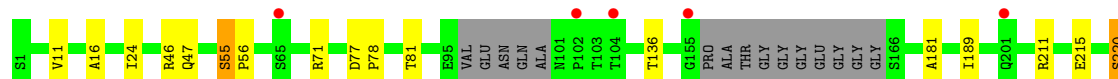
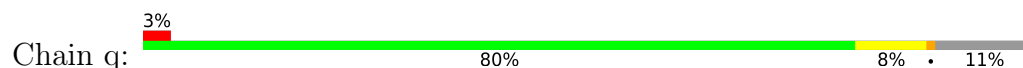
• Molecule 1: Capsid protein



• Molecule 1: Capsid protein



• Molecule 1: Capsid protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	208.51Å 255.50Å 260.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	81.44 – 2.80 81.44 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.7 (81.44-2.80) 99.7 (81.44-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.34 (at 2.82Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.196 , 0.234 0.197 , 0.235	Depositor DCC
R_{free} test set	8369 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	58.8	Xtriage
Anisotropy	0.275	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	38318	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.95% 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	a	0.14	0/2299	0.31	0/3136
1	b	0.13	0/2299	0.32	0/3136
1	c	0.13	0/2299	0.31	0/3136
1	d	0.15	0/2299	0.31	0/3136
1	e	0.13	0/2299	0.30	0/3136
1	f	0.13	0/2299	0.30	0/3136
1	g	0.13	0/2299	0.31	0/3136
1	h	0.13	0/2299	0.30	0/3136
1	i	0.13	0/2299	0.30	0/3136
1	j	0.13	0/2299	0.30	0/3136
1	k	0.14	0/2299	0.30	0/3136
1	l	0.13	0/2299	0.31	0/3136
1	m	0.13	0/2299	0.31	0/3136
1	n	0.13	0/2299	0.31	0/3136
1	o	0.13	0/2299	0.31	0/3136
1	p	0.14	0/2299	0.32	0/3136
1	q	0.14	0/2299	0.32	0/3136
All	All	0.13	0/39083	0.31	0/53312

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	2254	0	2212	23	0
1	b	2254	0	2212	27	0
1	c	2254	0	2212	19	0
1	d	2254	0	2212	24	0
1	e	2254	0	2212	21	0
1	f	2254	0	2212	20	0
1	g	2254	0	2212	18	0
1	h	2254	0	2212	19	0
1	i	2254	0	2212	18	0
1	j	2254	0	2212	20	0
1	k	2254	0	2212	19	0
1	l	2254	0	2212	17	0
1	m	2254	0	2212	25	0
1	n	2254	0	2212	24	0
1	o	2254	0	2212	24	0
1	p	2254	0	2212	16	1
1	q	2254	0	2212	18	0
All	All	38318	0	37604	258	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 3.

All (258) 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:b:72:TYR:HH	1:c:28:THR:HG1	1.21	0.83
1:o:72:TYR:HH	1:p:28:THR:HG1	1.33	0.76
1:d:16:ALA:HB1	1:d:71:ARG:HB3	1.71	0.73
1:m:193:THR:HG1	1:n:237:TYR:HH	1.35	0.73
1:o:92:ARG:HH12	1:o:101:ASN:HB3	1.55	0.72
1:a:193:THR:HG1	1:b:237:TYR:HH	1.39	0.71
1:e:16:ALA:HB1	1:e:71:ARG:HB3	1.73	0.70
1:m:72:TYR:HH	1:n:28:THR:HG1	1.38	0.69
1:l:16:ALA:HB1	1:l:71:ARG:HB3	1.75	0.68
1:i:193:THR:HG1	1:j:237:TYR:HH	1.43	0.67
1:o:16:ALA:HB1	1:o:71:ARG:HB3	1.76	0.66
1:g:16:ALA:HB1	1:g:71:ARG:HB3	1.77	0.66
1:f:16:ALA:HB1	1:f:71:ARG:HB3	1.76	0.66
1:a:16:ALA:HB1	1:a:71:ARG:HB3	1.78	0.66
1:i:16:ALA:HB1	1:i:71:ARG:HB3	1.78	0.65
1:c:16:ALA:HB1	1:c:71:ARG:HB3	1.79	0.65
1:a:211:ARG:NH2	1:b:33:ASN:OD1	2.28	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:m:16:ALA:HB1	1:m:71:ARG:HB3	1.80	0.64
1:n:16:ALA:HB1	1:n:71:ARG:HB3	1.78	0.64
1:m:92:ARG:HH12	1:m:101:ASN:HB3	1.62	0.64
1:e:193:THR:HG1	1:f:237:TYR:HH	1.46	0.64
1:p:16:ALA:HB1	1:p:71:ARG:HB3	1.80	0.64
1:d:72:TYR:HH	1:e:28:THR:HG1	1.42	0.64
1:p:193:THR:OG1	1:q:237:TYR:OH	2.15	0.64
1:j:193:THR:HG1	1:k:237:TYR:HH	1.46	0.63
1:f:211:ARG:NH2	1:g:33:ASN:OD1	2.24	0.63
1:e:72:TYR:HH	1:f:28:THR:HG1	1.47	0.62
1:m:211:ARG:NH2	1:n:33:ASN:OD1	2.28	0.62
1:k:16:ALA:HB1	1:k:71:ARG:HB3	1.81	0.62
1:l:193:THR:HG1	1:m:237:TYR:HH	1.45	0.61
1:n:193:THR:HG1	1:o:237:TYR:HH	1.47	0.60
1:d:211:ARG:NH2	1:e:33:ASN:OD1	2.21	0.60
1:b:16:ALA:HB1	1:b:71:ARG:HB3	1.83	0.60
1:d:215:GLU:OE1	1:e:47:GLN:NE2	2.36	0.59
1:e:72:TYR:OH	1:f:28:THR:OG1	2.21	0.59
1:j:16:ALA:HB1	1:j:71:ARG:HB3	1.83	0.59
1:h:16:ALA:HB1	1:h:71:ARG:HB3	1.85	0.58
1:o:193:THR:HG1	1:p:237:TYR:HH	1.51	0.58
1:a:215:GLU:OE1	1:b:47:GLN:NE2	2.36	0.58
1:e:193:THR:OG1	1:f:237:TYR:OH	2.16	0.58
1:b:181:ALA:HB1	1:b:236:ARG:HB3	1.86	0.58
1:j:72:TYR:HH	1:k:28:THR:HG1	1.49	0.58
1:j:181:ALA:HB1	1:j:236:ARG:HB3	1.84	0.58
1:l:72:TYR:OH	1:m:28:THR:OG1	2.21	0.57
1:a:72:TYR:HH	1:b:28:THR:HG1	1.50	0.57
1:a:72:TYR:OH	1:b:28:THR:OG1	2.23	0.56
1:b:193:THR:OG1	1:c:237:TYR:OH	2.21	0.56
1:l:55:SER:HB3	1:l:56:PRO:HD3	1.87	0.56
1:j:81:THR:HG21	1:k:126:ASN:HD21	1.71	0.56
1:f:318:THR:HG22	1:f:320:GLY:H	1.70	0.56
1:b:193:THR:HG1	1:c:237:TYR:HH	1.51	0.56
1:f:72:TYR:HH	1:g:28:THR:HG1	1.51	0.55
1:f:215:GLU:OE1	1:g:47:GLN:NE2	2.39	0.55
1:j:72:TYR:OH	1:k:28:THR:OG1	2.22	0.55
1:m:193:THR:OG1	1:n:237:TYR:OH	2.15	0.55
1:p:193:THR:HG1	1:q:237:TYR:HH	1.50	0.55
1:o:181:ALA:HB1	1:o:236:ARG:HB3	1.87	0.55
1:h:220:SER:HB3	1:h:221:PRO:HD3	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:e:181:ALA:HB1	1:e:236:ARG:HB3	1.90	0.54
1:k:55:SER:HB3	1:k:56:PRO:HD3	1.89	0.54
1:n:72:TYR:OH	1:o:28:THR:OG1	2.19	0.54
1:b:55:SER:HB3	1:b:56:PRO:HD3	1.90	0.54
1:n:211:ARG:NH2	1:o:33:ASN:OD1	2.28	0.54
1:q:55:SER:HB3	1:q:56:PRO:HD3	1.88	0.54
1:c:55:SER:HB3	1:c:56:PRO:HD3	1.90	0.54
1:i:215:GLU:OE1	1:j:47:GLN:NE2	2.41	0.54
1:q:16:ALA:HB1	1:q:71:ARG:HB3	1.88	0.54
1:e:11:VAL:HG23	1:f:24:ILE:HG12	1.90	0.54
1:m:181:ALA:HB1	1:m:236:ARG:HB3	1.89	0.54
1:o:92:ARG:NH1	1:o:101:ASN:HB3	2.20	0.54
1:g:55:SER:HB3	1:g:56:PRO:HD3	1.90	0.53
1:n:72:TYR:HH	1:o:28:THR:HG1	1.47	0.53
1:m:55:SER:HB3	1:m:56:PRO:HD3	1.89	0.53
1:a:220:SER:HB3	1:a:221:PRO:HD3	1.89	0.53
1:b:220:SER:HB3	1:b:221:PRO:HD3	1.90	0.53
1:h:181:ALA:HB1	1:h:236:ARG:HB3	1.90	0.53
1:j:55:SER:HB3	1:j:56:PRO:HD3	1.90	0.53
1:o:220:SER:HB3	1:o:221:PRO:HD3	1.90	0.53
1:f:220:SER:HB3	1:f:221:PRO:HD3	1.91	0.53
1:i:220:SER:HB3	1:i:221:PRO:HD3	1.91	0.52
1:c:193:THR:HG1	1:d:237:TYR:HH	1.54	0.52
1:i:55:SER:HB3	1:i:56:PRO:HD3	1.92	0.52
1:l:181:ALA:HB1	1:l:236:ARG:HB3	1.90	0.52
1:d:55:SER:HB3	1:d:56:PRO:HD3	1.92	0.51
1:h:55:SER:HB3	1:h:56:PRO:HD3	1.92	0.51
1:n:55:SER:HB3	1:n:56:PRO:HD3	1.91	0.51
1:n:220:SER:HB3	1:n:221:PRO:HD3	1.92	0.51
1:p:11:VAL:HG23	1:q:24:ILE:HG12	1.93	0.51
1:f:55:SER:HB3	1:f:56:PRO:HD3	1.93	0.51
1:f:181:ALA:HB1	1:f:236:ARG:HB3	1.92	0.51
1:c:11:VAL:HG23	1:d:24:ILE:HG12	1.92	0.51
1:j:318:THR:HG22	1:j:320:GLY:H	1.75	0.51
1:n:81:THR:HG21	1:o:126:ASN:HD21	1.76	0.51
1:g:77:ASP:HB3	1:g:78:PRO:HD3	1.93	0.51
1:a:81:THR:HG21	1:b:126:ASN:HD21	1.76	0.50
1:d:72:TYR:OH	1:e:28:THR:OG1	2.20	0.50
1:g:181:ALA:HB1	1:g:236:ARG:HB3	1.93	0.50
1:p:55:SER:HB3	1:p:56:PRO:HD3	1.91	0.50
1:a:47:GLN:NE2	1:q:215:GLU:OE1	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:a:55:SER:HB3	1:a:56:PRO:HD3	1.94	0.50
1:e:211:ARG:NH2	1:f:33:ASN:OD1	2.27	0.50
1:k:220:SER:HB3	1:k:221:PRO:HD3	1.94	0.50
1:g:11:VAL:HG23	1:h:24:ILE:HG12	1.92	0.50
1:h:215:GLU:OE1	1:i:47:GLN:NE2	2.45	0.50
1:k:11:VAL:HG23	1:l:24:ILE:HG12	1.93	0.50
1:m:318:THR:HG22	1:m:320:GLY:H	1.77	0.50
1:n:181:ALA:HB1	1:n:236:ARG:HB3	1.94	0.50
1:a:11:VAL:HG23	1:b:24:ILE:HG12	1.94	0.50
1:g:318:THR:HG22	1:g:320:GLY:H	1.76	0.50
1:l:211:ARG:NH2	1:m:33:ASN:OD1	2.31	0.50
1:i:193:THR:OG1	1:j:237:TYR:OH	2.18	0.50
1:l:220:SER:HB3	1:l:221:PRO:HD3	1.94	0.50
1:q:181:ALA:HB1	1:q:236:ARG:HB3	1.94	0.50
1:h:81:THR:HG21	1:i:126:ASN:HD21	1.77	0.49
1:k:92:ARG:HH12	1:k:101:ASN:HB3	1.77	0.49
1:f:109:ASP:OD1	1:f:112:ARG:NH2	2.44	0.49
1:g:193:THR:HG1	1:h:237:TYR:HH	1.52	0.49
1:e:220:SER:HB3	1:e:221:PRO:HD3	1.94	0.49
1:g:215:GLU:OE1	1:h:47:GLN:NE2	2.45	0.49
1:b:92:ARG:HH12	1:b:101:ASN:HB3	1.77	0.49
1:h:72:TYR:OH	1:i:28:THR:OG1	2.24	0.49
1:g:220:SER:HB3	1:g:221:PRO:HD3	1.94	0.49
1:m:220:SER:HB3	1:m:221:PRO:HD3	1.95	0.49
1:o:55:SER:HB3	1:o:56:PRO:HD3	1.93	0.49
1:p:181:ALA:HB1	1:p:236:ARG:HB3	1.94	0.49
1:b:215:GLU:OE1	1:c:47:GLN:NE2	2.46	0.48
1:j:77:ASP:HB3	1:j:78:PRO:HD3	1.95	0.48
1:h:11:VAL:HG23	1:i:24:ILE:HG12	1.94	0.48
1:l:11:VAL:HG23	1:m:24:ILE:HG12	1.95	0.48
1:m:72:TYR:OH	1:n:28:THR:OG1	2.21	0.48
1:m:215:GLU:OE1	1:n:47:GLN:NE2	2.46	0.48
1:n:90:ARG:NH1	1:o:34:GLN:OE1	2.46	0.48
1:q:220:SER:HB3	1:q:221:PRO:HD3	1.96	0.48
1:n:318:THR:HG22	1:n:320:GLY:H	1.77	0.48
1:f:92:ARG:HH12	1:f:101:ASN:HB3	1.79	0.48
1:q:295:VAL:HG13	1:q:299:ARG:HH12	1.79	0.48
1:d:220:SER:HB3	1:d:221:PRO:HD3	1.96	0.48
1:e:55:SER:HB3	1:e:56:PRO:HD3	1.96	0.47
1:c:242:ASP:HB3	1:c:243:PRO:HD3	1.96	0.47
1:e:46:ARG:HE	1:e:50:GLU:HG3	1.78	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:d:181:ALA:HB1	1:d:236:ARG:HB3	1.97	0.47
1:p:242:ASP:HB3	1:p:243:PRO:HD3	1.97	0.47
1:h:242:ASP:HB3	1:h:243:PRO:HD3	1.96	0.47
1:l:242:ASP:HB3	1:l:243:PRO:HD3	1.96	0.47
1:b:229:ASP:OD1	1:b:230:SER:N	2.47	0.46
1:l:318:THR:HG22	1:l:320:GLY:H	1.81	0.46
1:p:215:GLU:OE1	1:q:47:GLN:NE2	2.48	0.46
1:p:220:SER:HB3	1:p:221:PRO:HD3	1.96	0.46
1:l:77:ASP:HB3	1:l:78:PRO:HD3	1.97	0.46
1:k:181:ALA:HB1	1:k:236:ARG:HB3	1.97	0.46
1:g:81:THR:HG21	1:h:126:ASN:HD21	1.80	0.46
1:k:211:ARG:NH2	1:l:33:ASN:OD1	2.24	0.46
1:m:77:ASP:HB3	1:m:78:PRO:HD3	1.96	0.46
1:d:193:THR:HG1	1:e:237:TYR:HH	1.56	0.46
1:k:57:GLN:O	1:k:59:THR:N	2.48	0.46
1:p:77:ASP:HB3	1:p:78:PRO:HD3	1.97	0.46
1:d:242:ASP:HB3	1:d:243:PRO:HD3	1.98	0.46
1:f:77:ASP:HB3	1:f:78:PRO:HD3	1.97	0.46
1:a:318:THR:HG22	1:a:320:GLY:H	1.81	0.46
1:h:318:THR:HG22	1:h:320:GLY:H	1.81	0.46
1:o:77:ASP:HB3	1:o:78:PRO:HD3	1.98	0.46
1:d:92:ARG:HH12	1:d:101:ASN:HB3	1.81	0.45
1:j:211:ARG:NH2	1:k:33:ASN:OD1	2.37	0.45
1:c:318:THR:HG22	1:c:320:GLY:H	1.81	0.45
1:j:242:ASP:HB3	1:j:243:PRO:HD3	1.98	0.45
1:n:242:ASP:HB3	1:n:243:PRO:HD3	1.98	0.45
1:g:72:TYR:OH	1:h:28:THR:OG1	2.22	0.45
1:m:92:ARG:NH1	1:m:101:ASN:HB3	2.30	0.45
1:m:167:TYR:CD1	1:m:315:LEU:HB3	2.52	0.45
1:f:242:ASP:HB3	1:f:243:PRO:HD3	1.99	0.45
1:n:201:GLN:HG2	1:o:253:ASP:OD1	2.17	0.45
1:a:77:ASP:HB3	1:a:78:PRO:HD3	1.99	0.45
1:g:242:ASP:HB3	1:g:243:PRO:HD3	1.98	0.45
1:m:242:ASP:HB3	1:m:243:PRO:HD3	1.98	0.45
1:e:318:THR:HG22	1:e:320:GLY:H	1.82	0.45
1:j:220:SER:HB3	1:j:221:PRO:HD3	1.99	0.45
1:a:242:ASP:HB3	1:a:243:PRO:HD3	1.99	0.44
1:o:242:ASP:HB3	1:o:243:PRO:HD3	1.99	0.44
1:a:43:VAL:HG21	1:q:211:ARG:HD3	2.00	0.44
1:c:220:SER:HB3	1:c:221:PRO:HD3	1.99	0.44
1:d:77:ASP:HB3	1:d:78:PRO:HD3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:h:31:LEU:HD23	1:h:31:LEU:HA	1.85	0.44
1:d:318:THR:HG22	1:d:320:GLY:H	1.82	0.44
1:k:109:ASP:OD1	1:k:112:ARG:NH2	2.51	0.44
1:i:242:ASP:HB3	1:i:243:PRO:HD3	1.99	0.44
1:q:77:ASP:HB3	1:q:78:PRO:HD3	1.99	0.44
1:g:189:ILE:HD11	1:g:297:LEU:HB3	1.99	0.44
1:i:11:VAL:HG23	1:j:24:ILE:HG12	1.99	0.44
1:k:77:ASP:HB3	1:k:78:PRO:HD3	2.00	0.44
1:n:215:GLU:OE1	1:o:47:GLN:NE2	2.51	0.44
1:o:215:GLU:OE1	1:p:47:GLN:NE2	2.51	0.44
1:j:196:LEU:HD23	1:j:196:LEU:HA	1.87	0.43
1:i:188:LEU:HD13	1:i:216:VAL:HG11	2.00	0.43
1:b:46:ARG:HA	1:b:46:ARG:HD3	1.83	0.43
1:k:242:ASP:HB3	1:k:243:PRO:HD3	2.01	0.43
1:q:242:ASP:HB3	1:q:243:PRO:HD3	1.99	0.43
1:a:188:LEU:HD13	1:a:216:VAL:HG11	2.01	0.43
1:i:77:ASP:HB3	1:i:78:PRO:HD3	2.01	0.43
1:l:31:LEU:HD23	1:l:31:LEU:HA	1.83	0.43
1:b:318:THR:HG22	1:b:320:GLY:H	1.84	0.43
1:o:189:ILE:HD11	1:o:297:LEU:HB3	2.01	0.43
1:p:182:TRP:CD1	1:p:221:PRO:HD2	2.54	0.43
1:e:81:THR:HG21	1:f:126:ASN:HD21	1.84	0.43
1:g:92:ARG:NH1	1:g:101:ASN:HB3	2.34	0.43
1:q:318:THR:HG22	1:q:320:GLY:H	1.83	0.43
1:n:92:ARG:HH12	1:n:101:ASN:HB3	1.83	0.43
1:p:46:ARG:NE	1:p:50:GLU:OE2	2.52	0.43
1:a:93:ILE:HG13	1:b:111:THR:HA	2.01	0.42
1:a:126:ASN:HD21	1:q:81:THR:HG21	1.84	0.42
1:a:24:ILE:HG12	1:q:11:VAL:HG23	2.01	0.42
1:a:181:ALA:HB1	1:a:236:ARG:HB3	2.00	0.42
1:c:193:THR:OG1	1:d:237:TYR:OH	2.26	0.42
1:g:92:ARG:HH12	1:g:101:ASN:HB3	1.83	0.42
1:l:215:GLU:OE1	1:m:47:GLN:NE2	2.52	0.42
1:i:46:ARG:HD3	1:i:46:ARG:HA	1.87	0.42
1:h:77:ASP:HB3	1:h:78:PRO:HD3	2.00	0.42
1:j:26:LEU:HD23	1:j:26:LEU:HA	1.89	0.42
1:m:189:ILE:HD11	1:m:297:LEU:HB3	2.01	0.42
1:e:242:ASP:HB3	1:e:243:PRO:HD3	2.01	0.42
1:c:72:TYR:OH	1:d:28:THR:OG1	2.26	0.42
1:d:222:GLN:HB2	1:d:225:VAL:HB	2.02	0.42
1:b:109:ASP:OD1	1:b:112:ARG:NH2	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:e:77:ASP:HB3	1:e:78:PRO:HD3	2.02	0.42
1:b:189:ILE:HD11	1:b:297:LEU:HB3	2.02	0.42
1:c:215:GLU:OE1	1:d:47:GLN:NE2	2.53	0.42
1:j:81:THR:HG21	1:k:126:ASN:ND2	2.34	0.41
1:a:81:THR:HG21	1:b:126:ASN:ND2	2.34	0.41
1:c:93:ILE:HG13	1:d:111:THR:HA	2.02	0.41
1:d:79:LEU:HD11	1:d:131:GLU:HG3	2.01	0.41
1:i:72:TYR:OH	1:j:28:THR:OG1	2.26	0.41
1:n:193:THR:OG1	1:o:237:TYR:OH	2.25	0.41
1:q:46:ARG:HD3	1:q:46:ARG:HA	1.82	0.41
1:q:189:ILE:HD11	1:q:297:LEU:HB3	2.03	0.41
1:a:26:LEU:HD23	1:a:26:LEU:HA	1.91	0.41
1:h:196:LEU:HD23	1:h:196:LEU:HA	1.92	0.41
1:o:318:THR:HG22	1:o:320:GLY:H	1.85	0.41
1:c:77:ASP:HB3	1:c:78:PRO:HD3	2.03	0.41
1:d:31:LEU:HD23	1:d:31:LEU:HA	1.89	0.41
1:m:11:VAL:HG23	1:n:24:ILE:HG12	2.03	0.41
1:k:92:ARG:NH1	1:k:101:ASN:HB3	2.36	0.41
1:o:31:LEU:HD23	1:o:31:LEU:HA	1.84	0.41
1:o:201:GLN:HG2	1:p:253:ASP:OD1	2.20	0.41
1:i:31:LEU:HD23	1:i:31:LEU:HA	1.88	0.41
1:a:201:GLN:HG2	1:b:253:ASP:OD1	2.21	0.41
1:b:31:LEU:HD23	1:b:31:LEU:HA	1.83	0.41
1:m:31:LEU:HD23	1:m:31:LEU:HA	1.84	0.41
1:c:181:ALA:HB1	1:c:236:ARG:HB3	2.01	0.41
1:e:215:GLU:OE1	1:f:47:GLN:NE2	2.54	0.40
1:b:77:ASP:HB3	1:b:78:PRO:HD3	2.02	0.40
1:b:92:ARG:NH1	1:b:101:ASN:HB3	2.35	0.40
1:n:77:ASP:HB3	1:n:78:PRO:HD3	2.03	0.40
1:i:192:CYS:HB3	1:i:294:ILE:HD11	2.03	0.40
1:k:229:ASP:OD1	1:k:230:SER:N	2.54	0.40
1:c:201:GLN:HG2	1:d:253:ASP:OD1	2.21	0.40
1:c:211:ARG:NH2	1:d:33:ASN:OD1	2.29	0.40
1:l:189:ILE:HD11	1:l:297:LEU:HB3	2.03	0.40
1:l:201:GLN:HG2	1:m:253:ASP:OD1	2.20	0.40
1:h:26:LEU:HD23	1:h:26:LEU:HA	1.95	0.40

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:p:61:ARG:NH1	1:p:64:ASP:OD1[4_565]	2.11	0.09

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	279/323 (86%)	270 (97%)	7 (2%)	2 (1%)	18	47
1	b	279/323 (86%)	268 (96%)	9 (3%)	2 (1%)	18	47
1	c	279/323 (86%)	269 (96%)	8 (3%)	2 (1%)	18	47
1	d	279/323 (86%)	269 (96%)	8 (3%)	2 (1%)	18	47
1	e	279/323 (86%)	268 (96%)	9 (3%)	2 (1%)	18	47
1	f	279/323 (86%)	266 (95%)	11 (4%)	2 (1%)	18	47
1	g	279/323 (86%)	269 (96%)	8 (3%)	2 (1%)	18	47
1	h	279/323 (86%)	270 (97%)	7 (2%)	2 (1%)	18	47
1	i	279/323 (86%)	269 (96%)	8 (3%)	2 (1%)	18	47
1	j	279/323 (86%)	271 (97%)	6 (2%)	2 (1%)	18	47
1	k	279/323 (86%)	267 (96%)	9 (3%)	3 (1%)	11	36
1	l	279/323 (86%)	269 (96%)	8 (3%)	2 (1%)	18	47
1	m	279/323 (86%)	269 (96%)	8 (3%)	2 (1%)	18	47
1	n	279/323 (86%)	271 (97%)	6 (2%)	2 (1%)	18	47
1	o	279/323 (86%)	270 (97%)	7 (2%)	2 (1%)	18	47
1	p	279/323 (86%)	269 (96%)	8 (3%)	2 (1%)	18	47
1	q	279/323 (86%)	269 (96%)	8 (3%)	2 (1%)	18	47
All	All	4743/5491 (86%)	4573 (96%)	135 (3%)	35 (1%)	18	47

All (35) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	k	58	VAL
1	a	55	SER
1	a	220	SER
1	b	220	SER
1	c	55	SER
1	c	220	SER
1	d	55	SER
1	d	220	SER
1	e	220	SER
1	f	55	SER
1	f	220	SER
1	g	220	SER
1	h	220	SER
1	i	55	SER
1	i	220	SER
1	j	220	SER
1	k	220	SER
1	l	55	SER
1	l	220	SER
1	m	55	SER
1	m	220	SER
1	n	55	SER
1	n	220	SER
1	o	55	SER
1	o	220	SER
1	p	55	SER
1	p	220	SER
1	q	55	SER
1	q	220	SER
1	b	55	SER
1	e	55	SER
1	g	55	SER
1	h	55	SER
1	j	55	SER
1	k	55	SER

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	252/277 (91%)	251 (100%)	1 (0%)	84	94
1	b	252/277 (91%)	252 (100%)	0	100	100
1	c	252/277 (91%)	252 (100%)	0	100	100
1	d	252/277 (91%)	252 (100%)	0	100	100
1	e	252/277 (91%)	252 (100%)	0	100	100
1	f	252/277 (91%)	252 (100%)	0	100	100
1	g	252/277 (91%)	252 (100%)	0	100	100
1	h	252/277 (91%)	252 (100%)	0	100	100
1	i	252/277 (91%)	251 (100%)	1 (0%)	84	94
1	j	252/277 (91%)	252 (100%)	0	100	100
1	k	252/277 (91%)	252 (100%)	0	100	100
1	l	252/277 (91%)	252 (100%)	0	100	100
1	m	252/277 (91%)	252 (100%)	0	100	100
1	n	252/277 (91%)	252 (100%)	0	100	100
1	o	252/277 (91%)	251 (100%)	1 (0%)	84	94
1	p	252/277 (91%)	251 (100%)	1 (0%)	84	94
1	q	252/277 (91%)	251 (100%)	1 (0%)	84	94
All	All	4284/4709 (91%)	4279 (100%)	5 (0%)	88	97

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

Mol	Chain	Res	Type
1	a	136	THR
1	i	39	GLN
1	o	94	ILE
1	p	136	THR
1	q	136	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	a	29	ASN
1	a	57	GLN
1	b	29	ASN
1	b	198	ASN
1	b	201	GLN

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Mol	Chain	Res	Type
1	b	204	GLN
1	c	47	GLN
1	d	29	ASN
1	d	57	GLN
1	d	198	ASN
1	e	29	ASN
1	f	47	GLN
1	g	29	ASN
1	g	126	ASN
1	h	29	ASN
1	h	57	GLN
1	h	201	GLN
1	h	204	GLN
1	i	29	ASN
1	i	204	GLN
1	k	57	GLN
1	k	222	GLN
1	l	29	ASN
1	m	29	ASN
1	m	47	GLN
1	m	57	GLN
1	m	198	ASN
1	n	29	ASN
1	n	47	GLN
1	n	126	ASN
1	o	29	ASN
1	p	29	ASN
1	p	47	GLN
1	p	201	GLN
1	q	29	ASN
1	q	198	ASN

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

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	287/323 (88%)	-0.16	8 (2%) 55 45	34, 49, 105, 214	0
1	b	287/323 (88%)	-0.19	7 (2%) 59 49	36, 49, 100, 219	0
1	c	287/323 (88%)	-0.13	10 (3%) 47 38	37, 52, 106, 179	0
1	d	287/323 (88%)	-0.07	10 (3%) 47 38	39, 57, 112, 190	0
1	e	287/323 (88%)	0.11	12 (4%) 40 32	42, 62, 116, 189	0
1	f	287/323 (88%)	0.13	8 (2%) 55 45	42, 63, 112, 210	0
1	g	287/323 (88%)	0.07	9 (3%) 51 41	44, 62, 112, 206	0
1	h	287/323 (88%)	0.08	8 (2%) 55 45	44, 61, 111, 184	0
1	i	287/323 (88%)	-0.05	8 (2%) 55 45	43, 60, 114, 204	0
1	j	287/323 (88%)	-0.08	4 (1%) 73 64	42, 59, 106, 205	0
1	k	287/323 (88%)	-0.06	10 (3%) 47 38	41, 57, 107, 198	0
1	l	287/323 (88%)	-0.01	10 (3%) 47 38	43, 56, 116, 223	0
1	m	287/323 (88%)	-0.06	9 (3%) 51 41	40, 57, 108, 209	0
1	n	287/323 (88%)	-0.06	8 (2%) 55 45	42, 55, 111, 189	0
1	o	287/323 (88%)	-0.10	7 (2%) 59 49	38, 53, 107, 200	0
1	p	287/323 (88%)	-0.13	11 (3%) 44 35	38, 49, 103, 183	0
1	q	287/323 (88%)	-0.17	10 (3%) 47 38	35, 48, 100, 172	0
All	All	4879/5491 (88%)	-0.05	149 (3%) 51 41	34, 57, 111, 223	0

All (149) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	q	275	ALA	4.9
1	i	275	ALA	4.8
1	c	155	GLY	4.8
1	p	102	PRO	4.8

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Mol	Chain	Res	Type	RSRZ
1	q	320	GLY	4.7
1	k	320	GLY	4.7
1	l	320	GLY	4.7
1	o	102	PRO	4.3
1	g	104	THR	4.3
1	b	275	ALA	4.3
1	m	275	ALA	4.3
1	d	102	PRO	4.2
1	e	104	THR	4.2
1	e	275	ALA	4.1
1	c	104	THR	4.0
1	h	104	THR	4.0
1	a	320	GLY	4.0
1	l	155	GLY	4.0
1	l	275	ALA	4.0
1	f	102	PRO	3.9
1	f	275	ALA	3.9
1	j	275	ALA	3.9
1	k	275	ALA	3.9
1	p	275	ALA	3.9
1	p	104	THR	3.9
1	b	320	GLY	3.8
1	a	275	ALA	3.7
1	h	155	GLY	3.7
1	c	275	ALA	3.7
1	p	155	GLY	3.7
1	p	320	GLY	3.7
1	h	254	THR	3.6
1	i	320	GLY	3.6
1	d	104	THR	3.6
1	g	102	PRO	3.6
1	h	275	ALA	3.5
1	f	104	THR	3.5
1	g	275	ALA	3.5
1	e	102	PRO	3.5
1	b	155	GLY	3.4
1	d	155	GLY	3.4
1	g	279	VAL	3.4
1	m	102	PRO	3.4
1	a	102	PRO	3.4
1	n	102	PRO	3.4
1	b	102	PRO	3.3

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Mol	Chain	Res	Type	RSRZ
1	l	102	PRO	3.3
1	o	320	GLY	3.3
1	n	275	ALA	3.3
1	f	279	VAL	3.2
1	q	102	PRO	3.2
1	e	320	GLY	3.2
1	m	320	GLY	3.2
1	o	155	GLY	3.2
1	p	65	SER	3.1
1	m	104	THR	3.1
1	a	276	THR	3.1
1	o	275	ALA	3.1
1	c	3	SER	3.1
1	m	255	ARG	3.1
1	o	104	THR	3.0
1	c	320	GLY	3.0
1	f	201	GLN	3.0
1	c	102	PRO	3.0
1	n	155	GLY	2.9
1	l	280	ASP	2.9
1	p	154	SER	2.9
1	g	320	GLY	2.9
1	i	102	PRO	2.9
1	m	155	GLY	2.8
1	n	104	THR	2.8
1	n	254	THR	2.8
1	q	104	THR	2.8
1	f	320	GLY	2.8
1	f	171	THR	2.8
1	h	102	PRO	2.8
1	h	279	VAL	2.7
1	e	65	SER	2.7
1	b	104	THR	2.7
1	h	320	GLY	2.7
1	l	104	THR	2.6
1	a	279	VAL	2.6
1	k	3	SER	2.6
1	i	155	GLY	2.6
1	g	201	GLN	2.6
1	p	168	SER	2.6
1	d	278	ARG	2.6
1	k	102	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
1	b	279	VAL	2.6
1	e	171	THR	2.6
1	e	255	ARG	2.5
1	e	155	GLY	2.5
1	d	275	ALA	2.5
1	k	276	THR	2.5
1	j	320	GLY	2.5
1	b	276	THR	2.5
1	l	254	THR	2.5
1	e	278	ARG	2.4
1	d	201	GLN	2.4
1	a	280	ASP	2.4
1	k	94	ILE	2.4
1	q	65	SER	2.4
1	n	278	ARG	2.4
1	c	201	GLN	2.4
1	m	254	THR	2.4
1	o	276	THR	2.4
1	l	255	ARG	2.4
1	j	104	THR	2.4
1	n	320	GLY	2.3
1	o	65	SER	2.3
1	q	279	VAL	2.3
1	i	104	THR	2.3
1	f	278	ARG	2.3
1	k	155	GLY	2.3
1	m	276	THR	2.3
1	d	279	VAL	2.3
1	g	155	GLY	2.3
1	e	279	VAL	2.3
1	d	320	GLY	2.2
1	q	229	ASP	2.2
1	d	276	THR	2.2
1	g	276	THR	2.2
1	j	102	PRO	2.2
1	l	279	VAL	2.2
1	a	255	ARG	2.2
1	e	302	GLY	2.2
1	c	126	ASN	2.2
1	h	276	THR	2.2
1	p	94	ILE	2.1
1	c	279	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	p	254	THR	2.1
1	k	58	VAL	2.1
1	a	229	ASP	2.1
1	e	276	THR	2.1
1	i	254	THR	2.1
1	d	255	ARG	2.1
1	c	94	ILE	2.1
1	k	57	GLN	2.1
1	q	201	GLN	2.1
1	l	201	GLN	2.1
1	i	276	THR	2.1
1	k	104	THR	2.1
1	p	201	GLN	2.0
1	g	94	ILE	2.0
1	q	254	THR	2.0
1	m	95	GLU	2.0
1	n	255	ARG	2.0
1	q	155	GLY	2.0
1	i	279	VAL	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.