



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2026 – 12:01 AM UTC

PDB ID : 4PJF / pdb_00004pjf
Title : Structure of human MR1-Ac-6-FP in complex with human MAIT B-C10 TCR
Authors : Birkinshaw, R.W.; Rossjohn, J.
Deposited on : 2014-05-12
Resolution : 2.45 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

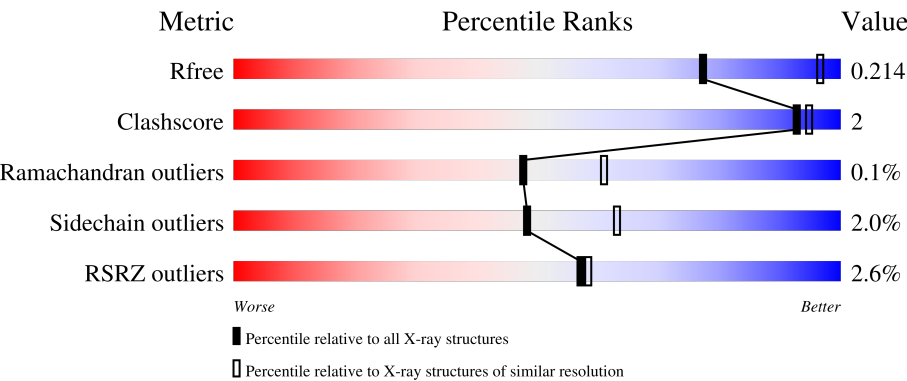
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	271	91%6%.
1	C	271	3% 87%7%5%
2	B	100	93%6%.
2	D	100	5% 91%6%.
3	E	205	89%7%.

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Mol	Chain	Length	Quality of chain
3	G	205	7% 83% 9% 7%
4	F	246	% 90% 7% ••
4	H	246	2% 92% 5% •

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 13209 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	263	Total	C	N	O	S	0	0	0
			2141	1371	372	388	10			
1	C	257	Total	C	N	O	S	0	0	0
			2065	1320	360	375	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP Q95460
A	261	SER	CYS	engineered mutation	UNP Q95460
C	0	MET	-	initiating methionine	UNP Q95460
C	261	SER	CYS	engineered mutation	UNP Q95460

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	0	0	0
			792	506	136	147	3			
2	D	97	Total	C	N	O	S	0	0	0
			762	491	125	144	2			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	initiating methionine	UNP P61769
D	0	MET	-	initiating methionine	UNP P61769

- Molecule 3 is a protein called TCR-alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	197	Total	C	N	O	S	0	0	0
			1491	947	235	300	9			

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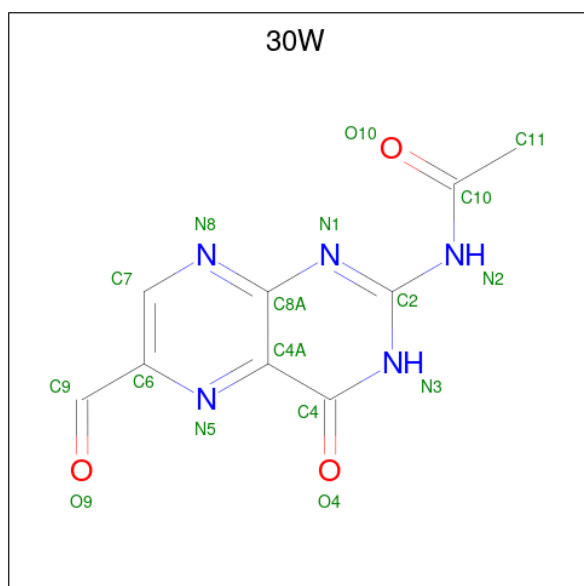
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	190	Total	C	N	O	S	0	0	0
			1407	902	221	275	9			

- Molecule 4 is a protein called TCR-beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	240	Total	C	N	O	S	0	0	0
			1842	1163	316	354	9			
4	H	240	Total	C	N	O	S	0	0	0
			1831	1153	320	349	9			

- Molecule 5 is N-(6-formyl-4-oxo-3,4-dihydropteridin-2-yl)acetamide (CCD ID: 30W) (formula: C₉H₇N₅O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			16	9	5	2		
5	C	1	Total	C	N	O	0	0
			16	9	5	2		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	160	Total	O	0	0
			160	160		
7	B	62	Total	O	0	0
			62	62		
7	C	135	Total	O	0	0
			135	135		
7	D	29	Total	O	0	0
			29	29		
7	E	91	Total	O	0	0
			91	91		
7	F	147	Total	O	0	0
			147	147		
7	G	78	Total	O	0	0
			78	78		
7	H	132	Total	O	0	0
			132	132		

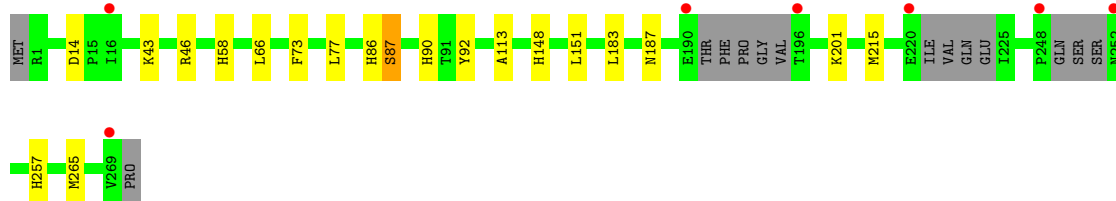
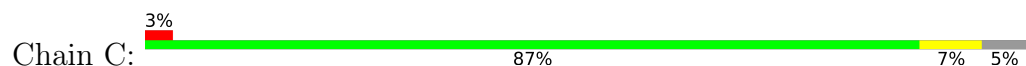
3 Residue-property plots [i](#)

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



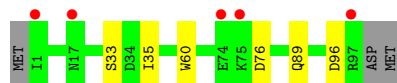
- Molecule 1: Major histocompatibility complex class I-related gene protein



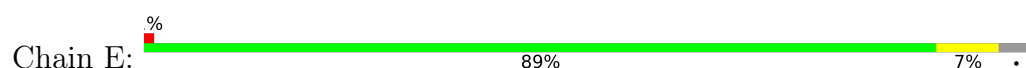
- Molecule 2: Beta-2-microglobulin

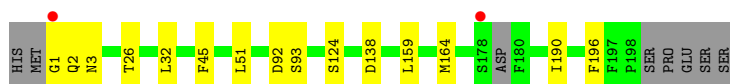


- Molecule 2: Beta-2-microglobulin

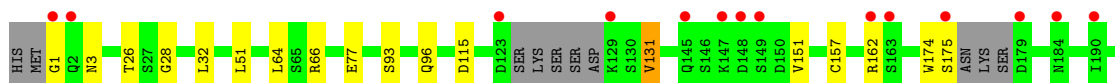
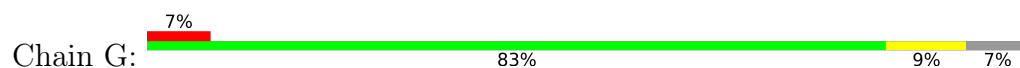


- Molecule 3: TCR-alpha

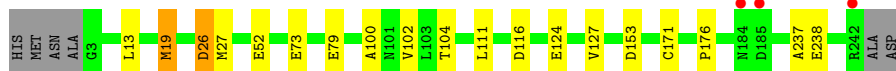
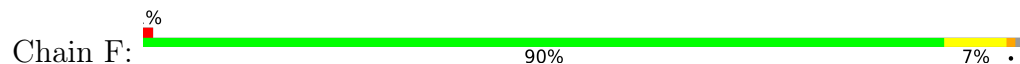




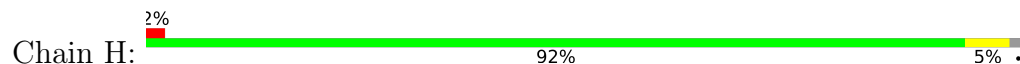
• Molecule 3: TCR-alpha



• Molecule 4: TCR-beta



• Molecule 4: TCR-beta



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	218.99Å 71.70Å 144.03Å 90.00° 104.85° 90.00°	Depositor
Resolution (Å)	35.87 – 2.45 35.87 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.0 (35.87-2.45) 99.0 (35.87-2.45)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.16 (at 2.45Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.178 , 0.212 0.179 , 0.214	Depositor DCC
R_{free} test set	3969 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	42.7	Xtriage
Anisotropy	0.460	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 57.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13209	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.58% 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: 30W, GOL

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	0.78	0/2206	1.23	2/3001 (0.1%)
1	C	0.77	1/2126 (0.0%)	1.22	1/2892 (0.0%)
2	B	0.73	0/815	1.15	3/1111 (0.3%)
2	D	0.69	0/785	1.16	2/1074 (0.2%)
3	E	0.77	0/1524	1.15	8/2073 (0.4%)
3	G	0.76	0/1438	1.10	4/1958 (0.2%)
4	F	0.73	0/1891	1.13	4/2582 (0.2%)
4	H	0.76	0/1879	1.12	5/2568 (0.2%)
All	All	0.76	1/12664 (0.0%)	1.16	29/17259 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	14	ASP	CA-C	6.60	1.57	1.53

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	87	SER	N-CA-C	8.90	123.45	112.23
4	F	26	ASP	CA-CB-CG	7.07	119.67	112.60
3	E	92	ASP	CA-CB-CG	6.46	119.06	112.60
4	H	116	ASP	CA-CB-CG	6.33	118.93	112.60
3	E	2	GLN	N-CA-C	6.31	121.12	112.04
3	G	3	ASN	N-CA-C	6.22	118.63	109.24
1	A	226	ASP	CA-CB-CG	5.84	118.44	112.60
1	A	31	HIS	N-CA-C	5.83	117.01	109.72
4	H	198	ALA	CA-C-N	5.83	128.09	120.28
4	H	198	ALA	C-N-CA	5.83	128.09	120.28
2	B	76	ASP	N-CA-C	5.73	118.02	109.25
4	F	116	ASP	CA-CB-CG	5.66	118.26	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	124	SER	CA-C-N	5.60	128.05	120.38
3	E	124	SER	C-N-CA	5.60	128.05	120.38
3	G	162	ARG	N-CA-C	5.55	118.04	111.33
4	F	124	GLU	N-CA-C	-5.39	99.94	108.73
3	E	3	ASN	N-CA-C	5.37	117.55	109.23
3	E	92	ASP	CA-C-N	5.36	127.72	120.38
3	E	92	ASP	C-N-CA	5.36	127.72	120.38
2	B	34	ASP	CA-CB-CG	5.29	117.89	112.60
2	D	76	ASP	N-CA-C	5.26	116.73	109.15
2	B	33	SER	N-CA-C	5.24	118.83	112.23
2	D	33	SER	N-CA-C	5.23	118.82	112.23
4	H	73	GLU	N-CA-C	5.14	117.86	109.59
3	E	196	PHE	CA-CB-CG	5.07	118.87	113.80
3	G	196	PHE	CA-CB-CG	5.05	118.85	113.80
4	H	124	GLU	N-CA-C	-5.04	100.31	108.52
3	G	115	ASP	CA-CB-CG	5.02	117.62	112.60
4	F	73	GLU	N-CA-C	5.01	117.66	109.59

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2141	0	2016	7	0
1	C	2065	0	1901	11	0
2	B	792	0	730	2	0
2	D	762	0	680	1	0
3	E	1491	0	1358	5	0
3	G	1407	0	1257	6	0
4	F	1842	0	1705	9	0
4	H	1831	0	1691	3	0
5	A	16	0	6	0	0
5	C	16	0	6	0	0
6	B	12	0	16	1	0
7	A	160	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	B	62	0	0	0	0
7	C	135	0	0	0	0
7	D	29	0	0	0	0
7	E	91	0	0	1	0
7	F	147	0	0	2	0
7	G	78	0	0	0	0
7	H	132	0	0	0	0
All	All	13209	0	11366	36	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 (36) 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)
4:F:13:LEU:HD21	4:F:19:MET:HG3	1.59	0.84
3:G:131:VAL:HG23	3:G:174:TRP:HB3	1.75	0.68
1:A:77:LEU:HD13	1:A:92:TYR:HB2	1.86	0.57
3:G:1:GLY:HA3	3:G:26:THR:HA	1.88	0.56
2:B:96:ASP:HB3	6:B:102:GOL:H12	1.87	0.55
3:G:28:GLY:HA3	3:G:93:SER:HB3	1.88	0.55
3:E:1:GLY:HA3	3:E:26:THR:HA	1.89	0.54
1:C:187:ASN:HD21	1:C:201:LYS:HE3	1.72	0.54
1:A:215:MET:HG3	1:A:257:HIS:CD2	2.42	0.54
1:C:215:MET:HG3	1:C:257:HIS:CD2	2.44	0.51
1:A:113:ALA:HB2	2:B:60:TRP:CE2	2.46	0.51
4:F:153:ASP:CG	4:F:176:PRO:HG3	2.36	0.51
1:C:43:LYS:HD2	1:C:66:LEU:HD13	1.93	0.50
1:A:186:VAL:HG11	1:A:269:VAL:HG22	1.93	0.50
1:A:189:LYS:HD3	1:C:87:SER:HB3	1.96	0.47
1:C:113:ALA:HB2	2:D:60:TRP:CE2	2.49	0.47
4:H:154:HIS:HB3	4:H:215:TYR:HB2	1.98	0.45
1:C:151:LEU:HD22	3:E:51:LEU:HD12	1.98	0.45
3:G:151:VAL:HG12	3:G:175:SER:HB2	1.99	0.45
3:G:64:LEU:HD21	3:G:66:ARG:HD3	1.99	0.45
3:E:45:PHE:HB2	4:F:102:VAL:HG13	2.01	0.43
4:F:79:GLU:HB2	7:F:422:HOH:O	2.17	0.43
1:C:148:HIS:HB3	4:F:100:ALA:HB1	2.01	0.43
4:F:19:MET:HE1	4:F:111:LEU:HD11	2.00	0.43
1:A:4:SER:HB3	1:A:99:GLU:HG2	2.00	0.43
1:A:151:LEU:HD22	3:G:51:LEU:HD12	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:86:HIS:HB3	1:C:90:HIS:NE2	2.34	0.42
4:F:27:MET:HE1	4:F:104:THR:HB	2.02	0.42
4:H:127:VAL:HG23	4:H:237:ALA:HB3	2.02	0.41
1:C:77:LEU:HD13	1:C:92:TYR:HB2	2.01	0.41
1:C:58:HIS:HD2	7:E:327:HOH:O	2.03	0.41
4:H:27:MET:HE1	4:H:104:THR:HB	2.02	0.41
3:E:159:LEU:HB3	4:F:171:CYS:HB2	2.02	0.41
1:C:46:ARG:HA	1:C:46:ARG:HD3	1.83	0.40
3:E:164:MET:HG3	7:F:409:HOH:O	2.21	0.40
4:F:127:VAL:HG23	4:F:237:ALA:HB3	2.02	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	259/271 (96%)	252 (97%)	7 (3%)	0	100	100
1	C	249/271 (92%)	244 (98%)	5 (2%)	0	100	100
2	B	97/100 (97%)	96 (99%)	1 (1%)	0	100	100
2	D	95/100 (95%)	93 (98%)	1 (1%)	1 (1%)	11	12
3	E	193/205 (94%)	192 (100%)	1 (0%)	0	100	100
3	G	184/205 (90%)	182 (99%)	2 (1%)	0	100	100
4	F	238/246 (97%)	233 (98%)	5 (2%)	0	100	100
4	H	238/246 (97%)	235 (99%)	2 (1%)	1 (0%)	30	37
All	All	1553/1644 (94%)	1527 (98%)	24 (2%)	2 (0%)	48	61

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	96	ASP
4	H	153	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	224/241 (93%)	220 (98%)	4 (2%)	51	65
1	C	209/241 (87%)	206 (99%)	3 (1%)	59	72
2	B	85/95 (90%)	84 (99%)	1 (1%)	63	74
2	D	78/95 (82%)	76 (97%)	2 (3%)	40	55
3	E	156/181 (86%)	152 (97%)	4 (3%)	40	55
3	G	137/181 (76%)	132 (96%)	5 (4%)	31	45
4	F	193/213 (91%)	189 (98%)	4 (2%)	47	62
4	H	190/213 (89%)	188 (99%)	2 (1%)	65	76
All	All	1272/1460 (87%)	1247 (98%)	25 (2%)	48	63

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	ARG
1	A	73	PHE
1	A	183	LEU
1	A	245	GLU
2	B	35	ILE
1	C	73	PHE
1	C	183	LEU
1	C	265	MET
2	D	35	ILE
2	D	89	GLN
3	E	32	LEU
3	E	93	SER
3	E	138	ASP
3	E	190	ILE

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Mol	Chain	Res	Type
4	F	19	MET
4	F	26	ASP
4	F	52	GLU
4	F	238	GLU
3	G	32	LEU
3	G	77	GLU
3	G	96	GLN
3	G	131	VAL
3	G	157	CYS
4	H	14	LYS
4	H	139	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	147	GLN
1	A	203	HIS
2	B	13	HIS
2	B	51	HIS
1	C	58	HIS
1	C	85	ASN
1	C	147	GLN
1	C	187	ASN
2	D	13	HIS
2	D	51	HIS
2	D	89	GLN
3	E	79	GLN
3	E	120	GLN
3	G	3	ASN
3	G	79	GLN
3	G	120	GLN
3	G	169	ASN
4	H	30	ASN
4	H	139	GLN
4	H	184	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 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$
5	30W	C	301	1	16,17,18	0.77	1 (6%)	18,24,25	2.14	3 (16%)
5	30W	A	301	1	16,17,18	0.70	1 (6%)	18,24,25	2.00	3 (16%)
6	GOL	B	101	-	5,5,5	0.10	0	5,5,5	0.36	0
6	GOL	B	102	-	5,5,5	0.15	0	5,5,5	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
5	30W	C	301	1	-	0/4/4/6	0/2/2/2
5	30W	A	301	1	-	0/4/4/6	0/2/2/2
6	GOL	B	101	-	-	0/4/4/4	-
6	GOL	B	102	-	-	0/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	301	30W	C2-N2	-2.15	1.33	1.37
5	C	301	30W	C4A-C8A	2.02	1.44	1.40

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	301	30W	N2-C2-N3	7.58	130.71	117.96
5	A	301	30W	N2-C2-N3	6.80	129.40	117.96
5	A	301	30W	N2-C2-N1	-3.55	112.15	117.87
5	C	301	30W	N2-C2-N1	-3.40	112.38	117.87
5	C	301	30W	C4A-C8A-N1	-2.51	119.67	123.19
5	A	301	30W	C4A-C8A-N1	-2.14	120.19	123.19

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	102	GOL	1	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	263/271 (97%)	-0.37	4 (1%) 72 74	28, 40, 66, 88	0
1	C	257/271 (94%)	-0.14	7 (2%) 56 57	29, 44, 76, 103	0
2	B	99/100 (99%)	-0.25	1 (1%) 79 80	31, 47, 72, 82	0
2	D	97/100 (97%)	0.39	5 (5%) 33 30	36, 65, 96, 109	0
3	E	197/205 (96%)	-0.07	2 (1%) 79 80	30, 45, 73, 83	0
3	G	190/205 (92%)	0.22	15 (7%) 18 16	25, 49, 97, 130	0
4	F	240/246 (97%)	-0.22	3 (1%) 75 77	32, 43, 66, 94	0
4	H	240/246 (97%)	-0.10	4 (1%) 69 71	26, 45, 87, 128	0
All	All	1583/1644 (96%)	-0.11	41 (2%) 57 58	25, 45, 82, 130	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	G	123	ASP	5.8
1	C	252	ASN	4.3
3	E	178	SER	4.1
1	C	190	GLU	3.9
3	G	149	SER	3.9
1	A	246	LEU	3.8
1	C	220	GLU	3.3
3	G	1	GLY	3.3
3	E	1	GLY	3.3
2	D	97	ARG	3.3
3	G	148	ASP	3.2
4	F	242	ARG	3.2
1	C	248	PRO	3.2
2	D	17	ASN	3.2
2	D	1	ILE	3.2
1	A	221	ILE	3.2

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Mol	Chain	Res	Type	RSRZ
3	G	129	LYS	3.0
3	G	175	SER	2.7
2	B	0	MET	2.6
1	C	269	VAL	2.4
4	F	184	ASN	2.4
2	D	75	LYS	2.4
4	H	200	PHE	2.4
2	D	74	GLU	2.4
3	G	179	ASP	2.3
1	C	196	THR	2.3
1	A	222	VAL	2.3
3	G	194	ASP	2.2
3	G	190	ILE	2.2
3	G	163	SER	2.2
4	F	185	ASP	2.2
1	A	253	LEU	2.1
4	H	242	ARG	2.1
1	C	16	ILE	2.1
3	G	2	GLN	2.1
3	G	162	ARG	2.1
4	H	238	GLU	2.1
3	G	147	LYS	2.1
3	G	184	ASN	2.0
3	G	145	GLN	2.0
4	H	226	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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($\text{\AA}^2$)	Q<0.9
6	GOL	B	102	6/6	0.89	0.17	52,60,62,62	0
6	GOL	B	101	6/6	0.95	0.08	35,38,43,45	0
5	30W	A	301	16/17	0.97	0.05	29,31,39,40	0
5	30W	C	301	16/17	0.97	0.05	29,35,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.