



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 03:25 pm BST

PDB ID : 5U16  
Title : Structure of human MR1-2-OH-1-NA in complex with human MAIT A-F7 TCR  
Authors : Keller, A.N.; Birkinshaw, R.W.; Rossjohn, J.  
Deposited on : 2016-11-27  
Resolution : 2.00 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](mailto: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.

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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

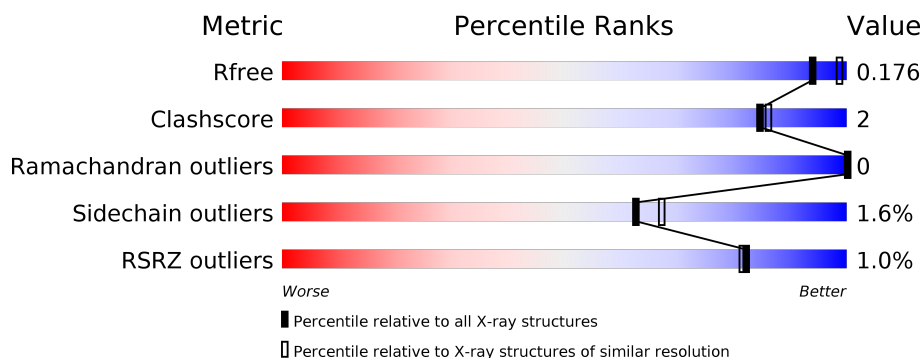
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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  $\geq 3$ , 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>%</div> <div> <div></div> <div>92%</div> <div>6%</div> <div>.</div> </div> </div>
1	C	271	<div> <div>92%</div> <div>6%</div> <div>.</div> </div>
2	B	99	<div> <div>91%</div> <div>6%</div> <div>.</div> </div>
2	D	99	<div> <div>%</div> <div>96%</div> <div>.</div> </div>
3	E	203	<div> <div>2%</div> <div>84%</div> <div>6%</div> <div>9%</div> <div>.</div> </div>
3	G	203	<div> <div>94%</div> <div>.</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
4	F	245	<div><div></div><div>2%</div><div>90%</div><div>7%</div><div>••</div></div>
4	H	245	<div><div></div><div>92%</div><div>6%</div><div>••</div></div>

## 2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 14076 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	1	0
			2191	1402	377	401	11			
1	C	265	Total	C	N	O	S	0	0	0
			2181	1399	374	398	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	conflict	UNP Q95460
C	0	MET	-	initiating methionine	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	96	Total	C	N	O	S	0	0	0
			797	511	134	150	2			
2	D	99	Total	C	N	O	S	0	0	0
			828	528	140	157	3			

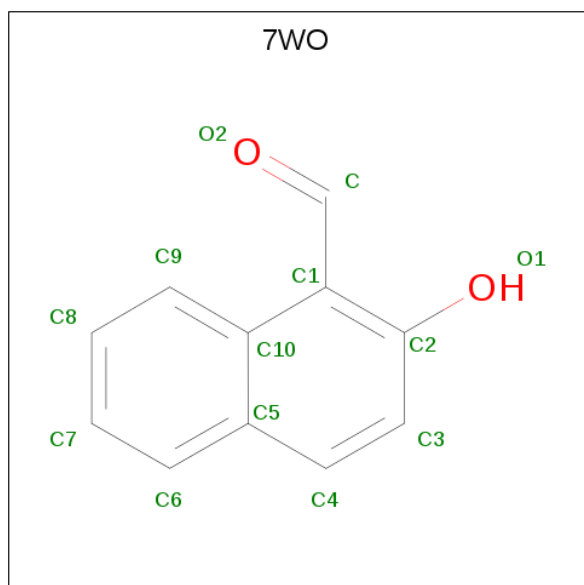
- Molecule 3 is a protein called MAIT T-cell receptor alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	185	Total	C	N	O	S	0	2	0
			1431	912	228	283	8			
3	G	200	Total	C	N	O	S	0	0	0
			1565	987	253	316	9			

- Molecule 4 is a protein called MAIT T-cell receptor beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	240	Total	C	N	O	S	0	0	0
			1877	1184	323	361	9			
4	H	243	Total	C	N	O	S	0	1	0
			1929	1212	335	373	9			

- Molecule 5 is 2-hydroxynaphthalene-1-carbaldehyde (three-letter code: 7WO) (formula:  $C_{11}H_8O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			12	11	1		
5	C	1	Total	C	O	0	0
			12	11	1		

- 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	B	1	Total	C	O	0	0
			6	3	3		
6	C	1	Total	C	O	0	0
			6	3	3		
6	D	1	Total	C	O	0	0
			6	3	3		
6	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	2	Total	Cl	0	0
			2	2		
7	C	1	Total	Cl	0	0
			1	1		

- Molecule 8 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	H	1	Total	Na	0	0
			1	1		
8	D	1	Total	Na	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	F	1	Total 1	Na 1	0	0

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	225	Total 225	O 225	0	0
9	B	59	Total 59	O 59	0	0
9	C	200	Total 200	O 200	0	0
9	D	90	Total 90	O 90	0	0
9	E	108	Total 108	O 108	0	0
9	F	101	Total 101	O 101	0	0
9	G	196	Total 196	O 196	0	0
9	H	238	Total 238	O 238	0	0

### 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.

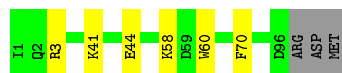
- Molecule 1: Major histocompatibility complex class I-related gene protein



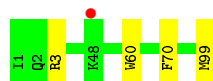
- Molecule 1: Major histocompatibility complex class I-related gene protein



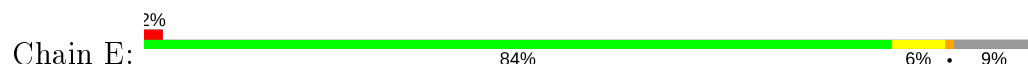
- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin



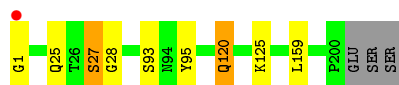
- Molecule 3: MAIT T-cell receptor alpha chain




- Molecule 3: MAIT T-cell receptor alpha chain



Chain G:  94% ..



- Molecule 4: MAIT T-cell receptor beta chain

Chain F:  90% 7% ..



- Molecule 4: MAIT T-cell receptor beta chain

Chain H:  92% 6% ..



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	216.44Å 69.85Å 143.12Å 90.00° 104.35° 90.00°	Depositor
Resolution (Å)	69.33 – 2.00 75.15 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (69.33-2.00) 99.8 (75.15-2.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.70 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.173 , 0.219 0.176 , 0.176	Depositor DCC
$R_{free}$ test set	7037 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	21.4	Xtriage
Anisotropy	0.460	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 48.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	14076	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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, 7WO, CL, NA

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	0.56	0/2260	0.69	4/3071 (0.1%)
1	C	0.48	0/2247	0.61	0/3055
2	B	0.40	0/820	0.54	0/1112
2	D	0.47	0/851	0.58	0/1152
3	E	0.47	0/1467	0.59	0/1990
3	G	0.54	0/1600	0.64	0/2168
4	F	0.43	0/1928	0.58	0/2626
4	H	0.54	0/1983	0.63	0/2696
All	All	0.50	0/13156	0.62	4/17870 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1	ARG	NE-CZ-NH2	-10.47	115.06	120.30
1	A	46	ARG	NE-CZ-NH2	-7.99	116.30	120.30
1	A	46	ARG	NE-CZ-NH1	6.88	123.74	120.30
1	A	1	ARG	NE-CZ-NH1	6.47	123.53	120.30

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	2191	0	2078	11	0
1	C	2181	0	2074	13	0
2	B	797	0	764	4	0
2	D	828	0	794	3	0
3	E	1431	0	1344	7	0
3	G	1565	0	1486	9	0
4	F	1877	0	1761	10	0
4	H	1929	0	1831	12	0
5	A	12	0	0	0	0
5	C	12	0	0	1	0
6	A	6	0	8	2	0
6	B	6	0	8	0	0
6	C	6	0	8	2	0
6	D	12	0	16	0	0
7	A	2	0	0	0	0
7	C	1	0	0	0	0
8	D	1	0	0	0	0
8	F	1	0	0	0	0
8	H	1	0	0	0	0
9	A	225	0	0	3	0
9	B	59	0	0	0	0
9	C	200	0	0	2	0
9	D	90	0	0	1	0
9	E	108	0	0	1	0
9	F	101	0	0	1	0
9	G	196	0	0	1	0
9	H	238	0	0	2	0
All	All	14076	0	12172	57	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 (57) 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:230:ILE:HD11	4:H:206:ARG:HD3	1.67	0.76
4:H:210:ARG:NH1	9:H:401:HOH:O	2.10	0.72
3:G:1:GLY:N	3:G:27:SER:H	1.87	0.72
1:C:152:TYR:CD1	4:F:100:GLY:HA3	2.26	0.70
1:A:149:GLU:OE1	9:A:401:HOH:O	2.09	0.69
4:F:208:HIS:NE2	4:F:239:GLU:OE1	2.16	0.69
9:C:435:HOH:O	3:G:125:LYS:HE2	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:95:TYR:OH	9:E:301:HOH:O	2.13	0.64
1:C:246:LEU:O	9:C:401:HOH:O	2.15	0.63
4:H:1:ASN:N	4:H:26:ASP:OD2	2.31	0.61
2:B:41:LYS:O	2:B:44:GLU:HG2	2.01	0.61
3:E:1:GLY:HA2	3:E:27:SER:H	1.66	0.61
3:G:25:GLN:NE2	9:G:302:HOH:O	2.34	0.60
3:G:1:GLY:H3	3:G:27:SER:H	1.50	0.59
4:F:154:ASP:OD1	4:F:177:PRO:HG2	2.03	0.58
1:C:113:ALA:HB2	2:D:60:TRP:CE2	2.38	0.58
1:A:113:ALA:HB2	2:B:60:TRP:CE2	2.39	0.57
3:G:159:LEU:HB3	4:H:172:CYS:HB2	1.88	0.55
1:A:215:MET:HG3	1:A:257:HIS:CD2	2.42	0.55
3:E:28:GLY:HA3	3:E:93[A]:SER:OG	2.06	0.55
3:G:28:GLY:HA3	3:G:93:SER:OG	2.06	0.54
1:A:6:ARG:NH2	2:B:58:LYS:HD3	2.22	0.54
1:C:96:ILE:HD13	6:C:302:GOL:H31	1.90	0.53
4:H:131:PRO:HD3	4:H:144:LEU:HG	1.93	0.51
1:A:152:TYR:CD2	4:H:100:GLY:HA2	2.46	0.51
1:C:215:MET:HG3	1:C:257:HIS:CD2	2.46	0.50
4:H:72:ARG:NH2	9:H:407:HOH:O	2.44	0.49
1:A:258:VAL:HB	1:A:265:MET:HG3	1.95	0.49
1:A:149:GLU:OE1	4:H:100:GLY:HA3	2.14	0.48
4:F:117:ASP:OD1	4:F:119:LYS:HG2	2.13	0.48
6:A:302:GOL:H31	3:G:95:TYR:OH	2.15	0.47
1:C:94:ARG:HH22	6:C:302:GOL:H12	1.79	0.47
4:F:131:PRO:HD3	4:F:144:LEU:HG	1.97	0.47
3:E:3:ASN:ND2	3:E:5:ASP:OD2	2.45	0.47
1:A:160:GLU:OE2	9:A:402:HOH:O	2.21	0.46
3:E:66[A]:ARG:HH11	3:E:66[A]:ARG:HG2	1.82	0.45
3:E:50:VAL:O	3:E:66[B]:ARG:HD3	2.17	0.45
1:C:116:GLY:O	2:D:3:ARG:NH2	2.50	0.45
4:F:155:HIS:HB3	4:F:216:TYR:HB2	1.98	0.44
3:E:134:PHE:HB2	3:E:186:PHE:CE2	2.53	0.43
4:F:224:TRP:CE2	4:F:226:GLN:HB2	2.54	0.43
1:C:152:TYR:CE1	4:F:100:GLY:HA3	2.53	0.43
4:F:9:LYS:HG3	4:F:10:PHE:CE1	2.55	0.42
6:A:302:GOL:H31	3:G:95:TYR:CZ	2.54	0.42
1:A:116:GLY:O	2:B:3:ARG:NH2	2.52	0.42
4:H:86:THR:HG23	4:H:113:THR:HA	2.01	0.42
4:H:218:LEU:HD22	4:H:231:PRO:HG2	2.02	0.42
1:A:265:MET:HG2	9:A:540:HOH:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:186:VAL:HG11	1:C:269:VAL:HG22	2.02	0.41
1:A:77:LEU:HD13	1:A:92:TYR:HB2	2.03	0.41
1:C:223:GLN:HG3	1:C:223:GLN:H	1.66	0.41
3:G:120:GLN:O	4:H:132:SER:HB2	2.21	0.41
1:C:43:LYS:NZ	5:C:301:7WO:O1	2.53	0.41
2:D:3:ARG:HD2	9:D:257:HOH:O	2.19	0.41
4:F:11:GLN:NE2	9:F:415:HOH:O	2.54	0.40
1:C:80:LEU:HA	1:C:80:LEU:HD23	1.87	0.40
4:H:224:TRP:CZ2	4:H:226:GLN:HB2	2.55	0.40

There are no symmetry-related clashes.

## 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	262/271 (97%)	257 (98%)	5 (2%)	0	100	100
1	C	261/271 (96%)	254 (97%)	7 (3%)	0	100	100
2	B	94/99 (95%)	93 (99%)	1 (1%)	0	100	100
2	D	97/99 (98%)	96 (99%)	1 (1%)	0	100	100
3	E	179/203 (88%)	175 (98%)	4 (2%)	0	100	100
3	G	198/203 (98%)	192 (97%)	6 (3%)	0	100	100
4	F	238/245 (97%)	236 (99%)	2 (1%)	0	100	100
4	H	242/245 (99%)	237 (98%)	5 (2%)	0	100	100
All	All	1571/1636 (96%)	1540 (98%)	31 (2%)	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	234/241 (97%)	232 (99%)	2 (1%)	78	83
1	C	232/241 (96%)	230 (99%)	2 (1%)	78	83
2	B	90/94 (96%)	89 (99%)	1 (1%)	73	78
2	D	94/94 (100%)	92 (98%)	2 (2%)	53	57
3	E	155/180 (86%)	149 (96%)	6 (4%)	32	30
3	G	176/180 (98%)	174 (99%)	2 (1%)	73	78
4	F	202/211 (96%)	197 (98%)	5 (2%)	47	49
4	H	211/211 (100%)	208 (99%)	3 (1%)	67	72
All	All	1394/1452 (96%)	1371 (98%)	23 (2%)	62	67

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	PHE
1	A	261	SER
2	B	70	PHE
1	C	73	PHE
1	C	224	GLU
2	D	70	PHE
2	D	99	MET
3	E	27	SER
3	E	66[A]	ARG
3	E	66[B]	ARG
3	E	124	SER
3	E	138	ASP
3	E	151	VAL
4	F	79	GLU
4	F	117	ASP
4	F	132	SER
4	F	194	ARG
4	F	226	GLN
3	G	27	SER

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Mol	Chain	Res	Type
3	G	120	GLN
4	H	132	SER
4	H	194	ARG
4	H	218	LEU

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

Mol	Chain	Res	Type
1	C	268	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 13 ligands modelled in this entry, 6 are monoatomic - leaving 7 for Mogul analysis.

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	C	302	-	5,5,5	0.37	0	5,5,5	0.32	0
6	GOL	D	103	-	5,5,5	0.40	0	5,5,5	0.53	0
5	7WO	C	301	1	13,13,14	0.74	1 (7%)	18,18,19	0.64	1 (5%)



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
6	GOL	D	102	-	5,5,5	0.52	0	5,5,5	0.39	0
6	GOL	B	101	-	5,5,5	0.34	0	5,5,5	0.68	0
5	7WO	A	301	1	13,13,14	0.75	1 (7%)	18,18,19	0.68	1 (5%)
6	GOL	A	302	-	5,5,5	0.27	0	5,5,5	0.54	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
6	GOL	C	302	-	-	4/4/4/4	-
6	GOL	D	103	-	-	0/4/4/4	-
5	7WO	C	301	1	-	-	0/2/2/2
6	GOL	D	102	-	-	2/4/4/4	-
6	GOL	B	101	-	-	3/4/4/4	-
5	7WO	A	301	1	-	-	0/2/2/2
6	GOL	A	302	-	-	2/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	301	7WO	C-C1	-2.52	1.45	1.51
5	C	301	7WO	C-C1	-2.48	1.45	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	301	7WO	O1-C2-C1	2.14	122.38	116.74
5	C	301	7WO	O1-C2-C1	2.01	122.04	116.74

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	101	GOL	C1-C2-C3-O3
6	A	302	GOL	O1-C1-C2-C3
6	C	302	GOL	O1-C1-C2-C3
6	C	302	GOL	C1-C2-C3-O3
6	D	102	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
6	B	101	GOL	O2-C2-C3-O3
6	A	302	GOL	O1-C1-C2-O2
6	B	101	GOL	O1-C1-C2-O2
6	D	102	GOL	O2-C2-C3-O3
6	C	302	GOL	O1-C1-C2-O2
6	C	302	GOL	O2-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	302	GOL	2	0
5	C	301	7WO	1	0
6	A	302	GOL	2	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	265/271 (97%)	-0.35	4 (1%) 73 72	14, 27, 70, 92	0
1	C	265/271 (97%)	-0.42	0 100 100	16, 27, 58, 92	0
2	B	96/99 (96%)	-0.34	0 100 100	20, 46, 70, 79	0
2	D	99/99 (100%)	-0.42	1 (1%) 82 81	16, 30, 60, 73	0
3	E	185/203 (91%)	0.01	4 (2%) 62 60	18, 37, 75, 95	0
3	G	200/203 (98%)	-0.33	1 (0%) 91 90	13, 24, 52, 73	0
4	F	240/245 (97%)	-0.07	6 (2%) 57 56	23, 41, 76, 93	0
4	H	243/245 (99%)	-0.40	0 100 100	15, 24, 47, 69	0
All	All	1593/1636 (97%)	-0.29	16 (1%) 82 81	13, 30, 68, 95	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	222	VAL	4.9
1	A	253	LEU	3.6
4	F	205	PRO	3.2
4	F	241	TRP	3.1
4	F	225	THR	3.1
3	E	124	SER	2.8
1	A	221	ILE	2.8
3	G	1	GLY	2.8
3	E	194	ASP	2.7
1	A	211	TYR	2.6
4	F	201	PHE	2.3
2	D	48	LYS	2.3
3	E	1	GLY	2.2
4	F	209	PHE	2.2
3	E	119	TYR	2.2
4	F	199	ALA	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 carbohydrates 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, 95<sup>th</sup> 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	101	6/6	0.85	0.22	39,44,52,54	0
6	GOL	C	302	6/6	0.89	0.23	47,54,56,60	0
7	CL	A	304	1/1	0.90	0.07	55,55,55,55	0
6	GOL	A	302	6/6	0.93	0.12	34,42,45,47	0
6	GOL	D	103	6/6	0.94	0.11	25,26,27,29	0
6	GOL	D	102	6/6	0.95	0.15	24,26,30,32	0
5	7WO	C	301	12/13	0.96	0.11	22,25,32,33	0
5	7WO	A	301	12/13	0.96	0.10	17,19,29,30	0
8	NA	F	301	1/1	0.96	0.08	34,34,34,34	0
8	NA	D	101	1/1	0.97	0.11	37,37,37,37	0
8	NA	H	301	1/1	0.98	0.07	30,30,30,30	0
7	CL	C	303	1/1	0.99	0.08	44,44,44,44	0
7	CL	A	303	1/1	0.99	0.06	27,27,27,27	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.