



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 10:58 AM UTC

PDB ID : 3BCP / pdb_00003bcp
Title : Crystal Structure of The Swapped non covalent form of P19A/L28Q/N67D
BS-RNase
Authors : Merlino, A.; Ercole, C.; Picone, D.; Pizzo, E.; Mazzarella, L.; Sica, F.
Deposited on : 2007-11-13
Resolution : 2.57 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

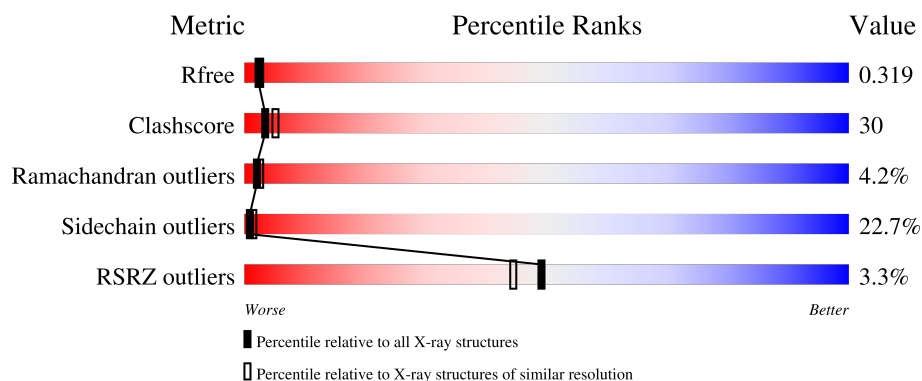
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.57 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	4770 (2.60-2.56)
Clashscore	190562	5124 (2.60-2.56)
Ramachandran outliers	187476	5046 (2.60-2.56)
Sidechain outliers	187428	5046 (2.60-2.56)
RSRZ outliers	180081	4770 (2.60-2.56)

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	124	4% 39% 48% 11%
1	B	124	2% 43% 43% 13%
1	C	124	2% 44% 40% 15%
1	D	124	5% 46% 37% 15%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3854 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 Seminal ribonuclease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	124	Total	C	N	O	S	0	0	0
			948	572	173	188	15			
1	B	124	Total	C	N	O	S	0	0	0
			948	572	173	188	15			
1	C	124	Total	C	N	O	S	0	0	0
			948	572	173	188	15			
1	D	124	Total	C	N	O	S	0	0	0
			948	572	173	188	15			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	ALA	PRO	engineered mutation	UNP P00669
A	28	GLN	LEU	engineered mutation	UNP P00669
A	67	ASP	ASN	engineered mutation	UNP P00669
B	19	ALA	PRO	engineered mutation	UNP P00669
B	28	GLN	LEU	engineered mutation	UNP P00669
B	67	ASP	ASN	engineered mutation	UNP P00669
C	19	ALA	PRO	engineered mutation	UNP P00669
C	28	GLN	LEU	engineered mutation	UNP P00669
C	67	ASP	ASN	engineered mutation	UNP P00669
D	19	ALA	PRO	engineered mutation	UNP P00669
D	28	GLN	LEU	engineered mutation	UNP P00669
D	67	ASP	ASN	engineered mutation	UNP P00669

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	14	Total	O	0	0
			14	14		
2	B	19	Total	O	0	0
			19	19		

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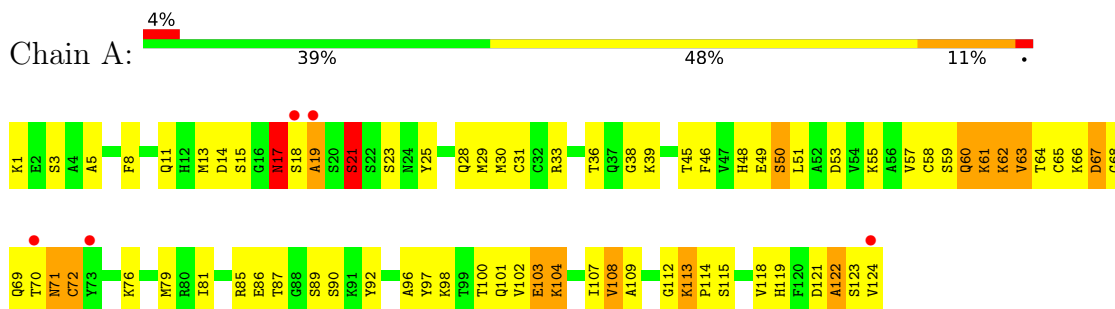
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	15	Total	O	0	0
			15	15		
2	D	14	Total	O	0	0
			14	14		

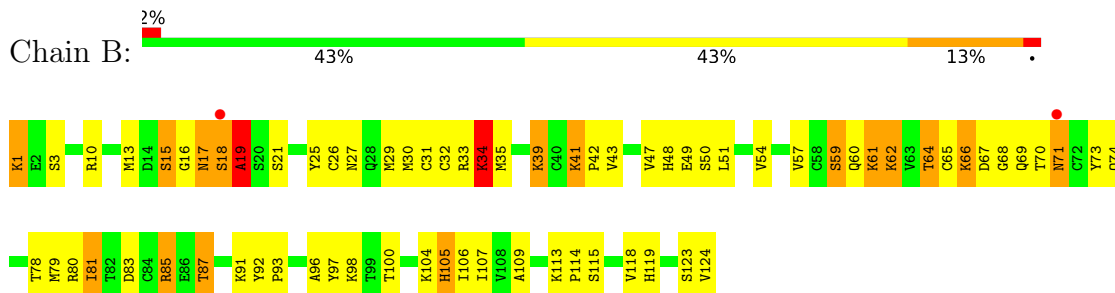
3 Residue-property plots [i](#)

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.

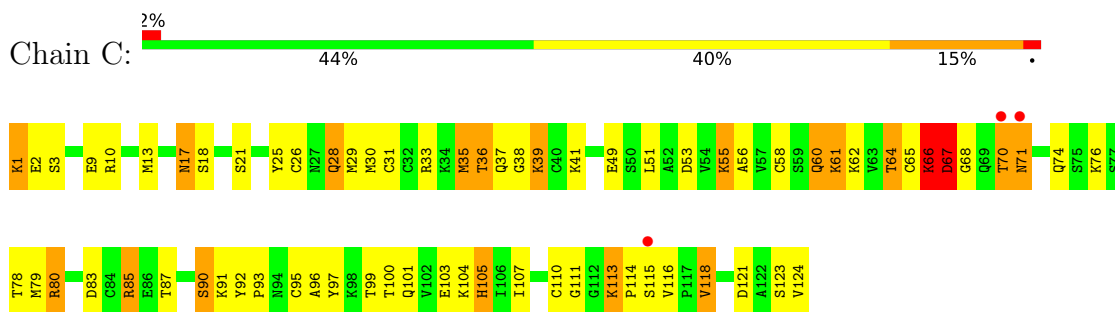
- Molecule 1: Seminal ribonuclease



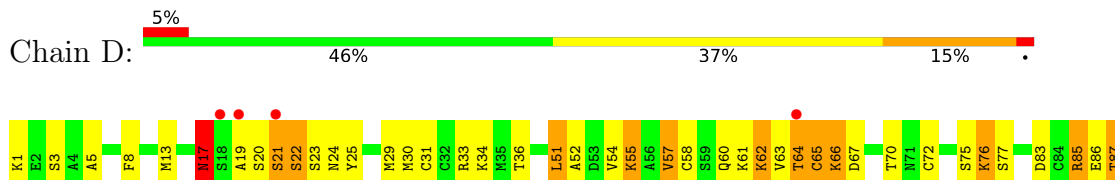
- Molecule 1: Seminal ribonuclease



- Molecule 1: Seminal ribonuclease



- Molecule 1: Seminal ribonuclease





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.72Å 86.40Å 99.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.57 20.00 – 2.57	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.57) 91.6 (20.00-2.57)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.41 (at 2.48Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.238 , 0.323 0.238 , 0.319	Depositor DCC
R_{free} test set	1162 reflections (6.52%)	wwPDB-VP
Wilson B-factor (Å ²)	42.1	Xtriage
Anisotropy	0.804	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 38.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3854	wwPDB-VP
Average B, all atoms (Å ²)	46.0	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 44.62 % 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 1.4867e-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

5.1 Standard geometry

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

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	0.54	0/941	0.94	2/1258 (0.2%)
1	B	0.50	0/941	0.96	4/1258 (0.3%)
1	C	0.62	2/941 (0.2%)	0.98	3/1258 (0.2%)
1	D	0.55	1/941 (0.1%)	1.03	5/1258 (0.4%)
All	All	0.56	3/3764 (0.1%)	0.98	14/5032 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	101	GLN	CD-OE1	5.68	1.34	1.23
1	C	17	ASN	CG-OD1	5.27	1.33	1.23
1	D	101	GLN	CD-OE1	5.15	1.33	1.23

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	19	ALA	N-CA-C	-9.69	101.46	113.28
1	D	66	LYS	N-CA-C	-7.45	104.32	113.41
1	B	41	LYS	CA-C-N	6.89	127.17	119.32
1	B	41	LYS	C-N-CA	6.89	127.17	119.32
1	A	19	ALA	N-CA-C	-6.19	104.54	113.21
1	C	105	HIS	N-CA-C	-5.96	101.67	110.48
1	D	118	VAL	N-CA-C	-5.61	107.82	113.20
1	D	104	LYS	N-CA-C	5.39	116.34	108.14
1	D	57	VAL	N-CA-C	-5.37	105.30	110.72
1	C	41	LYS	CA-C-N	5.28	124.92	119.05
1	C	41	LYS	C-N-CA	5.28	124.92	119.05
1	B	19	ALA	N-CA-C	-5.28	99.55	110.80
1	A	21	SER	N-CA-C	-5.14	105.79	111.71
1	B	105	HIS	N-CA-C	-5.01	103.79	110.55

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	948	0	925	64	0
1	B	948	0	925	69	0
1	C	948	0	925	70	0
1	D	948	0	925	59	0
2	A	14	0	0	3	0
2	B	19	0	0	1	0
2	C	15	0	0	3	0
2	D	14	0	0	1	0
All	All	3854	0	3700	225	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 30.

All (225) 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:D:104:LYS:HD2	1:D:123:SER:HB2	1.20	1.13
1:A:86:GLU:HB3	1:C:64:THR:HG21	1.31	1.11
1:A:86:GLU:HB3	1:C:64:THR:CG2	1.95	0.95
1:B:49:GLU:HG3	1:B:79:MET:HB3	1.60	0.83
1:B:27:ASN:HD21	1:B:97:TYR:H	1.28	0.78
1:A:57:VAL:HA	1:A:60:GLN:NE2	1.99	0.78
1:D:51:LEU:HD23	1:D:55:LYS:HE2	1.65	0.78
1:A:104:LYS:CD	1:A:123:SER:HB2	2.17	0.74
1:A:104:LYS:HD3	1:A:123:SER:HB2	1.68	0.74
1:D:65:CYS:HB3	1:D:121:ASP:OD2	1.87	0.74
1:A:98:LYS:CE	1:C:61:LYS:HD2	2.20	0.72
1:B:49:GLU:CG	1:B:79:MET:HB3	2.20	0.71
1:C:80:ARG:HH12	1:D:17:ASN:CB	2.04	0.70
1:A:87:THR:HG23	1:A:96:ALA:O	1.92	0.70
1:B:27:ASN:ND2	1:B:97:TYR:H	1.89	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:112:GLY:O	1:D:113:LYS:HB2	1.92	0.69
1:D:30:MET:HG3	1:D:97:TYR:CE2	2.27	0.69
1:B:39:LYS:HD3	1:B:39:LYS:N	2.09	0.68
1:D:104:LYS:CD	1:D:123:SER:HB2	2.12	0.68
1:C:30:MET:SD	1:C:35:MET:HE3	2.33	0.68
1:B:109:ALA:HB3	1:B:119:HIS:HB3	1.75	0.67
1:B:92:TYR:CD1	1:B:93:PRO:HA	2.30	0.67
1:B:79:MET:O	1:B:81:ILE:HG23	1.94	0.66
1:C:31:YCM:OZ1	1:C:31:YCM:HB3	1.95	0.66
1:B:85:ARG:HD2	2:B:658:HOH:O	1.95	0.65
1:D:54:VAL:O	1:D:57:VAL:HG23	1.97	0.65
1:A:86:GLU:O	1:C:64:THR:HG22	1.97	0.65
1:A:49:GLU:HG3	1:A:79:MET:HB3	1.79	0.64
1:A:18:SER:HB3	1:B:48:HIS:CE1	2.32	0.64
1:A:55:LYS:O	1:A:58:CYS:HB2	1.97	0.64
1:B:104:LYS:NZ	1:D:124:VAL:HA	2.13	0.64
1:D:109:ALA:HB3	1:D:119:HIS:HB3	1.80	0.64
1:D:51:LEU:O	1:D:55:LYS:HG2	1.98	0.64
1:D:87:THR:HG23	1:D:96:ALA:O	1.97	0.64
1:A:101:GLN:NE2	2:A:278:HOH:O	2.30	0.63
1:B:104:LYS:HZ2	1:D:124:VAL:HA	1.63	0.63
1:D:55:LYS:O	1:D:58:CYS:HB2	1.99	0.63
1:A:107:ILE:HD12	1:A:122:ALA:HB3	1.81	0.63
1:C:13:MET:HE2	1:D:51:LEU:HA	1.81	0.63
1:A:86:GLU:HG3	1:A:90:SER:CB	2.30	0.62
1:C:55:LYS:HE2	1:C:116:VAL:HG12	1.81	0.62
1:C:87:THR:HG23	1:C:96:ALA:C	2.25	0.61
1:D:60:GLN:HG2	1:D:76:LYS:HB2	1.81	0.61
1:B:87:THR:HG23	1:B:96:ALA:O	2.01	0.61
1:C:49:GLU:CG	1:C:79:MET:HB3	2.31	0.61
1:A:48:HIS:CE1	1:B:18:SER:HB3	2.36	0.61
1:C:39:LYS:HZ3	1:C:39:LYS:C	2.09	0.61
1:B:65:CYS:C	1:B:67:ASP:H	2.09	0.60
1:B:31:YCM:NZ2	1:B:31:YCM:HB3	2.17	0.60
1:D:25:TYR:CZ	1:D:29:MET:HG3	2.37	0.60
1:C:90:SER:HB2	1:C:96:ALA:HB3	1.85	0.59
1:A:19:ALA:O	1:A:23:SER:HB3	2.03	0.58
1:A:11:GLN:HG2	1:B:35:MET:SD	2.43	0.58
1:C:65:CYS:O	1:C:67:ASP:N	2.36	0.58
1:B:15:SER:C	1:B:17:ASN:H	2.10	0.58
1:B:104:LYS:HB3	1:B:124:VAL:C	2.29	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:65:CYS:C	1:C:67:ASP:N	2.60	0.58
1:A:109:ALA:HB3	1:A:119:HIS:HB3	1.85	0.58
1:D:111:GLY:O	1:D:115:SER:HA	2.04	0.58
1:C:53:ASP:O	1:C:56:ALA:HB3	2.04	0.58
1:A:19:ALA:O	1:A:23:SER:N	2.37	0.57
1:A:21:SER:HB3	1:B:21:SER:OG	2.04	0.57
1:D:104:LYS:NZ	1:D:123:SER:OG	2.30	0.57
1:C:87:THR:HG23	1:C:96:ALA:O	2.05	0.57
1:A:98:LYS:CD	1:C:61:LYS:HD2	2.36	0.56
1:A:51:LEU:HG	1:A:55:LYS:HD2	1.86	0.56
1:C:83:ASP:OD2	1:C:85:ARG:NH1	2.38	0.56
1:B:25:TYR:CE1	1:B:29:MET:HG3	2.40	0.56
1:A:50:SER:OG	1:A:53:ASP:HB2	2.06	0.56
1:A:25:TYR:CZ	1:A:29:MET:HG3	2.41	0.55
1:A:98:LYS:HD2	1:C:61:LYS:HD2	1.88	0.55
1:C:65:CYS:C	1:C:67:ASP:H	2.12	0.55
1:C:80:ARG:HH12	1:D:17:ASN:HB2	1.71	0.55
1:D:91:LYS:HE3	1:D:92:TYR:O	2.06	0.55
1:C:31:YCM:HA	1:C:36:THR:OG1	2.07	0.55
1:C:9:GLU:HB2	2:C:306:HOH:O	2.07	0.55
1:A:49:GLU:CG	1:A:79:MET:HB3	2.37	0.54
1:B:73:TYR:O	1:B:107:ILE:HA	2.08	0.54
1:C:80:ARG:HH12	1:D:17:ASN:HB3	1.72	0.54
1:C:28:GLN:HG3	1:C:29:MET:N	2.23	0.54
1:A:30:MET:HG3	1:A:97:TYR:CE2	2.42	0.54
1:C:1:LYS:HA	1:C:1:LYS:HZ3	1.72	0.54
1:B:59:SER:O	1:B:59:SER:OG	2.25	0.54
1:B:78:THR:HG22	1:B:105:HIS:CD2	2.43	0.53
1:B:18:SER:O	1:B:19:ALA:HB2	2.09	0.53
1:C:110:CYS:HA	1:C:118:VAL:HG22	1.90	0.53
1:C:71:ASN:O	1:C:71:ASN:CG	2.51	0.53
1:D:60:GLN:OE1	1:D:75:SER:HA	2.09	0.53
1:D:57:VAL:HG13	1:D:75:SER:HA	1.91	0.53
1:B:124:VAL:HG22	1:D:102:VAL:HG11	1.90	0.52
1:A:17:ASN:HB2	1:B:80:ARG:HH21	1.75	0.52
1:A:51:LEU:HA	1:B:13:MET:HE2	1.91	0.52
1:A:98:LYS:HD2	1:C:61:LYS:HZ3	1.74	0.52
1:B:104:LYS:HB3	1:B:124:VAL:OXT	2.09	0.52
1:A:113:LYS:HA	1:A:114:PRO:C	2.34	0.52
1:C:65:CYS:HA	1:C:121:ASP:OD2	2.11	0.51
1:A:5:ALA:O	1:A:8:PHE:HB3	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1:LYS:HB2	2:C:636:HOH:O	2.11	0.51
1:C:95:CYS:HB3	1:C:97:TYR:HE1	1.75	0.51
1:A:72:CYS:HA	1:A:108:VAL:O	2.11	0.50
1:B:92:TYR:CG	1:B:93:PRO:HA	2.47	0.50
1:A:86:GLU:CG	1:A:90:SER:HB3	2.42	0.50
1:C:55:LYS:HE2	1:C:116:VAL:CG1	2.42	0.50
1:C:49:GLU:HG2	1:C:79:MET:HB3	1.94	0.49
1:B:30:MET:SD	1:B:35:MET:HE3	2.53	0.49
1:B:16:GLY:C	1:B:17:ASN:HD22	2.20	0.49
1:B:57:VAL:CG2	1:B:79:MET:HE1	2.43	0.49
1:C:49:GLU:OE2	1:C:80:ARG:HD2	2.12	0.49
1:A:30:MET:O	1:A:36:THR:HG23	2.12	0.49
1:C:58:CYS:HB3	1:C:115:SER:OG	2.13	0.49
1:B:47:VAL:HG12	1:B:49:GLU:H	1.77	0.49
1:C:49:GLU:HG3	1:C:79:MET:CG	2.43	0.48
1:D:5:ALA:O	1:D:8:PHE:HB3	2.14	0.48
1:D:31:YCM:HA	1:D:36:THR:HG1	1.79	0.48
1:D:62:LYS:HD3	1:D:72:CYS:O	2.14	0.48
1:D:34:LYS:HA	2:D:268:HOH:O	2.14	0.48
1:A:98:LYS:HE2	1:C:61:LYS:HD2	1.96	0.48
1:C:124:VAL:O	1:C:124:VAL:HG13	2.13	0.48
1:A:79:MET:O	1:A:103:GLU:HA	2.14	0.48
1:D:83:ASP:OD2	1:D:85:ARG:NH1	2.47	0.48
1:B:39:LYS:HD3	1:B:39:LYS:H	1.75	0.47
1:D:64:THR:HG23	1:D:65:CYS:O	2.14	0.47
1:A:17:ASN:CB	1:B:80:ARG:HH21	2.28	0.47
1:A:30:MET:HG3	1:A:97:TYR:CZ	2.49	0.47
1:A:65:CYS:SG	1:A:69:GLN:HB2	2.54	0.47
1:B:83:ASP:OD2	1:B:85:ARG:NH1	2.47	0.47
1:B:113:LYS:HA	1:B:114:PRO:C	2.39	0.47
1:A:17:ASN:CB	1:B:80:ARG:NH2	2.77	0.47
1:C:1:LYS:HZ2	1:C:2:GLU:N	2.13	0.47
1:C:111:GLY:O	1:C:115:SER:HA	2.15	0.47
1:B:83:ASP:CG	1:D:124:VAL:HG21	2.40	0.47
1:A:87:THR:HG23	1:A:96:ALA:C	2.39	0.47
1:D:30:MET:HG3	1:D:97:TYR:CD2	2.50	0.46
1:B:25:TYR:CZ	1:B:29:MET:HG3	2.51	0.46
1:A:61:LYS:HG3	1:A:62:LYS:N	2.25	0.46
1:B:123:SER:O	1:B:124:VAL:HB	2.15	0.46
1:A:33:ARG:O	1:B:10:ARG:NH1	2.43	0.46
1:A:45:THR:HG23	1:A:81:ILE:HD12	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:ARG:NH2	2:A:298:HOH:O	2.45	0.45
1:B:87:THR:HG23	1:B:96:ALA:C	2.41	0.45
1:A:8:PHE:HA	2:A:205:HOH:O	2.16	0.45
1:C:37:GLN:C	1:C:39:LYS:H	2.24	0.45
1:C:104:LYS:HB3	1:C:124:VAL:C	2.41	0.45
1:D:104:LYS:HD2	1:D:123:SER:CB	2.15	0.45
1:D:110:CYS:HB3	1:D:115:SER:OG	2.17	0.45
1:A:62:LYS:HG3	1:A:63:VAL:N	2.31	0.45
1:A:63:VAL:O	1:A:71:ASN:O	2.35	0.45
1:C:51:LEU:HA	1:D:13:MET:HE2	1.98	0.45
1:C:78:THR:HB	1:C:104:LYS:C	2.42	0.45
1:D:92:TYR:CD1	1:D:93:PRO:HA	2.52	0.44
1:A:28:GLN:O	1:A:31:YCM:HB2	2.18	0.44
1:D:31:YCM:HA	1:D:36:THR:OG1	2.16	0.44
1:A:17:ASN:N	1:A:17:ASN:HD22	2.15	0.44
1:B:79:MET:HE3	1:B:106:ILE:HD11	1.98	0.44
1:C:33:ARG:HA	1:C:33:ARG:HD3	1.73	0.44
1:D:105:HIS:HB2	1:D:124:VAL:O	2.18	0.44
1:B:15:SER:C	1:B:17:ASN:N	2.75	0.44
1:B:33:ARG:O	1:B:34:LYS:C	2.60	0.44
1:C:49:GLU:HG3	1:C:79:MET:HB3	2.00	0.44
1:C:92:TYR:CE2	1:C:93:PRO:HB3	2.52	0.44
1:C:123:SER:O	1:C:124:VAL:HB	2.18	0.44
1:D:33:ARG:O	1:D:34:LYS:HB2	2.17	0.44
1:D:87:THR:HG23	1:D:96:ALA:C	2.42	0.44
1:A:38:GLY:C	1:A:92:TYR:CE1	2.96	0.44
1:C:13:MET:HE2	1:D:51:LEU:HG	1.99	0.44
1:B:65:CYS:C	1:B:67:ASP:N	2.75	0.44
1:B:70:THR:O	1:B:71:ASN:HB2	2.17	0.44
1:C:13:MET:CE	1:D:51:LEU:HA	2.47	0.44
1:C:78:THR:HG22	1:C:105:HIS:CD2	2.52	0.44
1:D:107:ILE:HD12	1:D:122:ALA:HB3	1.98	0.44
1:A:17:ASN:HB3	1:B:80:ARG:NH2	2.33	0.43
1:B:61:LYS:HG2	1:D:98:LYS:HE2	2.00	0.43
1:D:90:SER:HB2	1:D:96:ALA:HB3	2.00	0.43
1:A:86:GLU:CG	1:A:90:SER:CB	2.97	0.43
1:A:124:VAL:HG13	1:A:124:VAL:O	2.19	0.43
1:B:41:LYS:HA	1:B:42:PRO:HD2	1.78	0.43
1:C:49:GLU:HG3	1:C:79:MET:HG2	2.00	0.43
1:D:86:GLU:HG3	1:D:90:SER:OG	2.17	0.43
1:A:45:THR:CG2	1:A:81:ILE:HD12	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:15:SER:O	1:B:17:ASN:N	2.43	0.43
1:B:62:LYS:HD2	1:B:73:TYR:OH	2.18	0.43
1:B:18:SER:O	1:B:19:ALA:CB	2.66	0.43
1:B:26:CYS:O	1:B:30:MET:HB2	2.19	0.43
1:C:49:GLU:HB3	2:C:291:HOH:O	2.19	0.43
1:D:61:LYS:HE2	1:D:63:VAL:CG1	2.49	0.43
1:A:46:PHE:O	1:A:81:ILE:HA	2.18	0.43
1:B:32:YCM:C	1:B:34:LYS:H	2.31	0.43
1:C:10:ARG:NH1	1:D:34:LYS:HB2	2.34	0.43
1:B:1:LYS:HA	1:B:1:LYS:HZ3	1.83	0.43
1:A:13:MET:HE2	1:B:51:LEU:HA	2.00	0.42
1:A:66:LYS:HE2	1:A:121:ASP:OD1	2.19	0.42
1:B:31:YCM:HB3	1:B:31:YCM:HZ21	1.80	0.42
1:C:36:THR:HG22	1:C:95:CYS:SG	2.59	0.42
1:A:102:VAL:HG11	1:C:124:VAL:HG22	2.02	0.42
1:A:98:LYS:HD2	1:C:61:LYS:NZ	2.34	0.42
1:A:104:LYS:HD2	1:A:123:SER:HB2	1.98	0.42
1:C:74:GLN:HE21	1:C:107:ILE:HD11	1.83	0.42
1:C:10:ARG:NH1	1:D:33:ARG:O	2.50	0.42
1:C:65:CYS:O	1:C:66:LYS:C	2.62	0.42
1:B:65:CYS:O	1:B:67:ASP:N	2.51	0.41
1:B:54:VAL:O	1:B:57:VAL:HB	2.20	0.41
1:D:51:LEU:O	1:D:52:ALA:C	2.63	0.41
1:B:31:YCM:HZ21	1:B:31:YCM:CB	2.33	0.41
1:D:20:SER:HA	1:D:23:SER:HB3	2.01	0.41
1:C:113:LYS:HA	1:C:114:PRO:HA	1.94	0.41
1:A:69:GLN:HB3	1:A:71:ASN:OD1	2.20	0.41
1:D:57:VAL:HG13	1:D:75:SER:CA	2.50	0.41
1:A:14:ASP:O	1:B:48:HIS:HA	2.20	0.41
1:B:64:THR:HG22	1:B:65:CYS:O	2.21	0.41
1:D:20:SER:C	1:D:22:SER:N	2.78	0.41
1:C:1:LYS:NZ	1:C:2:GLU:H	2.19	0.41
1:C:26:CYS:HB2	1:C:99:THR:OG1	2.20	0.41
1:C:38:GLY:C	1:C:39:LYS:HG3	2.45	0.41
1:C:21:SER:OG	1:D:21:SER:HB2	2.21	0.41
1:C:60:GLN:HB3	1:C:61:LYS:H	1.74	0.41
1:D:124:VAL:O	1:D:124:VAL:HG13	2.21	0.41
1:B:79:MET:CE	1:B:106:ILE:HD11	2.52	0.40
1:B:31:YCM:HZ21	1:B:31:YCM:HA	1.85	0.40
1:D:55:LYS:HG2	1:D:55:LYS:H	1.60	0.40
1:D:121:ASP:O	1:D:122:ALA:HB2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:74:GLN:HA	1:B:106:ILE:O	2.22	0.40
1:C:66:LYS:H	1:C:66:LYS:HG3	1.43	0.40
1:B:43:VAL:HG22	1:B:85:ARG:HG3	2.02	0.40
1:C:25:TYR:CZ	1:C:29:MET:HG3	2.56	0.40
1:D:61:LYS:CG	1:D:62:LYS:N	2.85	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	120/124 (97%)	105 (88%)	8 (7%)	7 (6%)	1	1
1	B	120/124 (97%)	104 (87%)	11 (9%)	5 (4%)	2	2
1	C	120/124 (97%)	101 (84%)	14 (12%)	5 (4%)	2	2
1	D	120/124 (97%)	105 (88%)	12 (10%)	3 (2%)	4	7
All	All	480/496 (97%)	415 (86%)	45 (9%)	20 (4%)	2	2

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	71	ASN
1	B	19	ALA
1	B	71	ASN
1	C	17	ASN
1	C	67	ASP
1	D	113	LYS
1	A	72	CYS
1	B	66	LYS
1	C	66	LYS
1	C	68	GLY

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Mol	Chain	Res	Type
1	C	70	THR
1	A	17	ASN
1	A	67	ASP
1	A	68	GLY
1	B	68	GLY
1	D	17	ASN
1	D	122	ALA
1	B	34	LYS
1	A	122	ALA
1	A	112	GLY

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	107/107 (100%)	83 (78%)	24 (22%)	1	1
1	B	107/107 (100%)	84 (78%)	23 (22%)	1	1
1	C	107/107 (100%)	82 (77%)	25 (23%)	1	1
1	D	107/107 (100%)	82 (77%)	25 (23%)	1	1
All	All	428/428 (100%)	331 (77%)	97 (23%)	1	1

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	LYS
1	A	3	SER
1	A	15	SER
1	A	17	ASN
1	A	21	SER
1	A	39	LYS
1	A	50	SER
1	A	59	SER
1	A	60	GLN
1	A	61	LYS
1	A	62	LYS

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Mol	Chain	Res	Type
1	A	63	VAL
1	A	64	THR
1	A	67	ASP
1	A	70	THR
1	A	76	LYS
1	A	89	SER
1	A	100	THR
1	A	103	GLU
1	A	104	LYS
1	A	108	VAL
1	A	113	LYS
1	A	115	SER
1	A	118	VAL
1	B	1	LYS
1	B	3	SER
1	B	15	SER
1	B	17	ASN
1	B	18	SER
1	B	34	LYS
1	B	39	LYS
1	B	50	SER
1	B	59	SER
1	B	60	GLN
1	B	61	LYS
1	B	62	LYS
1	B	64	THR
1	B	66	LYS
1	B	69	GLN
1	B	81	ILE
1	B	85	ARG
1	B	87	THR
1	B	91	LYS
1	B	98	LYS
1	B	100	THR
1	B	115	SER
1	B	118	VAL
1	C	1	LYS
1	C	3	SER
1	C	18	SER
1	C	28	GLN
1	C	35	MET
1	C	36	THR

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Mol	Chain	Res	Type
1	C	39	LYS
1	C	55	LYS
1	C	60	GLN
1	C	61	LYS
1	C	62	LYS
1	C	64	THR
1	C	66	LYS
1	C	67	ASP
1	C	70	THR
1	C	71	ASN
1	C	76	LYS
1	C	80	ARG
1	C	85	ARG
1	C	90	SER
1	C	91	LYS
1	C	100	THR
1	C	103	GLU
1	C	113	LYS
1	C	118	VAL
1	D	1	LYS
1	D	3	SER
1	D	17	ASN
1	D	21	SER
1	D	22	SER
1	D	24	ASN
1	D	51	LEU
1	D	55	LYS
1	D	62	LYS
1	D	64	THR
1	D	65	CYS
1	D	66	LYS
1	D	67	ASP
1	D	70	THR
1	D	76	LYS
1	D	77	SER
1	D	85	ARG
1	D	87	THR
1	D	90	SER
1	D	91	LYS
1	D	100	THR
1	D	102	VAL
1	D	104	LYS

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Mol	Chain	Res	Type
1	D	107	ILE
1	D	118	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	ASN
1	A	101	GLN
1	A	105	HIS
1	B	17	ASN
1	B	27	ASN
1	B	69	GLN
1	B	74	GLN
1	B	94	ASN
1	B	101	GLN
1	C	69	GLN
1	C	74	GLN
1	C	119	HIS
1	D	17	ASN
1	D	94	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

8 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	$\# Z > 2$
1	YCM	D	32	1	7,9,10	0.51	0	5,10,12	0.68	0
1	YCM	C	31	1	7,9,10	0.53	0	5,10,12	0.24	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	YCM	B	32	1	7,9,10	0.53	0	5,10,12	0.37	0
1	YCM	A	32	1	7,9,10	0.66	0	5,10,12	0.67	0
1	YCM	D	31	1	7,9,10	0.65	0	5,10,12	0.29	0
1	YCM	A	31	1	7,9,10	0.63	0	5,10,12	0.27	0
1	YCM	C	32	1	7,9,10	0.42	0	5,10,12	0.29	0
1	YCM	B	31	1	7,9,10	0.62	0	5,10,12	0.73	0

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
1	YCM	D	32	1	-	3/6/8/10	-
1	YCM	C	31	1	-	0/6/8/10	-
1	YCM	B	32	1	-	2/6/8/10	-
1	YCM	A	32	1	-	2/6/8/10	-
1	YCM	D	31	1	-	1/6/8/10	-
1	YCM	A	31	1	-	0/6/8/10	-
1	YCM	C	32	1	-	2/6/8/10	-
1	YCM	B	31	1	-	1/6/8/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	32	YCM	SG-CD-CE-NZ2
1	B	31	YCM	SG-CD-CE-NZ2
1	B	32	YCM	SG-CD-CE-OZ1
1	B	32	YCM	SG-CD-CE-NZ2
1	C	32	YCM	SG-CD-CE-OZ1
1	C	32	YCM	SG-CD-CE-NZ2
1	D	32	YCM	SG-CD-CE-OZ1
1	D	32	YCM	SG-CD-CE-NZ2
1	A	32	YCM	SG-CD-CE-OZ1
1	D	32	YCM	CA-CB-SG-CD
1	D	31	YCM	SG-CD-CE-OZ1

There are no ring outliers.

5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	31	YCM	2	0
1	B	32	YCM	1	0
1	D	31	YCM	2	0
1	A	31	YCM	1	0
1	B	31	YCM	4	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 [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	122/124 (98%)	0.48	5 (4%) 41 37	27, 47, 75, 79	0
1	B	122/124 (98%)	0.35	2 (1%) 70 67	27, 45, 59, 66	0
1	C	122/124 (98%)	0.26	3 (2%) 58 54	25, 43, 57, 65	0
1	D	122/124 (98%)	0.51	6 (4%) 35 30	30, 50, 73, 80	0
All	All	488/496 (98%)	0.40	16 (3%) 49 44	25, 46, 70, 80	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	18	SER	4.8
1	A	19	ALA	4.0
1	C	71	ASN	3.7
1	C	70	THR	3.7
1	D	122	ALA	3.5
1	B	18	SER	3.1
1	A	70	THR	2.8
1	D	18	SER	2.7
1	D	19	ALA	2.5
1	D	21	SER	2.5
1	A	73	TYR	2.3
1	B	71	ASN	2.2
1	D	64	THR	2.2
1	C	115	SER	2.0
1	D	119	HIS	2.0
1	A	124	VAL	2.0

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
1	YCM	C	32	10/11	0.88	0.11	42,46,58,58	0
1	YCM	A	32	10/11	0.89	0.11	30,35,43,44	0
1	YCM	B	32	10/11	0.90	0.09	40,40,46,46	0
1	YCM	C	31	10/11	0.91	0.11	36,39,42,43	0
1	YCM	B	31	10/11	0.91	0.11	37,41,48,50	0
1	YCM	D	32	10/11	0.91	0.10	34,35,40,41	0
1	YCM	D	31	10/11	0.92	0.09	35,37,43,45	0
1	YCM	A	31	10/11	0.96	0.06	26,31,34,35	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.