



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:26 PM UTC

PDB ID : 5V3B / pdb_00005v3b
Title : Human A20 OTU domain (WT) with acetamidylated C103
Authors : Langley, D.B.; Christ, D.; Grey, S.
Deposited on : 2017-03-07
Resolution : 3.00 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

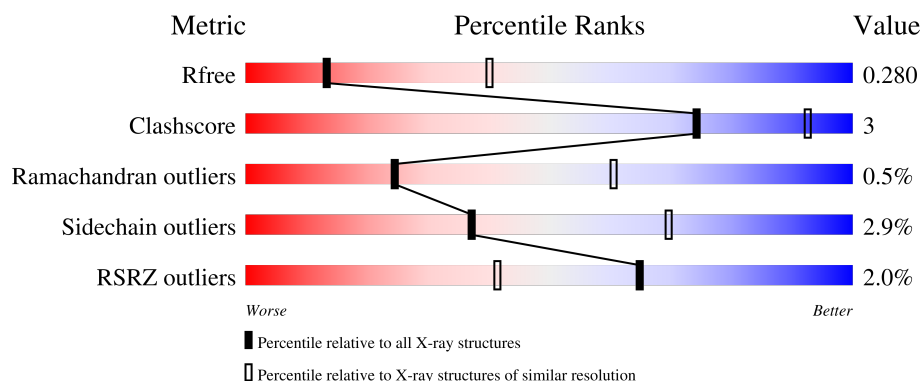
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	366	0% 77% 8% • 15%
1	B	366	83% 5% • 12%
1	C	366	0% 76% 9% • 14%
1	D	366	3% 80% 5% • 14%
1	E	366	2% 78% 8% • 13%

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Mol	Chain	Length	Quality of chain
1	F	366	4% 79% 8% 12%

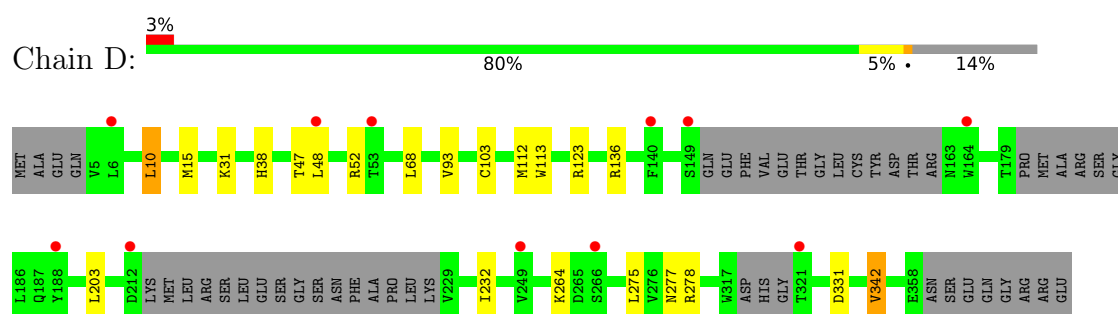
2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 14207 atoms, of which 0 are hydrogens and 0 are deuteriums.

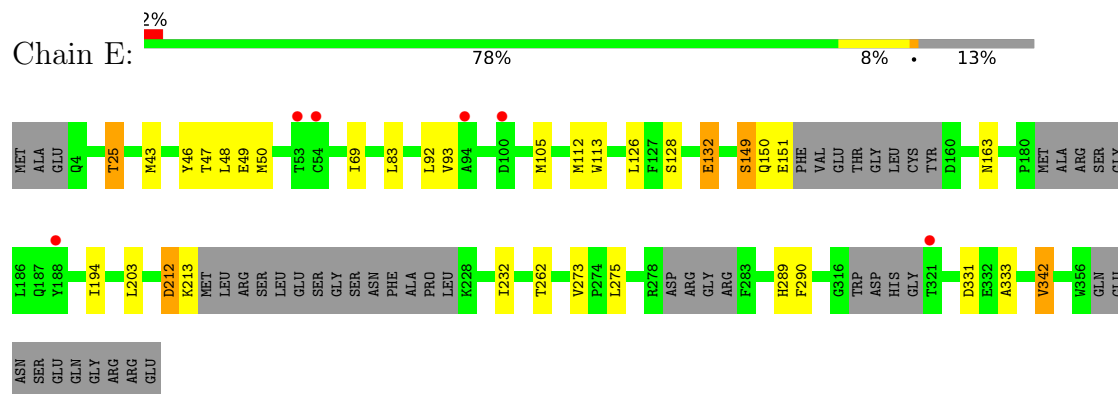
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Tumor necrosis factor alpha-induced protein 3.

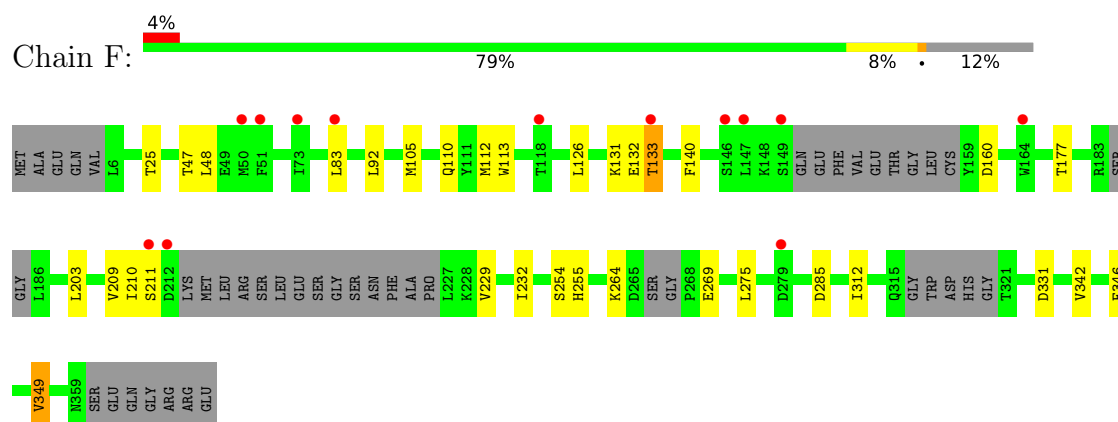
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	312	Total	C	N	O	S	0	0	0
			2405	1556	416	420	13			
1	B	323	Total	C	N	O	S	0	0	0
			2411	1559	420	420	12			
1	C	314	Total	C	N	O	S	0	0	0
			2301	1495	393	400	13			
1	D	316	Total	C	N	O	S	0	0	0
			2358	1527	415	404	12			
1	E	318	Total	C	N	O	S	0	0	0
			2361	1536	400	413	12			
1	F	322	Total	C	N	O	S	0	0	0
			2371	1538	412	407	14			



- Molecule 1: Tumor necrosis factor alpha-induced protein 3



- Molecule 1: Tumor necrosis factor alpha-induced protein 3



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	98.72Å 80.90Å 153.43Å 90.00° 102.46° 90.00°	Depositor
Resolution (Å)	34.56 – 3.00 34.56 – 3.00	Depositor EDS
% Data completeness (in resolution range)	97.3 (34.56-3.00) 97.3 (34.56-3.00)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 3.00Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.238 , 0.282 0.237 , 0.280	Depositor DCC
R_{free} test set	2341 reflections (3.96%)	wwPDB-VP
Wilson B-factor (Å ²)	81.9	Xtriage
Anisotropy	0.531	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 41.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	14207	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.15% 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](#)

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	0.85	1/2452 (0.0%)	0.96	3/3346 (0.1%)
1	B	0.81	0/2463	0.95	4/3376 (0.1%)
1	C	0.80	0/2348	0.95	2/3225 (0.1%)
1	D	0.82	0/2407	0.94	2/3291 (0.1%)
1	E	0.83	0/2409	0.97	3/3303 (0.1%)
1	F	0.86	2/2415 (0.1%)	1.05	8/3307 (0.2%)
All	All	0.83	3/14494 (0.0%)	0.97	22/19848 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	E	0	1
All	All	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	211	SER	C-O	-8.09	1.16	1.24
1	F	254	SER	N-CA	6.49	1.54	1.46
1	A	31	LYS	CA-C	5.01	1.59	1.52

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	210	ILE	N-CA-C	14.23	123.69	110.74
1	E	163	ASN	N-CA-C	9.58	122.72	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	163	ASN	N-CA-C	9.49	122.63	111.71
1	C	176	SER	N-CA-C	8.59	129.10	110.80
1	F	254	SER	N-CA-C	7.81	127.44	110.80
1	D	123	ARG	CB-CG-CD	7.47	128.49	111.30
1	E	163	ASN	CA-C-O	-7.00	112.19	120.10
1	A	264	LYS	N-CA-C	6.62	119.32	109.59
1	F	210	ILE	CB-CA-C	6.26	120.20	111.94
1	A	28	ASP	OD1-CG-OD2	6.20	137.78	122.90
1	B	163	ASN	CA-C-O	-6.19	113.11	120.10
1	F	264	LYS	N-CA-C	6.18	119.92	112.38
1	B	30	PHE	CB-CA-C	-5.90	101.24	110.74
1	E	50	MET	CG-SD-CE	-5.78	88.17	100.90
1	F	211	SER	CA-C-O	-5.73	113.03	120.71
1	C	163	ASN	N-CA-C	5.73	123.00	110.80
1	D	264	LYS	N-CA-C	5.61	118.31	109.96
1	F	254	SER	CA-C-N	5.43	130.27	122.40
1	F	254	SER	C-N-CA	5.43	130.27	122.40
1	A	28	ASP	CB-CG-OD2	-5.38	106.03	118.40
1	B	31	LYS	N-CA-CB	-5.32	100.90	110.37
1	F	255	HIS	N-CA-CB	5.07	120.00	112.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	5	VAL	Peptide
1	E	149	SER	Mainchain

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	2405	0	2233	10	0
1	B	2411	0	2144	7	0
1	C	2301	0	2007	25	0
1	D	2358	0	2102	10	0
1	E	2361	0	2101	20	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	2371	0	2117	13	0
All	All	14207	0	12704	82	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 3.

All (82) 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:C:169:ASP:O	1:C:172:ILE:HG22	1.81	0.81
1:C:241:GLN:HB2	1:E:46:TYR:OH	1.82	0.80
1:E:128:SER:O	1:E:132:GLU:HG2	1.82	0.79
1:C:346:PHE:HA	1:C:349:VAL:HG22	1.71	0.71
1:D:31:LYS:HG2	1:D:38:HIS:CE1	2.26	0.70
1:C:35:GLY:O	1:C:133:THR:HG23	1.91	0.69
1:C:133:THR:HG22	1:C:134:ASP:N	2.11	0.65
1:C:241:GLN:CB	1:E:46:TYR:OH	2.45	0.65
1:C:69:ILE:HD12	1:C:93:VAL:CG1	2.29	0.63
1:E:69:ILE:HD12	1:E:93:VAL:CG1	2.29	0.63
1:A:112:MET:HE1	1:A:203:LEU:HD22	1.80	0.62
1:A:16:ARG:O	1:A:19:VAL:HG22	2.01	0.60
1:E:289:HIS:HB3	1:E:290:PHE:CD1	2.37	0.59
1:C:69:ILE:HD12	1:C:93:VAL:HG12	1.84	0.59
1:D:112:MET:HE1	1:D:203:LEU:HD22	1.84	0.58
1:E:69:ILE:HD12	1:E:93:VAL:HG12	1.87	0.57
1:C:112:MET:HE1	1:C:203:LEU:HD22	1.87	0.57
1:E:112:MET:HE1	1:E:203:LEU:HD22	1.86	0.56
1:B:112:MET:HE1	1:B:203:LEU:HD22	1.85	0.56
1:F:131:LYS:O	1:F:133:THR:N	2.38	0.56
1:F:112:MET:HE1	1:F:203:LEU:HD22	1.88	0.54
1:C:133:THR:HG22	1:C:134:ASP:H	1.73	0.53
1:E:149:SER:O	1:E:151:GLU:N	2.42	0.53
1:A:16:ARG:O	1:A:19:VAL:CG2	2.57	0.53
1:E:48:LEU:HD11	1:E:342:VAL:HG23	1.92	0.51
1:A:83:LEU:HD11	1:A:92:LEU:HB2	1.93	0.50
1:F:48:LEU:HD11	1:F:342:VAL:HG23	1.93	0.50
1:C:83:LEU:HD11	1:C:92:LEU:HB2	1.94	0.50
1:B:275:LEU:C	1:B:286:LEU:HD23	2.37	0.50
1:E:49:GLU:HG3	1:E:262:THR:HB	1.94	0.49
1:B:48:LEU:HD11	1:B:342:VAL:HG23	1.94	0.49
1:D:48:LEU:HD11	1:D:342:VAL:HG23	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:83:LEU:HD11	1:E:92:LEU:HB2	1.94	0.49
1:D:31:LYS:CG	1:D:38:HIS:CE1	2.96	0.48
1:C:50:MET:HE1	1:C:51:PHE:CZ	2.49	0.48
1:C:346:PHE:O	1:C:349:VAL:HG22	2.13	0.48
1:C:133:THR:CG2	1:C:134:ASP:N	2.76	0.47
1:F:346:PHE:O	1:F:349:VAL:HG12	2.15	0.47
1:B:50:MET:HE1	1:B:51:PHE:CZ	2.50	0.47
1:F:83:LEU:HD11	1:F:92:LEU:HB2	1.96	0.47
1:A:48:LEU:HD11	1:A:342:VAL:HG23	1.98	0.46
1:C:176:SER:O	1:C:177:THR:OG1	2.24	0.46
1:A:50:MET:HE1	1:A:51:PHE:CZ	2.51	0.46
1:E:212:ASP:CG	1:E:213:LYS:N	2.74	0.46
1:A:209:VAL:HG11	1:A:229:VAL:HG12	1.99	0.45
1:F:140:PHE:CD1	1:F:140:PHE:C	2.94	0.45
1:B:232:ILE:HD11	1:B:275:LEU:HD22	1.99	0.45
1:A:47:THR:HG21	1:A:331:ASP:HB3	1.99	0.45
1:D:10:LEU:HD13	1:D:68:LEU:HD22	1.99	0.44
1:F:209:VAL:HG11	1:F:229:VAL:HG12	1.98	0.44
1:B:112:MET:HE3	1:B:113:TRP:CZ2	2.53	0.44
1:F:232:ILE:HD11	1:F:275:LEU:HD22	1.99	0.44
1:E:112:MET:HE3	1:E:113:TRP:CZ2	2.53	0.44
1:D:232:ILE:HD11	1:D:275:LEU:HD22	1.99	0.43
1:E:232:ILE:HD11	1:E:275:LEU:HD22	2.01	0.43
1:A:112:MET:HE3	1:A:113:TRP:CZ2	2.53	0.43
1:C:112:MET:HE3	1:C:113:TRP:CZ2	2.53	0.43
1:C:123:ARG:HG3	1:C:177:THR:CG2	2.49	0.43
1:F:112:MET:HE3	1:F:113:TRP:CZ2	2.54	0.43
1:C:232:ILE:HD11	1:C:275:LEU:HD22	2.00	0.43
1:D:112:MET:HE3	1:D:113:TRP:CZ2	2.53	0.43
1:E:49:GLU:HB2	1:E:333:ALA:HB2	1.99	0.43
1:F:177:THR:O	1:F:177:THR:CG2	2.66	0.43
1:B:105:MET:CE	1:B:126:LEU:HD22	2.49	0.42
1:C:209:VAL:HG11	1:C:229:VAL:HG12	2.00	0.42
1:C:133:THR:CG2	1:C:134:ASP:H	2.31	0.42
1:E:47:THR:HG21	1:E:331:ASP:HB3	2.02	0.42
1:F:112:MET:HE1	1:F:203:LEU:CD2	2.50	0.42
1:A:232:ILE:HD11	1:A:275:LEU:HD22	2.01	0.42
1:E:112:MET:HE1	1:E:203:LEU:CD2	2.49	0.42
1:C:231:GLY:HA2	1:C:286:LEU:HD22	2.01	0.41
1:C:121:VAL:HG13	1:C:122:LEU:N	2.35	0.41
1:C:348:LEU:HD11	1:D:15:MET:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:105:MET:CE	1:E:126:LEU:HD22	2.51	0.41
1:D:47:THR:HG21	1:D:331:ASP:HB3	2.02	0.41
1:D:112:MET:HE1	1:D:203:LEU:CD2	2.49	0.41
1:F:47:THR:HG21	1:F:331:ASP:HB3	2.03	0.41
1:E:194:ILE:HD13	1:E:290:PHE:CE2	2.56	0.41
1:C:112:MET:HE1	1:C:203:LEU:CD2	2.51	0.40
1:C:105:MET:CE	1:C:126:LEU:HD22	2.51	0.40
1:E:25:THR:HG23	1:E:43:MET:HE1	2.04	0.40
1:F:105:MET:CE	1:F:126:LEU:HD22	2.52	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	299/366 (82%)	287 (96%)	11 (4%)	1 (0%)	36	70
1	B	314/366 (86%)	297 (95%)	14 (4%)	3 (1%)	12	45
1	C	303/366 (83%)	289 (95%)	13 (4%)	1 (0%)	36	70
1	D	305/366 (83%)	292 (96%)	13 (4%)	0	100	100
1	E	305/366 (83%)	291 (95%)	13 (4%)	1 (0%)	36	70
1	F	309/366 (84%)	295 (96%)	11 (4%)	3 (1%)	12	45
All	All	1835/2196 (84%)	1751 (95%)	75 (4%)	9 (0%)	24	60

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	266	SER
1	E	150	GLN
1	F	132	GLU

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Mol	Chain	Res	Type
1	F	133	THR
1	B	265	ASP
1	B	159	TYR
1	B	264	LYS
1	C	163	ASN
1	F	160	ASP

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	235/334 (70%)	225 (96%)	10 (4%)	26	60
1	B	219/334 (66%)	215 (98%)	4 (2%)	51	77
1	C	201/334 (60%)	196 (98%)	5 (2%)	42	72
1	D	210/334 (63%)	203 (97%)	7 (3%)	33	67
1	E	215/334 (64%)	210 (98%)	5 (2%)	44	74
1	F	209/334 (63%)	203 (97%)	6 (3%)	37	70
All	All	1289/2004 (64%)	1252 (97%)	37 (3%)	37	70

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	THR
1	A	27	GLU
1	A	50	MET
1	A	54	CYS
1	A	58	PRO
1	A	143	GLN
1	A	212	ASP
1	A	266	SER
1	A	269	GLU
1	A	312	ILE
1	B	25	THR
1	B	50	MET

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Mol	Chain	Res	Type
1	B	163	ASN
1	B	329	LYS
1	C	50	MET
1	C	54	CYS
1	C	145	GLU
1	C	269	GLU
1	C	273	VAL
1	D	10	LEU
1	D	52	ARG
1	D	93	VAL
1	D	136	ARG
1	D	277	ASN
1	D	278	ARG
1	D	342	VAL
1	E	25	THR
1	E	132	GLU
1	E	212	ASP
1	E	273	VAL
1	E	342	VAL
1	F	25	THR
1	F	110	GLN
1	F	269	GLU
1	F	285	ASP
1	F	312	ILE
1	F	349	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	110	GLN
1	A	163	ASN
1	A	237	HIS
1	B	110	GLN
1	B	163	ASN
1	B	237	HIS
1	B	255	HIS
1	B	296	ASN
1	C	110	GLN
1	C	237	HIS
1	D	110	GLN
1	D	237	HIS
1	D	296	ASN

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Mol	Chain	Res	Type
1	E	110	GLN
1	E	237	HIS
1	E	255	HIS
1	F	110	GLN
1	F	237	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 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	C	103	1	7,9,10	0.86	0	5,10,12	0.55	0
1	YCM	E	103	1	7,9,10	0.97	0	5,10,12	0.56	0
1	YCM	A	103	1	7,9,10	1.01	1 (14%)	5,10,12	0.51	0
1	YCM	B	103	1	7,9,10	0.94	0	5,10,12	0.55	0
1	YCM	D	103	1	7,9,10	1.05	1 (14%)	5,10,12	0.48	0
1	YCM	F	103	1	7,9,10	0.84	0	5,10,12	0.49	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	C	103	1	-	1/6/8/10	-
1	YCM	E	103	1	-	1/6/8/10	-
1	YCM	A	103	1	-	1/6/8/10	-
1	YCM	B	103	1	-	1/6/8/10	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	YCM	D	103	1	-	1/6/8/10	-
1	YCM	F	103	1	-	1/6/8/10	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	103	YCM	CD-SG	-2.16	1.76	1.81
1	A	103	YCM	CD-SG	-2.06	1.76	1.81

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	E	103	YCM	CA-CB-SG-CD
1	F	103	YCM	CA-CB-SG-CD
1	A	103	YCM	CA-CB-SG-CD
1	B	103	YCM	CA-CB-SG-CD
1	C	103	YCM	CA-CB-SG-CD
1	D	103	YCM	CA-CB-SG-CD

There are no ring outliers.

No monomer is involved in short contacts.

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 ⓘ

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	311/366 (84%)	0.23	4 (1%)	75 53	48, 74, 109, 133	0
1	B	322/366 (87%)	0.40	1 (0%)	90 79	48, 86, 137, 182	0
1	C	313/366 (85%)	0.43	3 (0%)	79 59	49, 84, 127, 156	0
1	D	315/366 (86%)	0.55	11 (3%)	47 27	54, 90, 130, 176	0
1	E	317/366 (86%)	0.40	6 (1%)	66 43	49, 83, 122, 153	0
1	F	321/366 (87%)	0.50	13 (4%)	42 23	54, 89, 137, 188	0
All	All	1899/2196 (86%)	0.42	38 (2%)	65 41	48, 84, 130, 188	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	100	ASP	3.4
1	F	212	ASP	3.1
1	F	50	MET	3.1
1	D	321	THR	3.1
1	D	140	PHE	3.1
1	D	249	VAL	3.0
1	F	51	PHE	2.9
1	D	212	ASP	2.9
1	B	132	GLU	2.8
1	F	211	SER	2.8
1	D	149	SER	2.7
1	E	94	ALA	2.7
1	D	266	SER	2.6
1	F	118	THR	2.6
1	C	266	SER	2.6
1	C	336	PRO	2.5
1	A	277	ASN	2.5
1	D	53	THR	2.4
1	F	133	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	188	TYR	2.3
1	D	6	LEU	2.3
1	D	48	LEU	2.3
1	F	83	LEU	2.3
1	F	146	SER	2.3
1	F	164	TRP	2.3
1	E	54	CYS	2.3
1	A	312	ILE	2.2
1	F	147	LEU	2.2
1	A	313	PRO	2.2
1	D	164	TRP	2.1
1	F	73	ILE	2.1
1	F	149	SER	2.1
1	F	279	ASP	2.1
1	A	265	ASP	2.0
1	E	53	THR	2.0
1	C	53	THR	2.0
1	D	188	TYR	2.0
1	E	321	THR	2.0

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

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	B	103	10/11	0.80	0.20	68,86,91,106	0
1	YCM	D	103	10/11	0.82	0.15	76,88,125,128	0
1	YCM	C	103	10/11	0.86	0.10	65,76,93,98	0
1	YCM	E	103	10/11	0.88	0.10	58,70,108,116	0
1	YCM	A	103	10/11	0.91	0.11	66,79,152,160	0
1	YCM	F	103	10/11	0.91	0.09	57,75,114,125	0

6.3 Carbohydrates ⓘ

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.