



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:08 AM UTC

PDB ID : 1M06 / pdb_00001m06
Title : Structural Studies of Bacteriophage alpha3 Assembly, X-Ray Crystallography
Authors : Bernal, R.A.; Hafenstein, S.; Olson, N.H.; Bowman, V.; Chipman, P.R.; Baker, T.S.; Fane, B.A.; Rossmann, M.G.
Deposited on : 2002-06-12
Resolution : 3.50 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
Mogul : 2022.3.0, CSD as543be (2022)
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

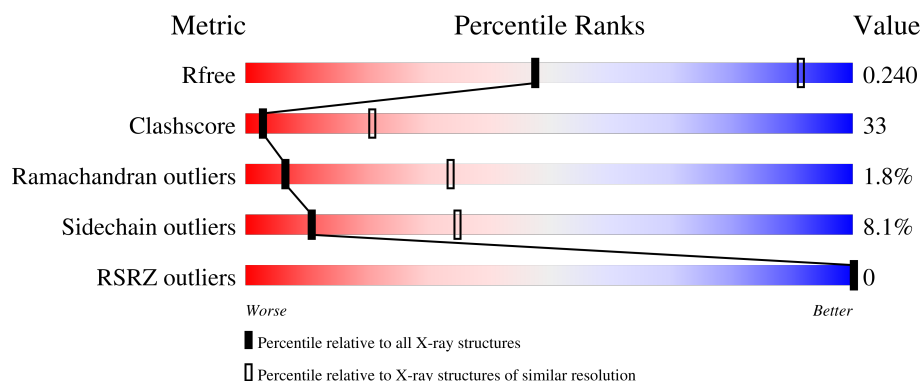
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

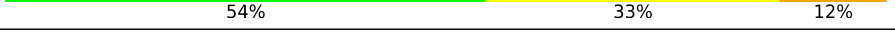
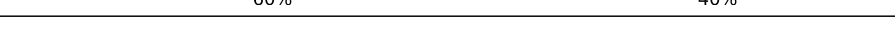
The reported resolution of this entry is 3.50 Å.

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	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	F	431	
2	G	187	
3	J	24	
4	X	10	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	3DR	X	1	-	-	X	-
4	3DR	X	10	-	-	X	-
4	3DR	X	2	X	-	X	-
4	3DR	X	3	-	-	X	-
4	3DR	X	7	-	-	X	-
4	3DR	X	8	-	-	X	-
4	3DR	X	9	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5093 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	F	422	Total	C	N	O	S	0	0	0
			3411	2168	599	628	16			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	160	GLY	ARG	SEE REMARK 999	UNP P08767

- Molecule 2 is a protein called Major spike protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	187	Total	C	N	O	S	0	0	0
			1373	855	235	275	8			

- Molecule 3 is a protein called Small core protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	J	24	Total	C	N	O	S	0	0	0
			199	124	44	30	1			

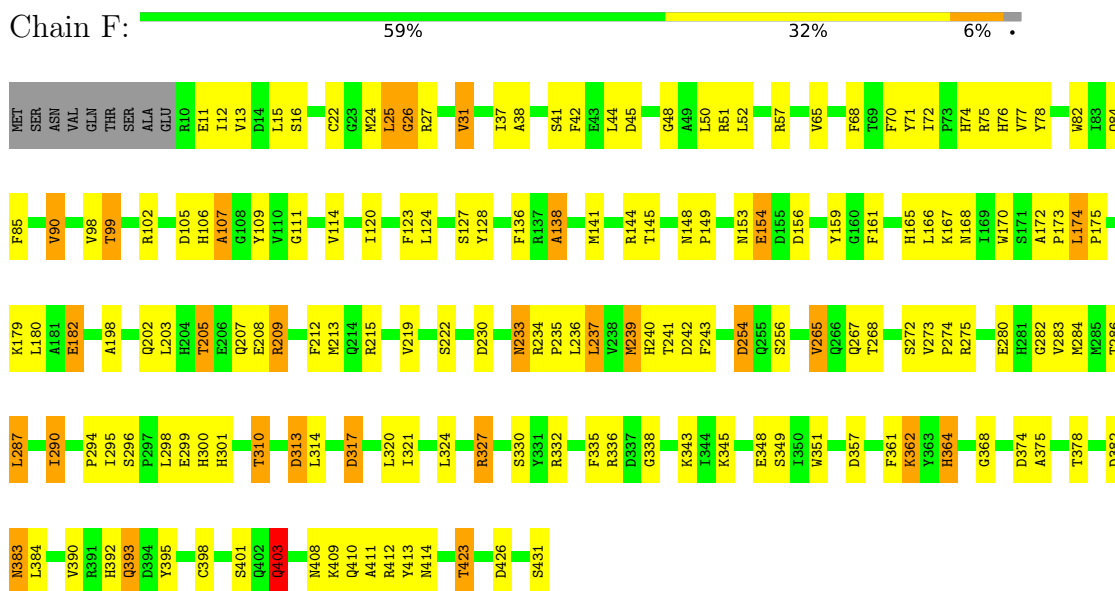
- Molecule 4 is a DNA chain called 5'-D(P*(3DR)P*(3DR)P*(3DR)P*(3DR)P*(3DR)P*(3DR)P*(3DR)P*(3DR)P*(3DR)P*(3DR))-3'.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	X	10	Total	C	O	P	0	0	0
			110	50	50	10			

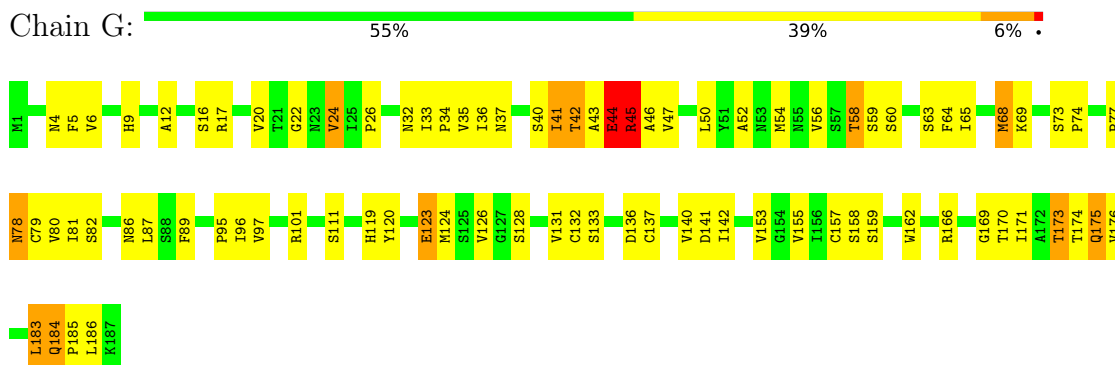
3 Residue-property plots

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: Capsid Protein



• Molecule 2: Major spike protein



• Molecule 3: Small core protein





- Molecule 4: 5'-D(P*(3DR)P*(3DR)P*(3DR)P*(3DR)P*(3DR)P*(3DR)P*(3DR)P*(3DR)P*(3DR)P*(3DR)P*(3DR))-3'

Chain X:  60% 40%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	290.26Å 332.12Å 337.70Å 90.00° 94.08° 90.00°	Depositor
Resolution (Å)	82.45 – 3.50 82.45 – 3.50	Depositor EDS
% Data completeness (in resolution range)	74.6 (82.45-3.50) 60.4 (82.45-3.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.67 (at 3.49Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.232 , 0.234 0.238 , 0.240	Depositor DCC
R_{free} test set	59662 reflections (9.99%)	wwPDB-VP
Wilson B-factor (Å ²)	30.2	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 597.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.35$, $\langle L^2 \rangle = 0.18$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.11	EDS
Total number of atoms	5093	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.32% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 3DR

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	F	0.57	0/3512	1.00	13/4775 (0.3%)
2	G	0.55	0/1399	1.13	7/1910 (0.4%)
3	J	0.73	0/203	1.05	2/265 (0.8%)
All	All	0.57	0/5114	1.04	22/6950 (0.3%)

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
4	X	1	0

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	119	HIS	N-CA-C	6.55	120.71	110.17
1	F	12	ILE	N-CA-C	6.43	117.95	109.21
2	G	120	TYR	CA-C-N	6.20	126.16	119.78
2	G	120	TYR	C-N-CA	6.20	126.16	119.78
1	F	314	LEU	N-CA-C	6.04	119.03	111.24
1	F	136	PHE	N-CA-C	6.04	120.81	112.90
1	F	138	ALA	N-CA-C	-6.04	102.16	109.83
2	G	184	GLN	CA-C-N	5.95	127.28	119.84
2	G	184	GLN	C-N-CA	5.95	127.28	119.84
3	J	2	LYS	N-CA-C	-5.85	104.90	112.68
1	F	111	GLY	N-CA-C	5.81	121.96	113.48
1	F	145	THR	N-CA-C	5.74	119.48	111.56
2	G	45	ARG	N-CA-C	-5.71	98.98	108.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	317	ASP	CA-C-N	5.41	124.87	119.24
1	F	317	ASP	C-N-CA	5.41	124.87	119.24
2	G	44	GLU	N-CA-C	-5.32	104.30	111.54
1	F	107	ALA	N-CA-C	-5.31	105.49	113.89
1	F	403	GLN	N-CA-C	-5.31	105.66	111.82
1	F	272	SER	N-CA-C	5.17	116.84	108.41
3	J	17	TRP	N-CA-C	5.11	116.94	108.20
1	F	398	CYS	N-CA-C	5.10	119.14	113.02
1	F	99	THR	N-CA-C	5.08	118.35	110.17

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	X	2	3DR	C4'

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	F	3411	0	3262	149	0
2	G	1373	0	1354	114	0
3	J	199	0	212	23	0
4	X	110	0	81	79	0
All	All	5093	0	4909	326	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 33.

All (326) 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)
2:G:46:ALA:CB	2:G:68:MET:CE	1.78	1.59
2:G:46:ALA:CB	2:G:68:MET:HE3	1.12	1.55
2:G:46:ALA:HB2	2:G:68:MET:CE	1.40	1.31

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:X:9:3DR:C1'	4:X:10:3DR:OP1	1.84	1.25
2:G:96:ILE:HA	2:G:124:MET:HE1	1.27	1.17
2:G:42:THR:C	2:G:44:GLU:H	1.49	1.15
2:G:46:ALA:N	2:G:68:MET:HE1	1.62	1.13
4:X:8:3DR:OP1	4:X:9:3DR:H1'2	1.49	1.13
1:F:167:LYS:HE3	4:X:2:3DR:H3'	1.31	1.10
2:G:46:ALA:CB	2:G:68:MET:HE1	1.80	1.10
2:G:46:ALA:HB3	2:G:68:MET:CE	1.59	1.10
3:J:23:GLN:O	4:X:2:3DR:H4'1	1.54	1.06
4:X:3:3DR:H3'	4:X:4:3DR:H5''	1.34	1.06
2:G:78:ASN:HB3	2:G:141:ASP:HA	1.29	1.06
1:F:393:GLN:H	1:F:393:GLN:NE2	1.54	1.05
4:X:9:3DR:C2'	4:X:10:3DR:OP1	2.04	1.05
4:X:9:3DR:H1'2	4:X:10:3DR:OP1	1.57	1.01
2:G:46:ALA:HB3	2:G:68:MET:HE3	1.09	0.97
1:F:393:GLN:HE21	1:F:393:GLN:N	1.63	0.96
2:G:95:PRO:HG2	2:G:132:CYS:HB3	1.47	0.96
2:G:41:ILE:H	2:G:41:ILE:CD1	1.80	0.94
1:F:51:ARG:HH22	1:F:410:GLN:HE21	1.07	0.94
2:G:4:ASN:HD21	2:G:6:VAL:HG12	1.33	0.93
1:F:167:LYS:CE	4:X:2:3DR:H3'	1.98	0.93
1:F:361:PHE:O	1:F:364:HIS:HB3	1.70	0.92
4:X:9:3DR:H2'	4:X:10:3DR:P	2.09	0.92
3:J:24:PHE:OXT	4:X:2:3DR:H2'	1.71	0.91
4:X:8:3DR:OP1	4:X:10:3DR:C5'	2.19	0.90
4:X:6:3DR:O3'	4:X:7:3DR:P	2.30	0.90
4:X:1:3DR:H2''	4:X:2:3DR:OP1	1.72	0.90
2:G:12:ALA:HB2	2:G:45:ARG:HG3	1.55	0.89
2:G:41:ILE:N	2:G:41:ILE:HD13	1.86	0.88
2:G:42:THR:C	2:G:44:GLU:N	2.24	0.88
4:X:9:3DR:H1'2	4:X:10:3DR:O5'	1.74	0.88
2:G:46:ALA:CA	2:G:68:MET:HE1	2.03	0.88
1:F:52:LEU:HB2	1:F:265:VAL:CG1	2.03	0.87
4:X:9:3DR:H1'2	4:X:10:3DR:P	2.14	0.87
4:X:8:3DR:OP1	4:X:9:3DR:C1'	2.21	0.87
4:X:8:3DR:OP1	4:X:10:3DR:O5'	1.91	0.87
1:F:52:LEU:HB2	1:F:265:VAL:HG11	1.55	0.85
1:F:241:THR:HG23	4:X:10:3DR:O3'	1.76	0.85
2:G:96:ILE:CA	2:G:124:MET:HE1	2.06	0.85
2:G:44:GLU:CG	2:G:175:GLN:HB3	2.07	0.85
2:G:46:ALA:HB2	2:G:68:MET:HE3	0.86	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:97:VAL:HG22	2:G:124:MET:HE3	1.59	0.84
1:F:243:PHE:HA	4:X:9:3DR:OP2	1.78	0.83
2:G:42:THR:O	2:G:44:GLU:N	2.11	0.83
1:F:182:GLU:HG3	1:F:203:LEU:HD13	1.59	0.82
2:G:44:GLU:HG3	2:G:175:GLN:HB3	1.62	0.82
4:X:3:3DR:H3'	4:X:4:3DR:C5'	2.11	0.81
1:F:393:GLN:H	1:F:393:GLN:HE21	0.82	0.80
1:F:403:GLN:HE21	1:F:403:GLN:HA	1.46	0.80
1:F:167:LYS:HZ2	4:X:2:3DR:C3'	1.95	0.79
2:G:45:ARG:C	2:G:68:MET:HE1	2.07	0.79
1:F:219:VAL:O	1:F:222:SER:HB3	1.83	0.79
2:G:126:VAL:HG22	2:G:133:SER:HB3	1.63	0.78
2:G:41:ILE:H	2:G:41:ILE:HD13	1.42	0.78
3:J:24:PHE:C	4:X:2:3DR:H2'	2.08	0.77
2:G:80:VAL:HG13	2:G:176:VAL:HB	1.65	0.77
2:G:68:MET:SD	2:G:173:THR:HG23	2.25	0.77
3:J:23:GLN:O	4:X:2:3DR:C4'	2.33	0.76
2:G:41:ILE:CD1	2:G:41:ILE:N	2.45	0.76
2:G:81:ILE:HG21	2:G:153:VAL:HG11	1.67	0.76
4:X:9:3DR:C2'	4:X:10:3DR:P	2.73	0.76
1:F:383:ASN:C	1:F:383:ASN:HD22	1.91	0.75
3:J:15:ARG:NH1	4:X:7:3DR:C3'	2.50	0.75
2:G:44:GLU:OE2	2:G:44:GLU:HA	1.87	0.75
4:X:8:3DR:P	4:X:9:3DR:O4'	2.45	0.75
1:F:287:LEU:N	1:F:287:LEU:HD12	2.02	0.74
2:G:17:ARG:HG3	2:G:17:ARG:HH11	1.52	0.73
3:J:16:LEU:HD13	4:X:4:3DR:OP1	1.88	0.73
1:F:310:THR:HG23	1:F:313:ASP:CG	2.13	0.73
3:J:15:ARG:HH12	4:X:7:3DR:C3'	2.01	0.73
1:F:310:THR:HG23	1:F:313:ASP:OD2	1.88	0.72
1:F:403:GLN:HA	1:F:403:GLN:NE2	2.02	0.72
2:G:32:ASN:O	2:G:34:PRO:HD3	1.89	0.72
4:X:1:3DR:P	4:X:7:3DR:H2''	2.30	0.72
3:J:1:MET:O	3:J:1:MET:HG2	1.91	0.71
3:J:24:PHE:HD1	4:X:2:3DR:H1'2	1.54	0.71
1:F:51:ARG:NH2	1:F:410:GLN:HE21	1.86	0.71
4:X:9:3DR:C1'	4:X:10:3DR:P	2.77	0.71
1:F:239:MET:HA	1:F:239:MET:HE2	1.73	0.70
1:F:31:VAL:HG11	1:F:409:LYS:HE3	1.73	0.69
2:G:17:ARG:HG2	2:G:47:VAL:HB	1.75	0.69
2:G:45:ARG:HD2	2:G:174:THR:HB	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:12:ALA:CB	2:G:45:ARG:HG3	2.22	0.68
2:G:44:GLU:OE1	2:G:68:MET:HG3	1.93	0.68
4:X:9:3DR:H2'	4:X:10:3DR:OP2	1.93	0.68
1:F:242:ASP:N	4:X:10:3DR:H2''	2.09	0.68
1:F:31:VAL:HG11	1:F:409:LYS:NZ	2.09	0.68
4:X:8:3DR:P	4:X:9:3DR:C1'	2.82	0.67
1:F:107:ALA:HB2	1:F:114:VAL:HG22	1.76	0.67
3:J:15:ARG:NH1	4:X:7:3DR:O3'	2.28	0.67
3:J:15:ARG:HH12	4:X:7:3DR:H3'	1.60	0.67
4:X:9:3DR:H1'1	4:X:10:3DR:OP1	1.92	0.67
2:G:78:ASN:HB3	2:G:141:ASP:CA	2.17	0.66
2:G:33:ILE:HG21	2:G:65:ILE:HG13	1.75	0.66
1:F:243:PHE:CA	4:X:9:3DR:OP2	2.44	0.66
3:J:8:PRO:HG2	3:J:10:ARG:NH1	2.10	0.66
1:F:31:VAL:HG11	1:F:409:LYS:CE	2.25	0.66
2:G:81:ILE:HD13	2:G:153:VAL:HG11	1.78	0.65
1:F:167:LYS:NZ	4:X:2:3DR:C3'	2.58	0.65
1:F:383:ASN:HD22	1:F:384:LEU:N	1.93	0.65
1:F:50:LEU:HD23	1:F:50:LEU:C	2.21	0.65
1:F:242:ASP:N	4:X:8:3DR:OP2	2.23	0.65
1:F:99:THR:H	1:F:148:ASN:HD22	1.43	0.64
2:G:44:GLU:O	2:G:174:THR:HA	1.97	0.64
1:F:383:ASN:C	1:F:383:ASN:ND2	2.56	0.64
3:J:5:ARG:HH11	3:J:5:ARG:HG2	1.63	0.64
2:G:46:ALA:HB2	2:G:68:MET:HE2	1.68	0.63
2:G:87:LEU:O	2:G:131:VAL:HG23	1.97	0.63
4:X:1:3DR:OP2	4:X:7:3DR:C3'	2.46	0.63
2:G:166:ARG:HG3	2:G:166:ARG:HH11	1.63	0.63
2:G:44:GLU:OE2	2:G:68:MET:HB2	1.99	0.62
3:J:16:LEU:HB2	4:X:4:3DR:OP1	2.00	0.62
1:F:22:CYS:HA	1:F:408:ASN:HA	1.82	0.62
1:F:78:TYR:HB2	1:F:82:TRP:HB2	1.80	0.62
2:G:64:PHE:CE1	2:G:157:CYS:HB3	2.35	0.62
1:F:167:LYS:HZ2	4:X:2:3DR:C4'	2.12	0.62
2:G:41:ILE:HG12	2:G:46:ALA:HB2	1.81	0.62
4:X:3:3DR:C3'	4:X:4:3DR:H5''	2.21	0.61
2:G:16:SER:HB3	2:G:45:ARG:O	2.00	0.61
1:F:90:VAL:HG22	1:F:90:VAL:O	1.99	0.61
1:F:167:LYS:NZ	4:X:2:3DR:H3'	2.16	0.61
3:J:15:ARG:NH1	4:X:7:3DR:H3'	2.16	0.61
1:F:403:GLN:HE21	1:F:403:GLN:CA	2.11	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:26:GLY:HA3	1:F:390:VAL:HB	1.83	0.60
3:J:17:TRP:HB2	4:X:3:3DR:H1'	1.82	0.60
1:F:208:GLU:HG2	1:F:212:PHE:CZ	2.37	0.60
1:F:242:ASP:H	4:X:10:3DR:H2''	1.67	0.59
1:F:75:ARG:NH2	1:F:230:ASP:OD1	2.36	0.59
2:G:123:GLU:HB2	2:G:136:ASP:OD2	2.03	0.59
1:F:167:LYS:NZ	4:X:2:3DR:C4'	2.66	0.59
1:F:209:ARG:HA	1:F:213:MET:HB2	1.85	0.59
1:F:332:ARG:HD2	1:F:338:GLY:O	2.02	0.58
2:G:183:LEU:HD23	2:G:183:LEU:C	2.28	0.58
1:F:412:ARG:HG2	1:F:414:ASN:HD21	1.68	0.58
2:G:33:ILE:N	2:G:33:ILE:HD12	2.18	0.58
1:F:71:TYR:C	1:F:72:ILE:HD12	2.29	0.58
1:F:412:ARG:HG2	1:F:414:ASN:ND2	2.19	0.57
4:X:1:3DR:P	4:X:7:3DR:O3'	2.62	0.57
1:F:74:HIS:CD2	1:F:282:GLY:HA2	2.40	0.57
1:F:138:ALA:HB2	1:F:141:MET:HE3	1.86	0.57
1:F:24:MET:HE3	1:F:392:HIS:ND1	2.20	0.56
2:G:46:ALA:HB3	2:G:68:MET:SD	2.43	0.56
2:G:24:VAL:HG13	2:G:54:MET:HA	1.86	0.56
3:J:15:ARG:HG3	4:X:3:3DR:H2''	1.87	0.56
2:G:56:VAL:O	2:G:56:VAL:HG12	2.04	0.56
1:F:51:ARG:HH22	1:F:410:GLN:NE2	1.90	0.56
2:G:41:ILE:H	2:G:41:ILE:HD12	1.65	0.56
1:F:27:ARG:HD3	1:F:159:TYR:HB2	1.87	0.56
2:G:63:SER:HB3	2:G:158:SER:OG	2.05	0.55
1:F:327:ARG:HD3	1:F:348:GLU:OE2	2.06	0.55
4:X:8:3DR:H5''	4:X:9:3DR:H5''	1.88	0.55
1:F:362:LYS:O	1:F:362:LYS:HD3	2.07	0.55
1:F:393:GLN:NE2	1:F:393:GLN:N	2.33	0.55
2:G:95:PRO:CG	2:G:132:CYS:HB3	2.28	0.54
1:F:27:ARG:NH1	1:F:392:HIS:HD2	2.05	0.54
2:G:17:ARG:HG3	2:G:17:ARG:NH1	2.22	0.54
2:G:124:MET:HE2	2:G:132:CYS:HB2	1.89	0.54
1:F:213:MET:SD	1:F:222:SER:HB2	2.47	0.54
2:G:77:PRO:HG2	2:G:78:ASN:H	1.74	0.53
2:G:33:ILE:HD12	2:G:33:ILE:H	1.74	0.53
1:F:254:ASP:OD1	1:F:254:ASP:C	2.52	0.53
1:F:208:GLU:HB3	1:F:213:MET:CE	2.39	0.53
2:G:81:ILE:HD12	2:G:140:VAL:CG2	2.39	0.52
1:F:332:ARG:HG2	1:F:332:ARG:HH11	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:22:GLY:O	2:G:52:ALA:HB1	2.09	0.52
2:G:81:ILE:HD12	2:G:140:VAL:HG21	1.92	0.52
4:X:1:3DR:OP2	4:X:7:3DR:O3'	2.27	0.52
2:G:35:VAL:HG23	2:G:111:SER:HA	1.91	0.52
1:F:57:ARG:HG3	1:F:368:GLY:HA3	1.92	0.52
1:F:242:ASP:H	4:X:10:3DR:C2'	2.22	0.52
1:F:65:VAL:HG22	1:F:290:ILE:HG23	1.92	0.52
4:X:2:3DR:H1'1	4:X:3:3DR:OP1	2.10	0.52
1:F:27:ARG:HH11	1:F:392:HIS:HD2	1.58	0.51
1:F:243:PHE:HB3	4:X:8:3DR:O4'	2.10	0.51
2:G:79:CYS:SG	2:G:142:ILE:HG21	2.50	0.51
1:F:105:ASP:OD1	1:F:105:ASP:N	2.43	0.51
2:G:35:VAL:O	2:G:37:ASN:N	2.35	0.51
1:F:172:ALA:N	1:F:173:PRO:CD	2.73	0.51
1:F:243:PHE:HB2	4:X:9:3DR:P	2.51	0.51
1:F:138:ALA:CB	1:F:141:MET:HE3	2.41	0.51
2:G:16:SER:CB	2:G:45:ARG:O	2.59	0.51
1:F:72:ILE:HD12	1:F:72:ILE:N	2.26	0.51
2:G:4:ASN:ND2	2:G:6:VAL:H	2.07	0.51
2:G:41:ILE:CG1	2:G:46:ALA:HB2	2.41	0.51
1:F:179:LYS:HD3	1:F:182:GLU:HG2	1.92	0.51
4:X:6:3DR:HO3'	4:X:7:3DR:P	2.33	0.50
2:G:166:ARG:HG3	2:G:166:ARG:NH1	2.27	0.50
1:F:174:LEU:HD22	1:F:345:LYS:O	2.11	0.50
2:G:56:VAL:HG13	2:G:59:SER:O	2.11	0.50
4:X:2:3DR:OP1	4:X:3:3DR:C5'	2.59	0.50
2:G:4:ASN:ND2	2:G:4:ASN:C	2.69	0.49
4:X:2:3DR:OP1	4:X:3:3DR:H2'	2.11	0.49
1:F:198:ALA:O	1:F:202:GLN:HG3	2.12	0.49
2:G:42:THR:HB	2:G:69:LYS:HB3	1.94	0.49
1:F:138:ALA:HB3	1:F:141:MET:HG3	1.94	0.49
2:G:4:ASN:ND2	2:G:6:VAL:HG12	2.14	0.49
2:G:44:GLU:CD	2:G:68:MET:HG3	2.38	0.49
2:G:58:THR:HG22	2:G:59:SER:N	2.28	0.49
1:F:70:PHE:CD1	1:F:237:LEU:HA	2.47	0.49
1:F:109:TYR:OH	1:F:149:PRO:HG3	2.13	0.49
1:F:202:GLN:O	1:F:205:THR:HG23	2.12	0.49
2:G:4:ASN:HD21	2:G:6:VAL:CG1	2.14	0.49
1:F:42:PHE:CD1	1:F:42:PHE:C	2.91	0.48
2:G:159:SER:HB3	2:G:162:TRP:CE2	2.48	0.48
3:J:24:PHE:HA	4:X:2:3DR:C1'	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:349:SER:HA	1:F:351:TRP:CE2	2.48	0.48
1:F:31:VAL:HG11	1:F:409:LYS:HZ1	1.77	0.48
2:G:101:ARG:HD2	2:G:140:VAL:CG1	2.44	0.48
1:F:300:HIS:HD2	1:F:301:HIS:O	1.97	0.48
2:G:4:ASN:C	2:G:4:ASN:HD22	2.20	0.48
4:X:1:3DR:C2'	4:X:2:3DR:O5'	2.54	0.48
1:F:239:MET:HB2	1:F:273:VAL:HG22	1.95	0.48
1:F:208:GLU:HG2	1:F:212:PHE:HZ	1.78	0.47
1:F:144:ARG:NH1	1:F:161:PHE:CD2	2.83	0.47
1:F:166:LEU:HD22	1:F:166:LEU:N	2.28	0.47
2:G:80:VAL:CG1	2:G:176:VAL:HB	2.42	0.47
1:F:167:LYS:NZ	4:X:2:3DR:H4'1	2.29	0.47
1:F:167:LYS:HZ1	4:X:2:3DR:H4'1	1.80	0.47
1:F:174:LEU:O	1:F:175:PRO:C	2.57	0.47
2:G:63:SER:HB3	2:G:158:SER:CB	2.44	0.47
4:X:1:3DR:H2''	4:X:2:3DR:O5'	2.10	0.47
1:F:37:ILE:HG22	1:F:38:ALA:N	2.29	0.47
1:F:70:PHE:CE1	1:F:237:LEU:HG	2.50	0.47
1:F:295:ILE:HD12	1:F:357:ASP:CB	2.45	0.47
1:F:336:ARG:HG3	1:F:336:ARG:HH11	1.80	0.46
2:G:86:ASN:O	2:G:169:GLY:HA3	2.15	0.46
1:F:283:VAL:HG12	1:F:284:MET:N	2.30	0.46
1:F:16:SER:HA	1:F:413:TYR:O	2.15	0.46
2:G:26:PRO:HD3	2:G:54:MET:HE3	1.97	0.46
2:G:81:ILE:HG21	2:G:153:VAL:CG1	2.41	0.46
2:G:73:SER:HB2	2:G:74:PRO:HD3	1.97	0.46
2:G:89:PHE:CE2	2:G:162:TRP:HZ3	2.34	0.46
1:F:239:MET:HE2	1:F:240:HIS:H	1.81	0.46
2:G:44:GLU:CB	2:G:175:GLN:HB3	2.45	0.46
4:X:1:3DR:OP2	4:X:7:3DR:C2'	2.63	0.46
2:G:44:GLU:OE1	2:G:68:MET:SD	2.74	0.46
1:F:75:ARG:HH12	1:F:280:GLU:CD	2.23	0.46
3:J:11:ARG:HG3	3:J:11:ARG:HH11	1.80	0.46
1:F:98:VAL:HG12	1:F:120:ILE:O	2.15	0.45
1:F:236:LEU:HD12	1:F:237:LEU:N	2.30	0.45
2:G:183:LEU:C	2:G:183:LEU:CD2	2.89	0.45
1:F:68:PHE:O	1:F:286:THR:HA	2.16	0.45
2:G:44:GLU:OE2	2:G:68:MET:CG	2.65	0.45
1:F:13:VAL:HG12	1:F:15:LEU:HD21	1.97	0.45
1:F:275:ARG:HG3	1:F:275:ARG:HH11	1.80	0.45
2:G:44:GLU:OE1	2:G:68:MET:CG	2.63	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:82:TRP:O	1:F:85:PHE:HB3	2.16	0.45
2:G:44:GLU:OE2	2:G:68:MET:CB	2.64	0.45
2:G:73:SER:N	2:G:74:PRO:CD	2.80	0.45
4:X:8:3DR:P	4:X:9:3DR:H1'2	2.49	0.45
2:G:87:LEU:HD11	2:G:171:ILE:HD11	1.99	0.45
1:F:38:ALA:HB2	1:F:280:GLU:HA	1.99	0.44
1:F:75:ARG:NH1	1:F:280:GLU:OE1	2.50	0.44
1:F:215:ARG:HG3	1:F:215:ARG:HH11	1.82	0.44
4:X:8:3DR:C5'	4:X:9:3DR:H5''	2.47	0.44
2:G:20:VAL:HB	2:G:50:LEU:HD23	1.98	0.44
4:X:2:3DR:P	4:X:3:3DR:H1'2	2.58	0.44
1:F:75:ARG:NH2	1:F:230:ASP:CG	2.76	0.44
1:F:76:HIS:HA	1:F:233:ASN:HD21	1.81	0.44
2:G:9:HIS:HB2	2:G:82:SER:HB2	1.98	0.44
1:F:48:GLY:HA3	1:F:411:ALA:HB2	2.00	0.44
1:F:50:LEU:HD11	1:F:290:ILE:HD12	1.98	0.44
1:F:144:ARG:HG2	1:F:144:ARG:HH11	1.83	0.44
2:G:68:MET:SD	2:G:173:THR:O	2.76	0.44
4:X:2:3DR:OP1	4:X:3:3DR:H1'2	2.18	0.44
1:F:274:PRO:O	1:F:275:ARG:C	2.61	0.44
1:F:98:VAL:HG23	1:F:148:ASN:ND2	2.32	0.44
2:G:136:ASP:OD1	2:G:137:CYS:N	2.40	0.43
1:F:44:LEU:HD12	1:F:45:ASP:N	2.33	0.43
1:F:84:GLN:O	1:F:85:PHE:C	2.61	0.43
1:F:287:LEU:HD12	1:F:287:LEU:H	1.77	0.43
1:F:127:SER:O	1:F:128:TYR:C	2.58	0.43
1:F:330:SER:HB3	1:F:343:LYS:HD3	2.00	0.43
1:F:207:GLN:O	1:F:208:GLU:C	2.61	0.43
1:F:409:LYS:C	1:F:409:LYS:CD	2.92	0.43
2:G:42:THR:CB	2:G:69:LYS:HB3	2.48	0.43
1:F:77:VAL:HG21	1:F:123:PHE:HA	2.00	0.43
1:F:317:ASP:O	1:F:321:ILE:HG13	2.18	0.43
1:F:153:ASN:O	1:F:154:GLU:C	2.61	0.43
3:J:5:ARG:HG2	3:J:5:ARG:NH1	2.32	0.43
1:F:213:MET:SD	1:F:219:VAL:HA	2.59	0.42
2:G:186:LEU:HA	2:G:186:LEU:HD23	1.86	0.42
1:F:124:LEU:HD23	1:F:124:LEU:HA	1.81	0.42
1:F:393:GLN:C	1:F:395:TYR:N	2.76	0.42
4:X:1:3DR:OP2	4:X:7:3DR:H2''	2.18	0.42
1:F:241:THR:HA	4:X:10:3DR:H2''	2.01	0.42
1:F:180:LEU:HD23	1:F:180:LEU:HA	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:234:ARG:HA	1:F:235:PRO:HD3	1.94	0.42
2:G:50:LEU:HD23	2:G:50:LEU:HA	1.84	0.42
2:G:128:SER:OG	2:G:131:VAL:HG12	2.20	0.42
1:F:241:THR:HG22	1:F:242:ASP:N	2.35	0.42
1:F:320:LEU:O	1:F:324:LEU:HG	2.19	0.42
2:G:126:VAL:HG22	2:G:133:SER:CB	2.43	0.42
4:X:1:3DR:C5'	4:X:7:3DR:O3'	2.68	0.42
2:G:41:ILE:CG1	2:G:46:ALA:CB	2.97	0.42
2:G:56:VAL:CG1	2:G:60:SER:HA	2.49	0.42
3:J:24:PHE:HA	4:X:2:3DR:O4'	2.20	0.42
1:F:213:MET:HB3	1:F:219:VAL:HG22	2.02	0.42
1:F:102:ARG:HG3	1:F:102:ARG:HH11	1.85	0.41
1:F:362:LYS:HD3	1:F:362:LYS:C	2.45	0.41
2:G:45:ARG:CD	2:G:174:THR:HB	2.47	0.41
1:F:267:GLN:OE1	1:F:267:GLN:HA	2.20	0.41
2:G:96:ILE:HA	2:G:124:MET:CE	2.20	0.41
1:F:166:LEU:H	1:F:166:LEU:CD2	2.33	0.41
1:F:296:SER:C	1:F:298:LEU:H	2.27	0.41
1:F:299:GLU:HB2	1:F:335:PHE:CE1	2.55	0.41
4:X:2:3DR:OP1	4:X:3:3DR:C1'	2.68	0.41
3:J:17:TRP:HD1	4:X:3:3DR:C4'	2.33	0.41
1:F:168:ASN:OD1	1:F:168:ASN:C	2.64	0.41
1:F:166:LEU:N	1:F:166:LEU:CD2	2.83	0.41
1:F:170:TRP:HZ2	1:F:294:PRO:HB2	1.86	0.41
2:G:45:ARG:HD2	2:G:45:ARG:HA	1.77	0.41
2:G:184:GLN:HA	2:G:185:PRO:HD2	1.81	0.41
2:G:17:ARG:NH1	2:G:17:ARG:CG	2.82	0.41
1:F:423:THR:HG23	1:F:426:ASP:CG	2.46	0.41
2:G:4:ASN:HD22	2:G:5:PHE:N	2.19	0.41
3:J:24:PHE:CD1	4:X:2:3DR:H1'2	2.43	0.40
1:F:153:ASN:ND2	1:F:156:ASP:OD2	2.55	0.40
1:F:242:ASP:H	4:X:10:3DR:C3'	2.34	0.40
1:F:242:ASP:OD2	4:X:9:3DR:H1'1	2.20	0.40
1:F:287:LEU:N	1:F:287:LEU:CD1	2.73	0.40
1:F:374:ASP:O	1:F:375:ALA:C	2.64	0.40
2:G:33:ILE:HA	2:G:34:PRO:HD2	1.87	0.40

There are no symmetry-related clashes.

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	F	420/431 (97%)	381 (91%)	33 (8%)	6 (1%)	9	38
2	G	185/187 (99%)	160 (86%)	21 (11%)	4 (2%)	5	30
3	J	22/24 (92%)	20 (91%)	1 (4%)	1 (4%)	2	17
All	All	627/642 (98%)	561 (90%)	55 (9%)	11 (2%)	6	34

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	25	LEU
2	G	36	ILE
2	G	43	ALA
1	F	106	HIS
2	G	42	THR
2	G	58	THR
3	J	8	PRO
1	F	382	ASP
1	F	154	GLU
1	F	26	GLY
1	F	90	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	F	368/376 (98%)	338 (92%)	30 (8%)	10	34

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	G	156/156 (100%)	143 (92%)	13 (8%)	10	34
3	J	19/19 (100%)	18 (95%)	1 (5%)	20	47
All	All	543/551 (98%)	499 (92%)	44 (8%)	11	35

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	11	GLU
1	F	25	LEU
1	F	31	VAL
1	F	41	SER
1	F	165	HIS
1	F	174	LEU
1	F	182	GLU
1	F	205	THR
1	F	209	ARG
1	F	233	ASN
1	F	237	LEU
1	F	239	MET
1	F	254	ASP
1	F	256	SER
1	F	265	VAL
1	F	268	THR
1	F	287	LEU
1	F	290	ILE
1	F	310	THR
1	F	313	ASP
1	F	327	ARG
1	F	362	LYS
1	F	364	HIS
1	F	378	THR
1	F	383	ASN
1	F	393	GLN
1	F	401	SER
1	F	403	GLN
1	F	423	THR
1	F	431	SER
2	G	24	VAL
2	G	40	SER
2	G	41	ILE
2	G	44	GLU

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Mol	Chain	Res	Type
2	G	45	ARG
2	G	68	MET
2	G	78	ASN
2	G	123	GLU
2	G	155	VAL
2	G	170	THR
2	G	173	THR
2	G	175	GLN
2	G	183	LEU
3	J	2	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	74	HIS
1	F	81	GLN
1	F	91	ASN
1	F	101	ASN
1	F	126	GLN
1	F	130	ASN
1	F	148	ASN
1	F	151	ASN
1	F	153	ASN
1	F	165	HIS
1	F	233	ASN
1	F	255	GLN
1	F	300	HIS
1	F	383	ASN
1	F	392	HIS
1	F	393	GLN
1	F	402	GLN
1	F	403	GLN
1	F	410	GLN
1	F	414	ASN
2	G	3	GLN
2	G	4	ASN
2	G	37	ASN
2	G	55	ASN
2	G	78	ASN
2	G	119	HIS
2	G	143	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

10 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	3DR	X	7	-	8,11,12	0.71	0	7,14,17	0.62	0
4	3DR	X	2	4	8,11,12	0.55	0	7,14,17	0.92	0
4	3DR	X	4	4	8,11,12	1.29	1 (12%)	7,14,17	0.68	0
4	3DR	X	6	-	8,11,12	0.55	0	7,14,17	0.92	0
4	3DR	X	8	-	8,11,12	0.68	0	7,14,17	2.66	4 (57%)
4	3DR	X	5	-	8,11,12	0.68	0	7,14,17	2.66	4 (57%)
4	3DR	X	10	-	8,11,12	0.70	0	7,14,17	0.62	0
4	3DR	X	3	4	8,11,12	0.70	0	7,14,17	0.62	0
4	3DR	X	1	4	8,11,12	0.69	0	7,14,17	2.64	4 (57%)
4	3DR	X	9	-	8,11,12	0.68	0	7,14,17	2.65	4 (57%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	3DR	X	7	-	-	1/3/15/16	0/1/1/1
4	3DR	X	2	4	1/1/3/3	3/3/15/16	0/1/1/1
4	3DR	X	4	4	-	2/3/15/16	0/1/1/1
4	3DR	X	6	-	-	2/3/15/16	0/1/1/1
4	3DR	X	8	-	-	2/3/15/16	0/1/1/1
4	3DR	X	5	-	-	0/3/15/16	0/1/1/1
4	3DR	X	10	-	-	2/3/15/16	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	3DR	X	3	4	-	2/3/15/16	0/1/1/1
4	3DR	X	1	4	-	2/3/15/16	0/1/1/1
4	3DR	X	9	-	-	2/3/15/16	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	X	4	3DR	C5'-C4'	2.76	1.59	1.51

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	X	5	3DR	O3'-C3'-C2'	4.65	122.35	111.43
4	X	1	3DR	O3'-C3'-C2'	4.64	122.32	111.43
4	X	8	3DR	O3'-C3'-C2'	4.63	122.30	111.43
4	X	9	3DR	O3'-C3'-C2'	4.63	122.28	111.43
4	X	5	3DR	C1'-C2'-C3'	3.21	106.70	103.26
4	X	8	3DR	C1'-C2'-C3'	3.21	106.70	103.26
4	X	9	3DR	C1'-C2'-C3'	3.17	106.66	103.26
4	X	1	3DR	C1'-C2'-C3'	3.15	106.63	103.26
4	X	9	3DR	O3'-C3'-C4'	-2.92	98.99	110.07
4	X	8	3DR	O3'-C3'-C4'	-2.90	99.07	110.07
4	X	1	3DR	O3'-C3'-C4'	-2.89	99.10	110.07
4	X	5	3DR	O3'-C3'-C4'	-2.89	99.13	110.07
4	X	5	3DR	C1'-O4'-C4'	2.49	112.28	108.37
4	X	8	3DR	C1'-O4'-C4'	2.45	112.23	108.37
4	X	1	3DR	C1'-O4'-C4'	2.43	112.19	108.37
4	X	9	3DR	C1'-O4'-C4'	2.41	112.17	108.37

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	X	2	3DR	C4'

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	X	1	3DR	C3'-C4'-C5'-O5'
4	X	6	3DR	O4'-C4'-C5'-O5'
4	X	8	3DR	C3'-C4'-C5'-O5'
4	X	9	3DR	C3'-C4'-C5'-O5'

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Mol	Chain	Res	Type	Atoms
4	X	1	3DR	O4'-C4'-C5'-O5'
4	X	2	3DR	C3'-C4'-C5'-O5'
4	X	8	3DR	O4'-C4'-C5'-O5'
4	X	9	3DR	O4'-C4'-C5'-O5'
4	X	2	3DR	O4'-C4'-C5'-O5'
4	X	6	3DR	C3'-C4'-C5'-O5'
4	X	7	3DR	O4'-C4'-C5'-O5'
4	X	4	3DR	O4'-C4'-C5'-O5'
4	X	4	3DR	C3'-C4'-C5'-O5'
4	X	3	3DR	C3'-C4'-C5'-O5'
4	X	10	3DR	C3'-C4'-C5'-O5'
4	X	2	3DR	C4'-C5'-O5'-P
4	X	3	3DR	O4'-C4'-C5'-O5'
4	X	10	3DR	O4'-C4'-C5'-O5'

There are no ring outliers.

9 monomers are involved in 79 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	X	7	3DR	14	0
4	X	2	3DR	26	0
4	X	4	3DR	5	0
4	X	6	3DR	2	0
4	X	8	3DR	11	0
4	X	10	3DR	18	0
4	X	3	3DR	12	0
4	X	1	3DR	10	0
4	X	9	3DR	21	0

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

The following chains have linkage breaks:

Mol	Chain	Number of breaks
4	X	7

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	X	4:3DR	O3'	5:3DR	P	20.26
1	X	7:3DR	O3'	8:3DR	P	19.47
1	X	5:3DR	O3'	6:3DR	P	15.55
1	X	8:3DR	O3'	9:3DR	P	6.19
1	X	9:3DR	O3'	10:3DR	P	5.27
1	X	6:3DR	O3'	7:3DR	P	2.30
1	X	2:3DR	O3'	3:3DR	P	1.77

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	F	422/431 (97%)	-0.53	0 100 100	1, 4, 36, 91	0
2	G	187/187 (100%)	-0.48	0 100 100	1, 17, 52, 96	0
3	J	24/24 (100%)	-0.38	0 100 100	1, 30, 91, 103	0
4	X	0/10	-	-	-	-
All	All	633/652 (97%)	-0.51	0 100 100	1, 8, 47, 103	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [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	3DR	X	3	11/12	0.97	0.14	1,1,1,1	0
4	3DR	X	1	11/12	0.98	0.10	1,1,1,1	0
4	3DR	X	7	11/12	0.98	0.09	1,1,1,1	0
4	3DR	X	9	11/12	0.98	0.08	1,1,1,1	0
4	3DR	X	5	11/12	0.99	0.07	1,1,1,1	0
4	3DR	X	6	11/12	0.99	0.08	1,1,1,1	0
4	3DR	X	2	11/12	0.99	0.09	1,1,1,1	0
4	3DR	X	8	11/12	0.99	0.06	1,1,1,1	0
4	3DR	X	4	11/12	0.99	0.07	1,1,1,1	0
4	3DR	X	10	11/12	0.99	0.07	1,1,1,1	0

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.