



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 10, 2020 – 07:31 AM BST

PDB ID : 6W9V
Title : Structure of human MAIT A-F7 TCR in complex with patient MR1-R9H without ligand
Authors : Awad, W.; Rossjohn, J.
Deposited on : 2020-03-23
Resolution : 1.95 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 : 2.13.1
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) : 2.13.1

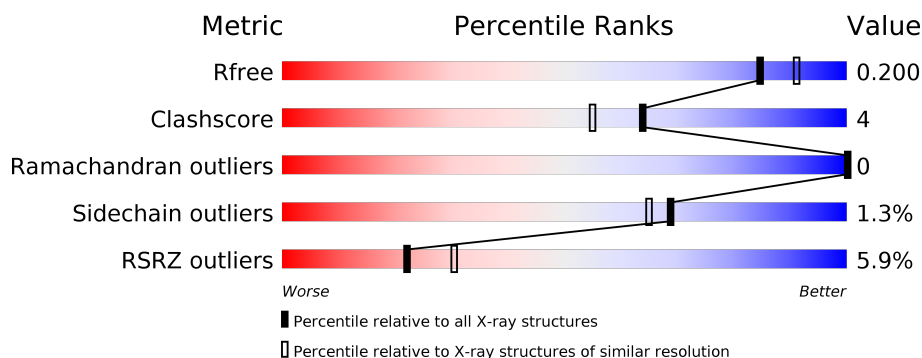
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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.



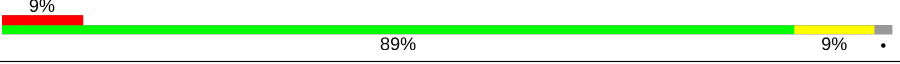

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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	271	<div> <div>3%</div> <div> <div></div> <div>87%</div> <div>11%</div> <div></div> </div> <div></div> </div>
1	C	271	<div> <div>3%</div> <div> <div></div> <div>87%</div> <div>10%</div> <div></div> </div> <div></div> </div>
2	B	100	<div> <div>21%</div> <div> <div></div> <div>91%</div> <div>6%</div> <div></div> </div> <div></div> </div>
2	F	100	<div> <div></div> <div> <div></div> <div>95%</div> <div>5%</div> <div></div> </div> <div></div> </div>
3	D	204	<div> <div>17%</div> <div> <div></div> <div>84%</div> <div>8%</div> <div>8%</div> </div> <div></div> </div>
3	G	204	<div> <div></div> <div> <div></div> <div>88%</div> <div>10%</div> <div></div> </div> <div></div> </div>

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Mol	Chain	Length	Quality of chain
4	E	246	
4	H	246	

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
6	GOL	C	301	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 14539 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 Major histocompatibility complex class I-related gene protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	20	0
			2276	1467	390	406	13			
1	C	267	Total	C	N	O	S	0	16	0
			2268	1466	385	405	12			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	expression tag	UNP Q95460
A	9	HIS	ARG	engineered mutation	UNP Q95460
A	261	SER	CYS	conflict	UNP Q95460
C	0	MET	-	expression tag	UNP Q95460
C	9	HIS	ARG	engineered mutation	UNP Q95460
C	261	SER	CYS	conflict	UNP Q95460

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	98	Total	C	N	O	S	0	2	0
			783	504	131	145	3			
2	F	100	Total	C	N	O	S	0	3	0
			822	529	139	150	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	expression tag	UNP P61769
F	0	MET	-	expression tag	UNP P61769

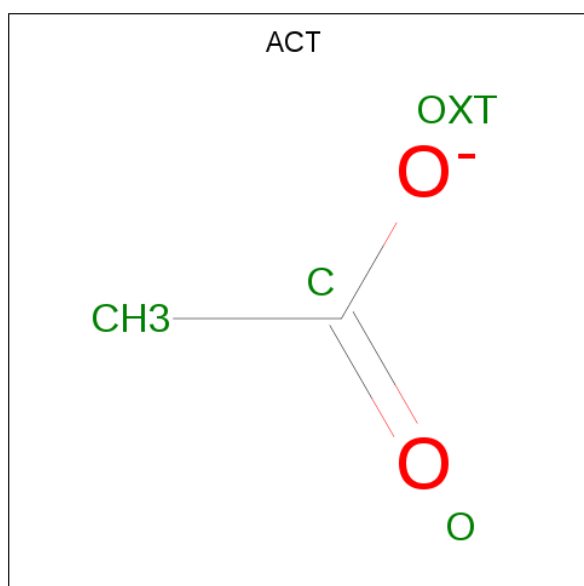
- Molecule 3 is a protein called TCR-alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	188	Total	C	N	O	S	0	9	0
			1460	944	227	280	9			
3	G	201	Total	C	N	O	S	0	20	0
			1644	1048	254	329	13			

- Molecule 4 is a protein called TCR-beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	241	Total	C	N	O	S	0	9	0
			1901	1205	323	361	12			
4	H	244	Total	C	N	O	S	0	16	0
			1967	1249	334	371	13			

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	H	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	C	1	Total	C	O	0	0
			6	3	3		
6	F	1	Total	C	O	0	0
			6	3	3		

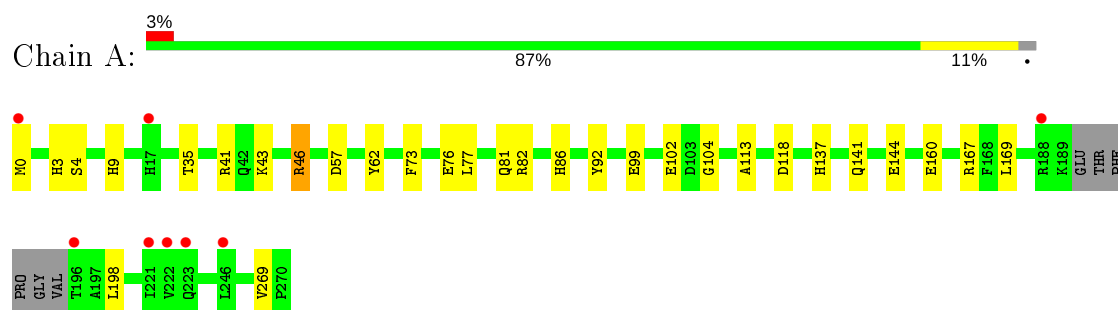
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	247	Total	O	0	0
			247	247		
7	B	67	Total	O	0	0
			67	67		
7	C	229	Total	O	0	0
			229	229		
7	D	121	Total	O	0	0
			121	121		
7	E	138	Total	O	0	0
			138	138		
7	F	105	Total	O	0	0
			105	105		
7	G	202	Total	O	0	0
			202	202		
7	H	277	Total	O	0	0
			277	277		

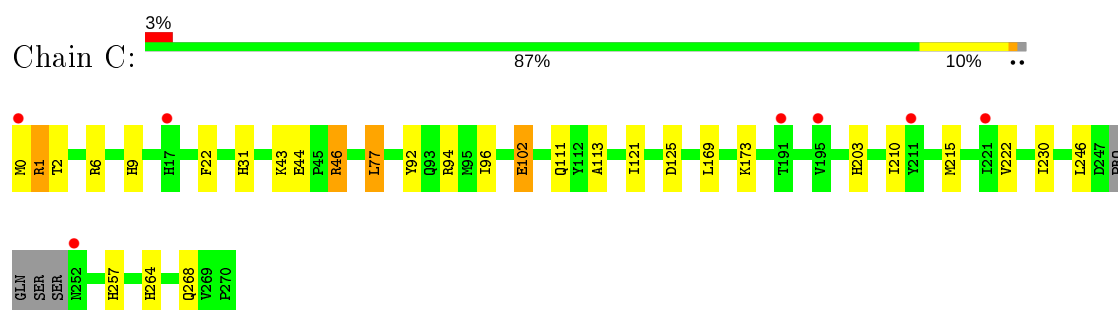
3 Residue-property plots [i](#)

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.

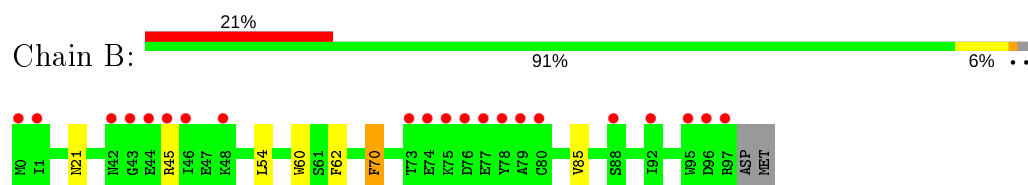
- Molecule 1: Major histocompatibility complex class I-related gene protein



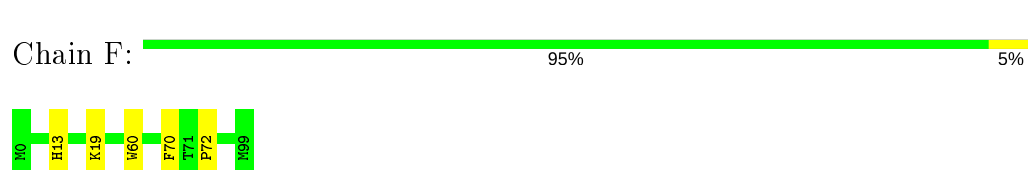
- Molecule 1: Major histocompatibility complex class I-related gene protein



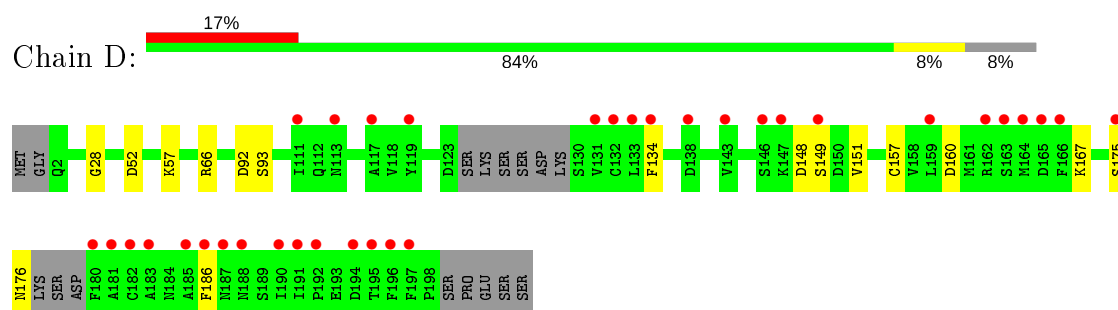
- Molecule 2: Beta-2-microglobulin



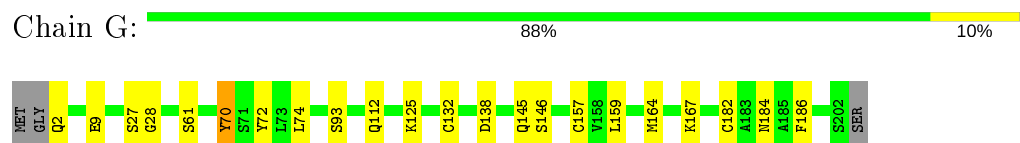
- Molecule 2: Beta-2-microglobulin



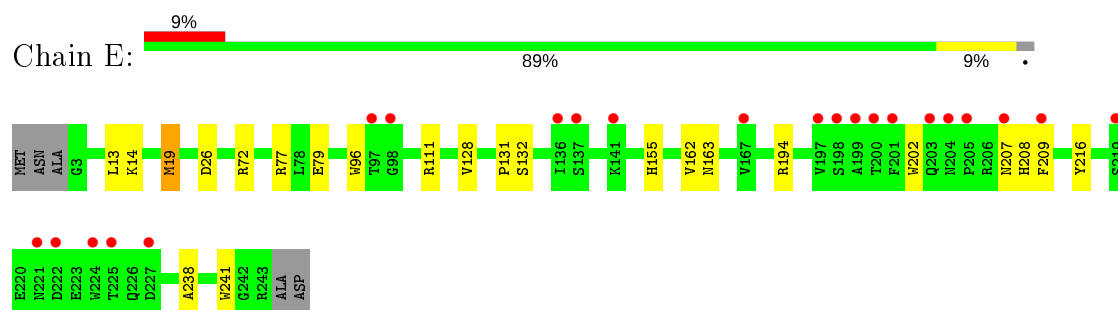
- Molecule 3: TCR-alpha chain



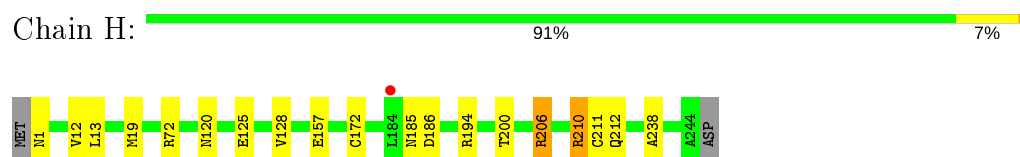
• Molecule 3: TCR-alpha chain



• Molecule 4: TCR-beta chain



• Molecule 4: TCR-beta chain



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	216.99Å 70.41Å 142.90Å 90.00° 104.15° 90.00°	Depositor
Resolution (Å)	46.19 – 1.95 47.91 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.6 (46.19-1.95) 99.6 (47.91-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 1.95Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, R_{free}	0.160 , 0.200 0.162 , 0.200	Depositor DCC
R_{free} test set	7601 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	35.9	Xtriage
Anisotropy	0.325	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 47.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	14539	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% 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: GOL, ACT

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	0.65	0/2400	0.72	3/3261 (0.1%)
1	C	0.60	0/2383	0.73	3/3239 (0.1%)
2	B	0.50	0/812	0.61	0/1107
2	F	0.60	0/854	0.69	0/1161
3	D	0.59	0/1520	0.73	2/2068 (0.1%)
3	G	0.70	0/1743	0.78	1/2365 (0.0%)
4	E	0.54	0/1976	0.66	1/2692 (0.0%)
4	H	0.67	0/2067	0.72	1/2814 (0.0%)
All	All	0.62	0/13755	0.71	11/18707 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	46[A]	ARG	NE-CZ-NH1	-7.00	116.80	120.30
1	A	46[B]	ARG	NE-CZ-NH1	-7.00	116.80	120.30
1	C	1	ARG	NE-CZ-NH1	6.93	123.77	120.30
3	G	157	CYS	CA-CB-SG	6.70	126.06	114.00
4	H	210	ARG	NE-CZ-NH1	-6.16	117.22	120.30
1	C	1	ARG	NE-CZ-NH2	-6.13	117.24	120.30
3	D	92	ASP	CB-CG-OD1	6.09	123.78	118.30
4	E	77	ARG	NE-CZ-NH1	-5.88	117.36	120.30
3	D	157	CYS	CA-CB-SG	5.51	123.92	114.00
1	A	118	ASP	CB-CG-OD1	5.32	123.09	118.30
1	C	77	LEU	CA-CB-CG	5.09	127.00	115.30

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	2276	0	2182	22	0
1	C	2268	0	2177	28	0
2	B	783	0	716	5	0
2	F	822	0	782	3	0
3	D	1460	0	1347	8	0
3	G	1644	0	1594	15	0
4	E	1901	0	1778	14	0
4	H	1967	0	1881	17	0
5	A	4	0	3	0	0
5	H	4	0	3	0	0
6	A	12	0	16	0	0
6	C	6	0	8	5	0
6	F	6	0	8	0	0
7	A	247	0	0	10	2
7	B	67	0	0	2	0
7	C	229	0	0	14	0
7	D	121	0	0	2	2
7	E	138	0	0	4	0
7	F	105	0	0	1	0
7	G	202	0	0	7	0
7	H	277	0	0	8	0
All	All	14539	0	12495	107	2

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 4.

All (107) 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:44:GLU:OE1	7:C:401:HOH:O	1.76	1.00
4:E:111:ARG:NH2	7:E:301:HOH:O	1.99	0.95
1:C:94:ARG:HH12	6:C:301:GOL:H12	1.35	0.92
1:A:92:TYR:OH	7:A:401:HOH:O	1.88	0.90
1:C:173:LYS:NZ	7:C:402:HOH:O	2.00	0.89
4:H:210:ARG:NH1	4:H:212:GLN:OE1	2.11	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:264[B]:HIS:ND1	7:C:406:HOH:O	2.16	0.79
1:C:215[B]:MET:HG3	1:C:257:HIS:CD2	2.19	0.78
4:E:14:LYS:NZ	7:E:303:HOH:O	2.19	0.74
3:G:184[B]:ASN:ND2	7:G:305:HOH:O	2.22	0.72
2:B:45:ARG:NH1	7:B:101:HOH:O	2.21	0.71
1:C:268[B]:GLN:OE1	7:C:403:HOH:O	2.08	0.70
6:C:301:GOL:O1	7:C:404:HOH:O	2.10	0.68
1:C:246:LEU:O	7:C:405:HOH:O	2.11	0.68
4:H:128[A]:VAL:HG23	4:H:238:ALA:HB3	1.76	0.67
3:G:182[B]:CYS:SG	7:G:410:HOH:O	2.52	0.66
1:C:9[B]:HIS:ND1	7:C:408:HOH:O	2.20	0.66
1:C:0:MET:SD	1:C:169[B]:LEU:HD21	2.36	0.66
1:C:43:LYS:NZ	7:C:412:HOH:O	2.27	0.65
1:C:222:VAL:HA	4:H:200[B]:THR:HG21	1.79	0.65
1:C:94:ARG:NH1	6:C:301:GOL:H12	2.11	0.64
4:H:210:ARG:HD2	7:H:1854:HOH:O	1.97	0.64
1:C:94:ARG:HH12	6:C:301:GOL:C1	2.09	0.64
3:G:138[B]:ASP:OD1	7:G:302:HOH:O	2.15	0.64
3:G:9:GLU:OE1	7:G:303:HOH:O	2.15	0.63
1:C:1:ARG:HD2	7:C:595:HOH:O	1.98	0.62
4:H:12[B]:VAL:HG23	7:H:1708:HOH:O	1.98	0.62
2:B:85:VAL:HG13	7:B:109:HOH:O	2.00	0.61
3:D:28:GLY:HA3	3:D:93[A]:SER:HB3	1.84	0.59
3:G:70[A]:TYR:HH	3:G:72:TYR:HD2	1.50	0.59
1:A:0:MET:HE1	1:A:169[B]:LEU:HG	1.83	0.59
1:A:3:HIS:CD2	1:A:169[B]:LEU:HD21	2.38	0.59
1:A:76[B]:GLU:OE1	7:A:401:HOH:O	2.17	0.58
4:E:79:GLU:OE1	7:E:302:HOH:O	2.17	0.57
1:C:125:ASP:OD2	7:C:407:HOH:O	2.18	0.56
4:E:155:HIS:HB3	4:E:216:TYR:HB2	1.88	0.56
1:A:160:GLU:OE2	7:A:402:HOH:O	2.18	0.55
3:D:148:ASP:HB3	3:D:151:VAL:HG12	1.88	0.55
1:A:113:ALA:HB2	2:B:60:TRP:CE2	2.41	0.55
1:A:9[B]:HIS:ND1	7:A:405:HOH:O	2.20	0.55
3:D:66[B]:ARG:NH2	7:D:301:HOH:O	2.20	0.55
1:C:77:LEU:HD13	1:C:92:TYR:HB2	1.89	0.55
3:G:159[A]:LEU:HB3	4:H:172[A]:CYS:HB2	1.91	0.53
1:C:6[A]:ARG:HH11	1:C:6[A]:ARG:HG3	1.73	0.53
1:A:43:LYS:HD2	1:A:62:TYR:HB3	1.90	0.52
1:C:230:ILE:HD11	4:H:206[B]:ARG:HD2	1.92	0.52
4:H:206[A]:ARG:HG3	7:H:1801:HOH:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:GLU:O	7:A:404:HOH:O	2.20	0.51
3:G:145:GLN:OE1	7:G:304:HOH:O	2.19	0.50
4:H:1:ASN:HA	7:H:1896:HOH:O	2.11	0.50
4:E:131:PRO:HD2	4:E:202:TRP:CZ2	2.46	0.50
3:G:132[B]:CYS:HG	3:G:186:PHE:HE1	1.59	0.49
4:E:208:HIS:HB2	4:E:241:TRP:CZ3	2.47	0.49
1:C:1:ARG:HA	1:C:102[B]:GLU:HG2	1.94	0.48
1:C:113:ALA:HB2	2:F:60:TRP:CE2	2.49	0.47
4:E:163:ASN:OD1	4:E:207:ASN:HA	2.13	0.47
2:F:13:HIS:HE1	7:F:246:HOH:O	1.97	0.47
3:D:175:SER:OG	3:D:176:ASN:N	2.47	0.47
1:C:203:HIS:HE1	7:C:592:HOH:O	1.95	0.47
1:C:210:ILE:O	4:H:206[A]:ARG:HD2	2.15	0.47
1:C:9[B]:HIS:HD2	1:C:22:PHE:HZ	1.63	0.47
4:E:13:LEU:HD21	4:E:19[B]:MET:HG3	1.97	0.47
4:H:125:GLU:OE2	7:H:1701:HOH:O	2.20	0.47
1:A:104:GLY:O	7:A:403:HOH:O	2.19	0.47
1:A:4:SER:HB3	1:A:99:GLU:HG2	1.96	0.47
4:E:128:VAL:HG23	4:E:238:ALA:HB3	1.97	0.46
3:G:28:GLY:HA3	3:G:93[A]:SER:HB3	1.97	0.46
3:D:160:ASP:HA	3:D:167:LYS:HG2	1.98	0.46
1:A:41[A]:ARG:HD2	7:A:447:HOH:O	2.16	0.46
7:C:572:HOH:O	3:G:125:LYS:HD2	2.16	0.46
4:H:72:ARG:NH2	7:H:1711:HOH:O	2.48	0.45
1:C:111:GLN:HG2	1:C:121:ILE:HG23	1.97	0.45
4:E:19[A]:MET:HE3	4:E:19[A]:MET:HB2	1.84	0.45
1:C:2[B]:THR:HG22	1:C:102[B]:GLU:H	1.82	0.45
1:A:137:HIS:O	1:A:141[B]:GLN:HG2	2.17	0.45
1:A:35:THR:HB	1:A:46[A]:ARG:HD3	1.98	0.45
1:A:167:ARG:HD2	7:A:593:HOH:O	2.17	0.44
4:H:120:ASN:HB3	7:H:1808:HOH:O	2.18	0.44
1:C:96:ILE:HG21	6:C:301:GOL:H11	2.00	0.44
3:G:28:GLY:HA3	3:G:93[B]:SER:OG	2.18	0.44
1:A:77:LEU:HD13	1:A:92:TYR:HB2	1.99	0.44
3:D:134:PHE:HB2	3:D:186:PHE:CE2	2.54	0.43
2:B:54:LEU:HD11	2:B:62:PHE:HB3	2.00	0.43
2:B:21:ASN:HB3	2:B:70:PHE:CE1	2.53	0.43
3:G:2:GLN:OE1	3:G:27:SER:N	2.48	0.43
4:H:128[B]:VAL:HG22	4:H:238:ALA:HB1	2.01	0.43
4:E:96:TRP:O	7:E:304:HOH:O	2.22	0.43
3:G:112:GLN:NE2	7:G:301:HOH:O	2.06	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:0:MET:CE	1:A:169[B]:LEU:HG	2.47	0.42
1:A:76[A]:GLU:HG2	7:A:401:HOH:O	2.18	0.42
2:F:19:LYS:O	2:F:72:PRO:HD2	2.19	0.42
1:C:46[B]:ARG:NE	7:C:414:HOH:O	2.35	0.42
3:D:52:ASP:OD1	7:D:302:HOH:O	2.21	0.42
1:C:31:HIS:HE1	7:C:537:HOH:O	2.02	0.42
4:H:13:LEU:HD11	4:H:19[B]:MET:SD	2.60	0.42
1:A:57[B]:ASP:OD1	7:A:406:HOH:O	2.21	0.41
4:E:128:VAL:HG23	4:E:238:ALA:CB	2.50	0.41
4:E:26[A]:ASP:OD1	4:E:72[A]:ARG:NE	2.41	0.41
1:A:81:GLN:HG2	1:A:86:HIS:O	2.21	0.41
4:H:157:GLU:OE2	7:H:1702:HOH:O	2.22	0.41
1:A:0:MET:O	1:A:102:GLU:HG3	2.21	0.41
4:H:128[B]:VAL:HG13	4:H:211[B]:CYS:SG	2.61	0.41
3:D:148:ASP:OD2	3:D:149:SER:N	2.53	0.41
3:G:61[B]:SER:OG	3:G:74:LEU:HB3	2.21	0.41
1:A:198:LEU:HD13	1:A:269:VAL:HG21	2.02	0.41
4:E:162:VAL:HG22	4:E:209:PHE:HD2	1.87	0.40
3:G:167:LYS:NZ	7:G:314:HOH:O	2.54	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:632:HOH:O	7:D:396:HOH:O 2_658	2.03	0.17
7:A:493:HOH:O	7:D:321:HOH:O 2_658	2.17	0.03

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	A	281/271 (104%)	279 (99%)	2 (1%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	279/271 (103%)	274 (98%)	5 (2%)	0	100	100
2	B	98/100 (98%)	98 (100%)	0	0	100	100
2	F	101/100 (101%)	101 (100%)	0	0	100	100
3	D	191/204 (94%)	186 (97%)	5 (3%)	0	100	100
3	G	220/204 (108%)	216 (98%)	4 (2%)	0	100	100
4	E	248/246 (101%)	245 (99%)	3 (1%)	0	100	100
4	H	258/246 (105%)	255 (99%)	3 (1%)	0	100	100
All	All	1676/1642 (102%)	1654 (99%)	22 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	245/241 (102%)	242 (99%)	3 (1%)	71	68
1	C	242/241 (100%)	238 (98%)	4 (2%)	60	55
2	B	82/95 (86%)	81 (99%)	1 (1%)	71	68
2	F	90/95 (95%)	89 (99%)	1 (1%)	73	71
3	D	150/181 (83%)	149 (99%)	1 (1%)	84	82
3	G	194/181 (107%)	188 (97%)	6 (3%)	40	28
4	E	202/212 (95%)	198 (98%)	4 (2%)	55	48
4	H	214/212 (101%)	208 (97%)	6 (3%)	43	33
All	All	1419/1458 (97%)	1393 (98%)	26 (2%)	69	53

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73[A]	PHE
1	A	73[B]	PHE

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Mol	Chain	Res	Type
1	A	82	ARG
2	B	70	PHE
1	C	46[A]	ARG
1	C	46[B]	ARG
1	C	102[A]	GLU
1	C	102[B]	GLU
3	D	57	LYS
4	E	19[A]	MET
4	E	19[B]	MET
4	E	132	SER
4	E	194	ARG
2	F	70	PHE
3	G	70[A]	TYR
3	G	70[B]	TYR
3	G	146[A]	SER
3	G	146[B]	SER
3	G	146[C]	SER
3	G	164	MET
4	H	185	ASN
4	H	186[A]	ASP
4	H	186[B]	ASP
4	H	194	ARG
4	H	206[A]	ARG
4	H	206[B]	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
4	E	28	ASN
3	G	188	ASN
4	H	120	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 ligands 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	$\# Z > 2$
6	GOL	A	303	-	5,5,5	0.93	0	5,5,5	1.38	1 (20%)
6	GOL	A	302	-	5,5,5	1.29	1 (20%)	5,5,5	0.93	0
6	GOL	C	301	-	5,5,5	0.87	0	5,5,5	0.95	0
5	ACT	A	301	-	1,3,3	7.50	1 (100%)	0,3,3	0.00	-
6	GOL	F	101	-	5,5,5	1.23	0	5,5,5	0.99	0
5	ACT	H	1601	-	1,3,3	8.19	1 (100%)	0,3,3	0.00	-

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
6	GOL	A	303	-	-	2/4/4/4	-
6	GOL	C	301	-	-	2/4/4/4	-
6	GOL	A	302	-	-	0/4/4/4	-
6	GOL	F	101	-	-	0/4/4/4	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	H	1601	ACT	CH3-C	8.19	1.59	1.48
5	A	301	ACT	CH3-C	7.50	1.58	1.48
6	A	302	GOL	C3-C2	2.34	1.61	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	303	GOL	C3-C2-C1	-2.45	102.18	111.70

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	303	GOL	C1-C2-C3-O3
6	C	301	GOL	O1-C1-C2-C3
6	A	303	GOL	O2-C2-C3-O3
6	C	301	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	301	GOL	5	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
1	A	265/271 (97%)	-0.10	8 (3%)	50	59	27, 39, 72, 89	11 (4%)
1	C	267/271 (98%)	-0.03	7 (2%)	56	65	27, 37, 62, 83	15 (5%)
2	B	98/100 (98%)	0.82	21 (21%)	0	1	33, 60, 83, 93	5 (5%)
2	F	100/100 (100%)	-0.19	0	100	100	29, 41, 58, 65	5 (5%)
3	D	188/204 (92%)	0.73	35 (18%)	1	1	29, 49, 89, 114	7 (3%)
3	G	201/204 (98%)	-0.25	0	100	100	26, 36, 56, 71	15 (7%)
4	E	241/246 (97%)	0.17	22 (9%)	9	15	31, 50, 91, 107	10 (4%)
4	H	244/246 (99%)	-0.07	1 (0%)	92	95	26, 36, 54, 78	15 (6%)
All	All	1604/1642 (97%)	0.09	94 (5%)	22	30	26, 41, 80, 114	83 (5%)

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	180	PHE	6.0
4	E	225	THR	5.8
3	D	162	ARG	5.6
3	D	197	PHE	5.1
3	D	163	SER	4.9
3	D	117	ALA	4.8
3	D	182	CYS	4.7
4	E	199	ALA	4.4
4	E	200	THR	4.4
3	D	149	SER	4.4
3	D	119	TYR	4.3
3	D	181	ALA	4.2
2	B	78	TYR	4.2
4	E	205	PRO	4.2
3	D	194	ASP	4.1
4	E	167	VAL	4.1

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Mol	Chain	Res	Type	RSRZ
2	B	76	ASP	3.9
1	A	222	VAL	3.9
4	E	201	PHE	3.9
3	D	147	LYS	3.6
1	A	223	GLN	3.6
1	A	221	ILE	3.5
4	E	219	SER	3.5
2	B	43	GLY	3.4
2	B	79	ALA	3.4
2	B	73	THR	3.3
3	D	138[A]	ASP	3.3
3	D	190	ILE	3.3
1	C	211[A]	TYR	3.3
4	E	198	SER	3.3
2	B	92	ILE	3.2
1	A	196	THR	3.2
4	E	136	ILE	3.2
4	E	209	PHE	3.2
4	E	137	SER	3.2
4	E	204	ASN	3.2
4	E	227	ASP	3.1
3	D	143	VAL	3.1
4	E	141	LYS	3.1
3	D	186	PHE	3.0
3	D	132	CYS	3.0
1	C	252	ASN	3.0
3	D	165	ASP	3.0
2	B	0	MET	3.0
3	D	131	VAL	2.9
2	B	1	ILE	2.9
2	B	46	ILE	2.9
3	D	113	ASN	2.9
1	A	17	HIS	2.9
1	A	246	LEU	2.9
4	H	184	LEU	2.9
3	D	183	ALA	2.8
3	D	191	ILE	2.8
4	E	221	ASN	2.8
2	B	45	ARG	2.8
3	D	196	PHE	2.8
2	B	42	ASN	2.8
3	D	133	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
4	E	222	ASP	2.7
2	B	88	SER	2.7
1	C	0	MET	2.7
3	D	111	ILE	2.7
3	D	187	ASN	2.7
1	A	188	ARG	2.6
3	D	164	MET	2.6
3	D	185	ALA	2.6
1	C	17	HIS	2.6
1	C	195	VAL	2.6
4	E	197	VAL	2.5
4	E	224	TRP	2.5
3	D	146	SER	2.4
3	D	166	PHE	2.4
2	B	74	GLU	2.4
3	D	195	THR	2.4
3	D	175	SER	2.4
2	B	44	GLU	2.3
2	B	95	TRP	2.3
4	E	203	GLN	2.3
4	E	98	GLY	2.3
2	B	48	LYS	2.3
2	B	77	GLU	2.3
1	A	0	MET	2.2
2	B	75	LYS	2.2
3	D	188	ASN	2.2
1	C	221	ILE	2.2
2	B	96	ASP	2.2
3	D	159	LEU	2.2
2	B	97	ARG	2.2
2	B	80	CYS	2.2
1	C	191	THR	2.2
4	E	97	THR	2.1
3	D	192	PRO	2.1
3	D	134	PHE	2.0
4	E	207	ASN	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [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
5	ACT	A	301	4/4	0.82	0.19	72,75,75,75	0
6	GOL	A	302	6/6	0.87	0.19	42,45,45,46	6
6	GOL	C	301	6/6	0.90	0.13	58,60,61,62	0
5	ACT	H	1601	4/4	0.94	0.10	52,57,57,60	0
6	GOL	A	303	6/6	0.96	0.11	35,38,43,44	0
6	GOL	F	101	6/6	0.97	0.08	34,35,37,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.