



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 04:32 pm BST

PDB ID : 5V3B
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

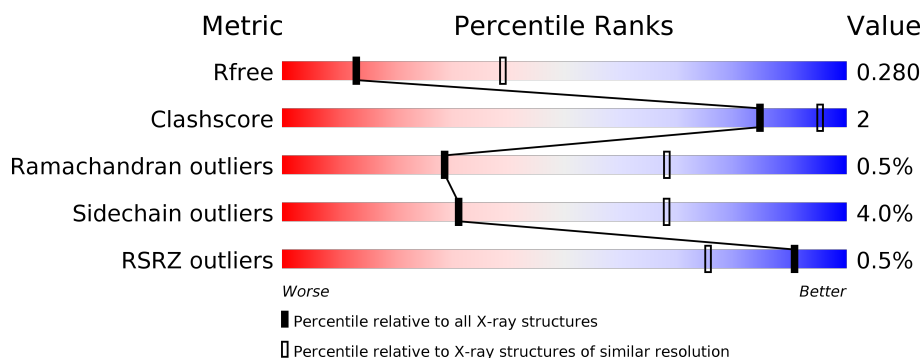
1 Overall quality at a glance

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}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (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 on 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	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>7%</div> <div>15%</div> </div> </div>
1	B	366	<div> <div>82%</div> <div>6%</div> <div>12%</div> </div>
1	C	366	<div> <div>%</div> <div> <div></div> <div>76%</div> <div>9%</div> <div>14%</div> </div> </div>
1	D	366	<div> <div>80%</div> <div>6%</div> <div>14%</div> </div>
1	E	366	<div> <div>78%</div> <div>8%</div> <div>13%</div> </div>
1	F	366	<div> <div>%</div> <div> <div></div> <div>80%</div> <div>7%</div> <div>12%</div> </div> </div>

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.

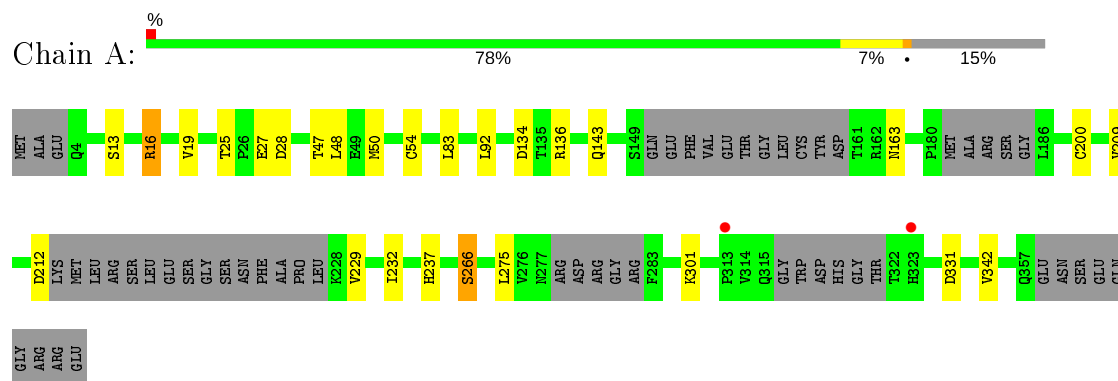
- Molecule 1 is a protein called Tumor necrosis factor alpha-induced protein 3.

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			

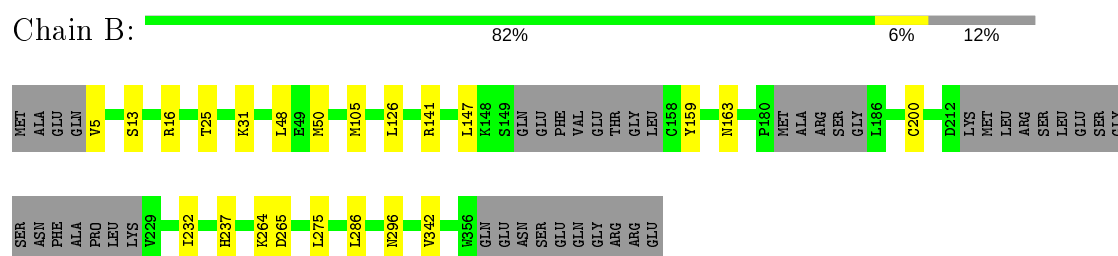
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

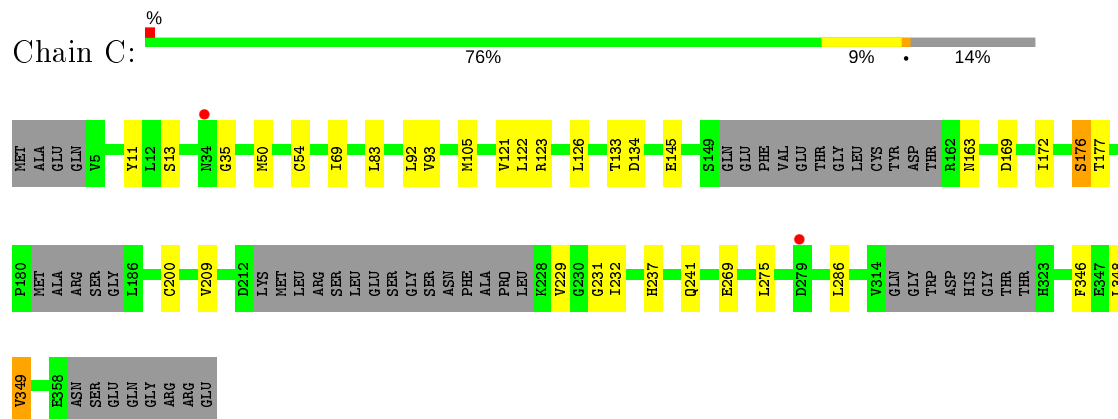
- Molecule 1: Tumor necrosis factor alpha-induced protein 3




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

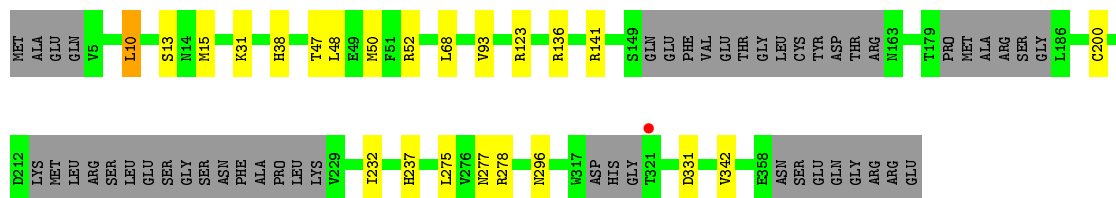


- Molecule 1: Tumor necrosis factor alpha-induced protein 3



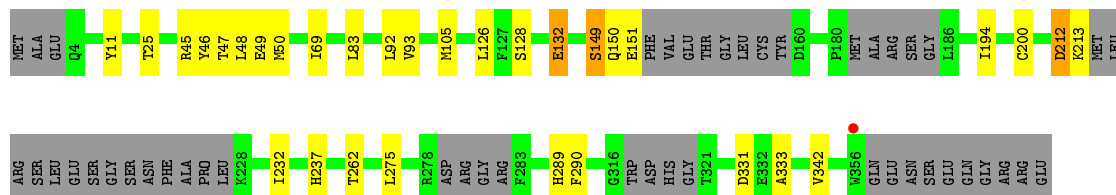
- Molecule 1: Tumor necrosis factor alpha-induced protein 3

Chain D:  80% 6% 14%




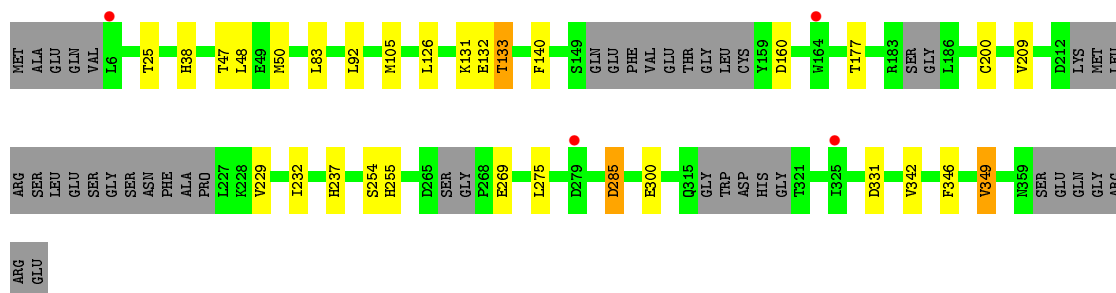
- Molecule 1: Tumor necrosis factor alpha-induced protein 3

Chain E:  78% 8% 13%



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

Chain F:  80% 7% 12%



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) 97.4 (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	2340 reflections (5.04%)	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)	0.23 , 42.9	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

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	0.60	0/2452	0.83	7/3346 (0.2%)
1	B	0.56	0/2463	0.76	4/3376 (0.1%)
1	C	0.55	0/2348	0.73	2/3225 (0.1%)
1	D	0.57	0/2407	0.75	4/3291 (0.1%)
1	E	0.58	0/2409	0.75	2/3303 (0.1%)
1	F	0.59	0/2415	0.81	4/3307 (0.1%)
All	All	0.58	0/14494	0.77	23/19848 (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

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	28	ASP	CB-CG-OD2	-13.63	106.03	118.30
1	A	28	ASP	OD1-CG-OD2	7.62	137.78	123.30
1	E	50	MET	CG-SD-CE	-7.52	88.17	100.20
1	F	285	ASP	CB-CG-OD1	-6.79	112.19	118.30
1	C	176	SER	N-CA-C	6.70	129.10	111.00
1	E	45	ARG	NE-CZ-NH2	-6.62	116.99	120.30
1	D	123	ARG	CB-CG-CD	6.50	128.49	111.60
1	F	254	SER	N-CA-C	6.09	127.44	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	28	ASP	CB-CG-OD1	-6.07	112.83	118.30
1	D	50	MET	CG-SD-CE	-5.93	90.71	100.20
1	A	134	ASP	CB-CG-OD1	5.64	123.37	118.30
1	B	147	LEU	CA-CB-CG	5.48	127.91	115.30
1	B	141	ARG	NE-CZ-NH2	-5.45	117.57	120.30
1	F	50	MET	CG-SD-CE	-5.42	91.53	100.20
1	B	31	LYS	N-CA-CB	-5.39	100.90	110.60
1	D	123	ARG	NE-CZ-NH1	5.37	122.98	120.30
1	A	136	ARG	NE-CZ-NH2	5.33	122.96	120.30
1	D	141	ARG	NE-CZ-NH2	5.25	122.93	120.30
1	F	255	HIS	N-CA-CB	5.22	120.00	110.60
1	C	349	VAL	CA-CB-CG2	5.15	118.62	110.90
1	B	16	ARG	NE-CZ-NH2	-5.15	117.73	120.30
1	A	134	ASP	CB-CG-OD2	-5.08	113.73	118.30
1	A	16	ARG	NE-CZ-NH1	5.02	122.81	120.30

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	7	0
1	B	2411	0	2144	4	0
1	C	2301	0	2007	22	0
1	D	2358	0	2102	7	0
1	E	2361	0	2101	17	0
1	F	2371	0	2117	10	0
All	All	14207	0	12704	63	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 2.

All (63) 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: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:E:69:ILE:HD12	1:E:93:VAL:HG12	1.87	0.57
1:F:131:LYS:O	1:F:133:THR:N	2.38	0.56
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: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
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: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: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:209:VAL:HG11	1:A:229:VAL:HG12	1.99	0.45
1:B:275:LEU:C	1:B:286:LEU:HD23	2.37	0.45
1:C:241:GLN:HB2	1:E:46:TYR:HH	1.80	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:232:ILE:HD11	1:F:275:LEU:HD22	1.99	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:C:123:ARG:HG3	1:C:177:THR:CG2	2.49	0.43
1:C:232:ILE:HD11	1:C:275:LEU:HD22	2.00	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:A:232:ILE:HD11	1:A:275:LEU:HD22	2.01	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
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:E:212:ASP:CG	1:E:213:LYS:N	2.74	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:105:MET:CE	1:C:126:LEU:HD22	2.51	0.40
1:F:140:PHE:C	1:F:140:PHE:CD1	2.94	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%)	41 76
1	B	314/366 (86%)	297 (95%)	14 (4%)	3 (1%)	15 53

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	303/366 (83%)	289 (95%)	13 (4%)	1 (0%)	41	76
1	D	305/366 (83%)	292 (96%)	13 (4%)	0	100	100
1	E	305/366 (83%)	291 (95%)	13 (4%)	1 (0%)	41	76
1	F	309/366 (84%)	295 (96%)	11 (4%)	3 (1%)	15	53
All	All	1835/2196 (84%)	1751 (95%)	75 (4%)	9 (0%)	29	68

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	266	SER
1	E	150	GLN
1	F	132	GLU
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%)	223 (95%)	12 (5%)	24	60
1	B	219/334 (66%)	212 (97%)	7 (3%)	39	74
1	C	201/334 (60%)	193 (96%)	8 (4%)	31	68
1	D	210/334 (63%)	200 (95%)	10 (5%)	25	62
1	E	215/334 (64%)	209 (97%)	6 (3%)	43	77
1	F	209/334 (63%)	201 (96%)	8 (4%)	33	69
All	All	1289/2004 (64%)	1238 (96%)	51 (4%)	31	68

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	SER
1	A	25	THR
1	A	27	GLU
1	A	50	MET
1	A	54	CYS
1	A	143	GLN
1	A	163	ASN
1	A	200	CYS
1	A	212	ASP
1	A	237	HIS
1	A	266	SER
1	A	301	LYS
1	B	13	SER
1	B	25	THR
1	B	50	MET
1	B	163	ASN
1	B	200	CYS
1	B	237	HIS
1	B	296	ASN
1	C	11	TYR
1	C	13	SER
1	C	50	MET
1	C	54	CYS
1	C	145	GLU
1	C	200	CYS
1	C	237	HIS
1	C	269	GLU
1	D	10	LEU
1	D	13	SER
1	D	52	ARG
1	D	93	VAL
1	D	136	ARG
1	D	200	CYS
1	D	237	HIS
1	D	277	ASN
1	D	278	ARG
1	D	296	ASN
1	E	11	TYR
1	E	25	THR
1	E	132	GLU
1	E	200	CYS
1	E	212	ASP
1	E	237	HIS

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Mol	Chain	Res	Type
1	F	25	THR
1	F	38	HIS
1	F	200	CYS
1	F	237	HIS
1	F	269	GLU
1	F	285	ASP
1	F	300	GLU
1	F	349	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	163	ASN

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	A	103	1	7,9,10	0.98	1 (14%)	4,10,12	0.47	0
1	YCM	C	103	1	7,9,10	0.83	0	4,10,12	0.51	0
1	YCM	E	103	1	7,9,10	0.94	0	4,10,12	0.55	0
1	YCM	B	103	1	7,9,10	0.91	0	4,10,12	0.52	0
1	YCM	D	103	1	7,9,10	1.02	1 (14%)	4,10,12	0.45	0
1	YCM	F	103	1	7,9,10	0.82	0	4,10,12	0.51	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	A	103	1	-	0/6/8/10	-
1	YCM	C	103	1	-	0/6/8/10	-
1	YCM	E	103	1	-	1/6/8/10	-
1	YCM	B	103	1	-	0/6/8/10	-
1	YCM	D	103	1	-	0/6/8/10	-
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	-2.10	1.76	1.81
1	A	103	YCM	CD-SG	-2.00	1.76	1.81

There are no bond angle outliers.

There are no chirality outliers.

All (2) 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

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates 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 [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	311/366 (84%)	-0.46	2 (0%) 89 72	48, 74, 109, 133	0
1	B	322/366 (87%)	-0.34	0 100 100	48, 86, 137, 182	0
1	C	313/366 (85%)	-0.37	2 (0%) 89 72	49, 84, 127, 156	0
1	D	315/366 (86%)	-0.38	1 (0%) 94 84	53, 90, 130, 176	0
1	E	317/366 (86%)	-0.49	1 (0%) 94 84	49, 83, 122, 153	0
1	F	321/366 (87%)	-0.20	4 (1%) 79 54	54, 89, 137, 188	0
All	All	1899/2196 (86%)	-0.37	10 (0%) 91 75	48, 84, 130, 188	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	279	ASP	3.9
1	C	279	ASP	2.6
1	F	164	TRP	2.6
1	A	313	PRO	2.6
1	F	6	LEU	2.5
1	A	323	HIS	2.4
1	C	34	ASN	2.3
1	E	356	TRP	2.3
1	D	321	THR	2.3
1	F	325	ILE	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(\AA^2)	Q<0.9
1	YCM	D	103	10/11	0.80	0.22	76,88,125,128	0
1	YCM	B	103	10/11	0.81	0.28	68,86,91,106	0
1	YCM	C	103	10/11	0.90	0.18	65,76,93,98	0
1	YCM	F	103	10/11	0.91	0.20	57,75,114,125	0
1	YCM	E	103	10/11	0.95	0.11	58,70,108,116	0
1	YCM	A	103	10/11	0.95	0.18	66,79,152,160	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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.