



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 13, 2022 – 12:08 PM EDT

PDB ID : 7NA5
Title : Structure of the H2DB-TCR ternary complex with HSF2 melanoma neoantigen
Authors : Patskovsky, Y.; Finnigan, J.; Patskovska, L.; Newman, J.; Bhardwaj, N.; Krogsgaard, M.
Deposited on : 2021-06-19
Resolution : 2.50 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 : 2.28.1
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) : 2.28.1

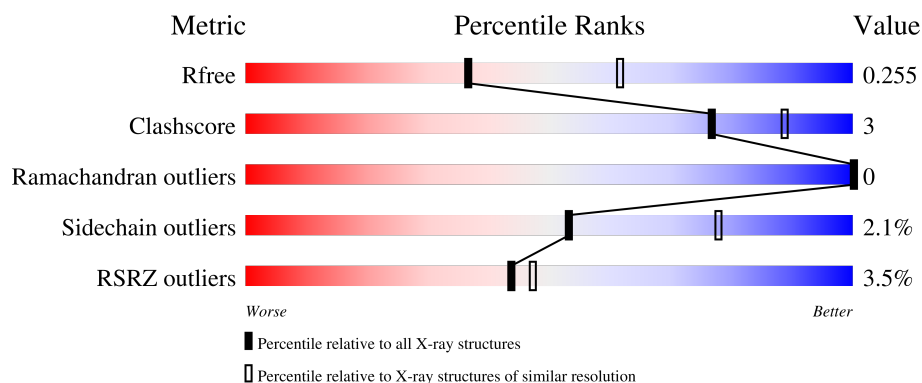
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	281	<div> <div>7%</div> <div> <div></div> <div>93%</div> <div>6%</div> </div> </div>
2	B	100	<div> <div></div> <div> <div></div> <div>90%</div> <div>9%</div> </div> </div>
3	C	9	<div> <div></div> <div> <div></div> <div>89%</div> <div>11%</div> </div> </div>
4	D	191	<div> <div>7%</div> <div> <div></div> <div>81%</div> <div>12%</div> <div>6%</div> </div> </div>
5	E	263	<div> <div>4%</div> <div> <div></div> <div>83%</div> <div>8%</div> <div>10%</div> </div> </div>

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6689 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 H-2 class I histocompatibility antigen, D-B alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	278	Total	C	N	O	S	0	1	0
			2249	1423	391	425	10			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP P01899

- 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
			825	525	139	158	3			

There is a discrepancy between the modelled and reference sequences:

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

- Molecule 3 is a protein called Heat shock factor protein 2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			79	52	15	12			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	5	ASN	LYS	engineered mutation	UNP P38533

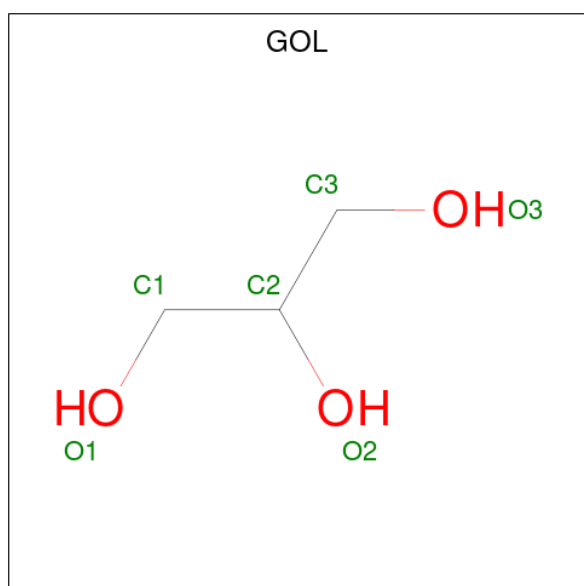
- Molecule 4 is a protein called 47BE7 TCR alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	180	Total	C	N	O	S	0	1	0
			1372	863	229	272	8			

- Molecule 5 is a protein called 47BE7 TCR beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	238	Total	C	N	O	S	0	0	0
			1858	1182	319	349	8			

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



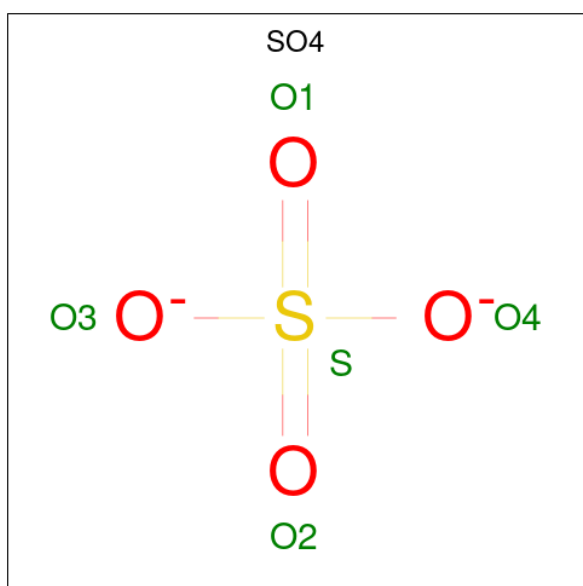
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	B	1	Total	C	O	0	0
			6	3	3		
6	B	1	Total	C	O	0	0
			6	3	3		

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

- Molecule 7 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	O	S	0	0
			5	4	1		
7	D	1	Total	O	S	0	0
			5	4	1		
7	E	1	Total	O	S	0	0
			5	4	1		
7	E	1	Total	O	S	0	0
			5	4	1		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	82	Total 82	O 82	0	0
8	B	43	Total 43	O 43	0	0
8	C	4	Total 4	O 4	0	0
8	D	41	Total 41	O 41	0	0
8	E	38	Total 38	O 38	0	0

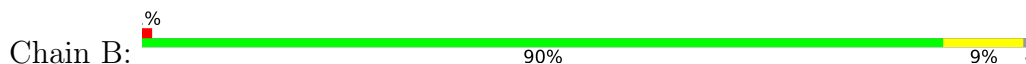
3 Residue-property plots [i](#)

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.

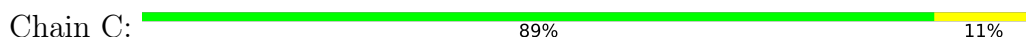
- Molecule 1: H-2 class I histocompatibility antigen, D-B alpha chain



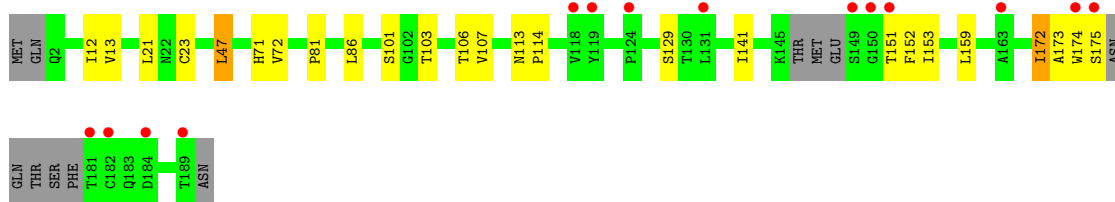
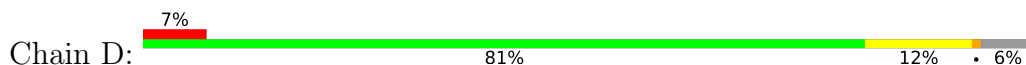
- Molecule 2: Beta-2-microglobulin



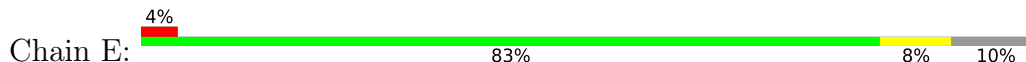
- Molecule 3: Heat shock factor protein 2



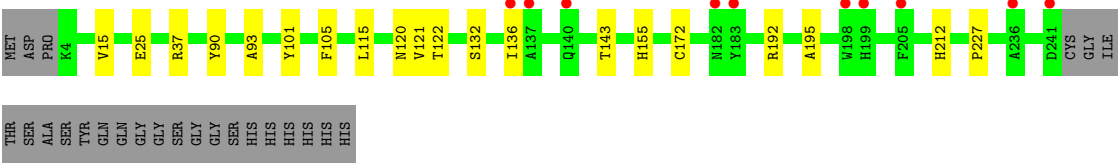
- Molecule 4: 47BE7 TCR alpha chain



- Molecule 5: 47BE7 TCR beta chain



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THR
SER
ALA
SER
TYR
GLN
GLN
GLY
GLY
SER
GLY
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HIS
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4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	155.44Å 191.71Å 67.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.01 – 2.50 50.02 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.3 (50.01-2.50) 99.3 (50.02-2.50)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.49 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.184 , 0.249 0.192 , 0.255	Depositor DCC
R_{free} test set	1073 reflections (3.04%)	wwPDB-VP
Wilson B-factor (Å ²)	47.2	Xtriage
Anisotropy	0.314	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.5	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	0.95	EDS
Total number of atoms	6689	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% 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, SO4

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	0.63	0/2318	0.80	0/3152
2	B	0.64	0/848	0.82	0/1148
3	C	0.53	0/81	0.72	0/107
4	D	0.70	0/1408	0.88	2/1913 (0.1%)
5	E	0.65	0/1914	0.81	0/2610
All	All	0.65	0/6569	0.82	2/8930 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	71[A]	HIS	CA-C-O	5.33	131.28	120.10
4	D	71[B]	HIS	CA-C-O	5.33	131.28	120.10

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	2249	0	2104	10	0
2	B	825	0	783	4	0
3	C	79	0	78	1	0
4	D	1372	0	1253	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	E	1858	0	1721	11	0
6	A	36	0	48	0	0
6	B	12	0	16	0	0
6	C	6	0	8	0	0
6	D	6	0	8	0	0
6	E	18	0	24	0	0
7	A	5	0	0	0	0
7	D	5	0	0	0	0
7	E	10	0	0	0	0
8	A	82	0	0	1	0
8	B	43	0	0	1	0
8	C	4	0	0	0	0
8	D	41	0	0	0	0
8	E	38	0	0	0	0
All	All	6689	0	6043	38	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (38) 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:D:23:CYS:HB3	4:D:72:VAL:HG13	1.81	0.61
1:A:11:ALA:HA	1:A:21:ARG:O	2.04	0.58
4:D:21:LEU:HD22	4:D:103:THR:HG21	1.88	0.55
2:B:3:ARG:NH1	2:B:59:ASP:O	2.39	0.55
5:E:155:HIS:HB3	5:E:212:HIS:HB2	1.89	0.55
5:E:136:ILE:HG23	5:E:195:ALA:HB1	1.90	0.54
5:E:37:ARG:HG3	5:E:90:TYR:CE2	2.44	0.53
4:D:172:ILE:HD11	4:D:174:TRP:CE3	2.44	0.53
4:D:159:LEU:C	4:D:159:LEU:HD12	2.31	0.51
1:A:187:ALA:HA	1:A:204:TRP:O	2.11	0.50
5:E:121:VAL:HG12	5:E:227:PRO:HB2	1.93	0.50
2:B:83:ASN:HA	8:B:216:HOH:O	2.12	0.50
1:A:193:PRO:HD3	8:A:455:HOH:O	2.11	0.48
2:B:36:GLU:HB3	2:B:83:ASN:HB2	1.97	0.46
5:E:143:THR:OG1	5:E:192:ARG:NH1	2.49	0.46
1:A:21:ARG:NH2	1:A:37:ASP:OD2	2.50	0.45
4:D:81:PRO:HA	4:D:107:VAL:HB	1.98	0.45
4:D:23:CYS:HB3	4:D:72:VAL:CG1	2.47	0.44
4:D:159:LEU:HB3	5:E:172:CYS:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:23:LEU:O	2:B:67:TYR:HA	2.18	0.44
4:D:153:ILE:HG12	4:D:173:ALA:HB2	1.99	0.44
5:E:93:ALA:HA	5:E:105:PHE:O	2.17	0.44
1:A:201:LEU:O	1:A:246:SER:HA	2.18	0.43
5:E:15:VAL:HA	5:E:115:LEU:O	2.17	0.43
5:E:136:ILE:HG23	5:E:195:ALA:CB	2.48	0.43
4:D:151:THR:HG23	4:D:174:TRP:O	2.19	0.43
4:D:47:LEU:HA	4:D:47:LEU:HD23	1.82	0.43
1:A:260:ARG:HA	1:A:270:LEU:O	2.19	0.42
1:A:156:TYR:CE2	1:A:160:LEU:HD21	2.53	0.42
5:E:120:ASN:O	5:E:122:THR:HG23	2.19	0.42
4:D:113:ASN:N	4:D:114:PRO:HD3	2.35	0.42
1:A:15:PRO:HB3	1:A:90:GLY:O	2.20	0.41
4:D:152:PHE:O	4:D:173:ALA:HA	2.20	0.41
4:D:86:LEU:HD12	4:D:86:LEU:HA	1.83	0.41
3:C:7:VAL:HA	5:E:101:TYR:CE1	2.56	0.41
1:A:21:ARG:CZ	1:A:23:ILE:HD11	2.51	0.41
4:D:12:ILE:HA	4:D:106:THR:O	2.21	0.40
1:A:156:TYR:CE2	1:A:160:LEU:HD11	2.56	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	275/281 (98%)	267 (97%)	8 (3%)	0	100	100
2	B	97/100 (97%)	96 (99%)	1 (1%)	0	100	100
3	C	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
4	D	175/191 (92%)	172 (98%)	3 (2%)	0	100	100
5	E	236/263 (90%)	228 (97%)	8 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	790/844 (94%)	769 (97%)	21 (3%)	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	231/239 (97%)	229 (99%)	2 (1%)	78	92
2	B	93/95 (98%)	90 (97%)	3 (3%)	39	65
3	C	8/8 (100%)	8 (100%)	0	100	100
4	D	149/171 (87%)	142 (95%)	7 (5%)	26	49
5	E	195/228 (86%)	193 (99%)	2 (1%)	76	90
All	All	676/741 (91%)	662 (98%)	14 (2%)	53	78

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	225	THR
1	A	255	GLN
2	B	2	GLN
2	B	70	PHE
2	B	88	SER
4	D	13	VAL
4	D	47	LEU
4	D	101	SER
4	D	129	SER
4	D	141	ILE
4	D	172	ILE
4	D	175	SER
5	E	25	GLU
5	E	132	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	ASN
1	A	220	ASN
1	A	226	GLN
2	B	2	GLN
5	E	210	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

17 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	$\# Z > 2$
7	SO4	E	304	-	4,4,4	0.39	0	6,6,6	0.05	0
6	GOL	B	101	-	5,5,5	0.12	0	5,5,5	0.33	0
6	GOL	E	302	-	5,5,5	0.09	0	5,5,5	0.31	0
6	GOL	E	303	-	5,5,5	0.11	0	5,5,5	0.30	0
7	SO4	A	307	-	4,4,4	0.33	0	6,6,6	0.10	0
6	GOL	A	305	-	5,5,5	0.13	0	5,5,5	0.30	0
6	GOL	B	102	-	5,5,5	0.08	0	5,5,5	0.32	0
7	SO4	D	202	-	4,4,4	0.89	0	6,6,6	0.10	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	GOL	C	101	-	5,5,5	0.11	0	5,5,5	0.13	0
6	GOL	A	301	-	5,5,5	0.14	0	5,5,5	0.45	0
6	GOL	A	302	-	5,5,5	0.14	0	5,5,5	0.32	0
6	GOL	D	201	-	5,5,5	0.17	0	5,5,5	0.48	0
6	GOL	E	301	-	5,5,5	0.16	0	5,5,5	0.51	0
6	GOL	A	306	-	5,5,5	0.09	0	5,5,5	0.33	0
6	GOL	A	303	-	5,5,5	0.09	0	5,5,5	0.29	0
6	GOL	A	304	-	5,5,5	0.11	0	5,5,5	0.38	0
7	SO4	E	305	-	4,4,4	0.34	0	6,6,6	0.13	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	B	101	-	-	2/4/4/4	-
6	GOL	E	302	-	-	0/4/4/4	-
6	GOL	E	303	-	-	0/4/4/4	-
6	GOL	A	305	-	-	0/4/4/4	-
6	GOL	B	102	-	-	0/4/4/4	-
6	GOL	C	101	-	-	1/4/4/4	-
6	GOL	A	301	-	-	3/4/4/4	-
6	GOL	A	302	-	-	2/4/4/4	-
6	GOL	D	201	-	-	4/4/4/4	-
6	GOL	E	301	-	-	2/4/4/4	-
6	GOL	A	306	-	-	0/4/4/4	-
6	GOL	A	303	-	-	0/4/4/4	-
6	GOL	A	304	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	301	GOL	C1-C2-C3-O3
6	A	302	GOL	C1-C2-C3-O3
6	A	302	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
6	A	304	GOL	C1-C2-C3-O3
6	D	201	GOL	O1-C1-C2-C3
6	D	201	GOL	C1-C2-C3-O3
6	D	201	GOL	O1-C1-C2-O2
6	B	101	GOL	C1-C2-C3-O3
6	E	301	GOL	O1-C1-C2-C3
6	A	304	GOL	O2-C2-C3-O3
6	B	101	GOL	O2-C2-C3-O3
6	D	201	GOL	O2-C2-C3-O3
6	A	301	GOL	O2-C2-C3-O3
6	E	301	GOL	O1-C1-C2-O2
6	C	101	GOL	O1-C1-C2-C3
6	A	301	GOL	O1-C1-C2-C3

There are no ring outliers.

No monomer is involved in short contacts.

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	278/281 (98%)	-0.14	3 (1%) 80 82	32, 47, 93, 153	3 (1%)
2	B	99/100 (99%)	-0.05	1 (1%) 82 84	34, 48, 78, 114	0
3	C	9/9 (100%)	-0.21	0 100 100	33, 37, 40, 41	0
4	D	180/191 (94%)	0.31	14 (7%) 13 13	32, 52, 121, 150	0
5	E	238/263 (90%)	0.08	10 (4%) 36 39	31, 59, 93, 126	0
All	All	804/844 (95%)	0.04	28 (3%) 44 47	31, 50, 101, 153	3 (0%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	163	ALA	6.1
5	E	241	ASP	5.4
1	A	279	SER	5.3
4	D	151	THR	4.9
4	D	149	SER	4.8
4	D	174	TRP	3.9
2	B	1	ILE	3.9
1	A	278	PRO	3.3
4	D	184	ASP	3.2
5	E	183	TYR	3.2
5	E	182	ASN	3.2
4	D	181	THR	3.1
1	A	277	PRO	3.1
4	D	189	THR	3.0
4	D	175	SER	3.0
5	E	140	GLN	2.8
4	D	118	VAL	2.7
5	E	236	ALA	2.7
4	D	182	CYS	2.7
5	E	205	PHE	2.6

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Mol	Chain	Res	Type	RSRZ
4	D	131	LEU	2.6
5	E	136	ILE	2.5
5	E	199	HIS	2.3
5	E	137	ALA	2.3
5	E	198	TRP	2.3
4	D	150	GLY	2.1
4	D	119	TYR	2.0
4	D	124	PRO	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 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(Å ²)	Q<0.9
6	GOL	A	306	6/6	0.80	0.29	82,98,110,110	0
6	GOL	B	102	6/6	0.85	0.16	79,90,91,92	0
6	GOL	D	201	6/6	0.85	0.17	72,82,83,84	0
7	SO4	A	307	5/5	0.87	0.23	82,86,107,107	0
6	GOL	E	302	6/6	0.88	0.14	70,79,89,89	0
6	GOL	E	301	6/6	0.88	0.27	58,60,65,68	0
6	GOL	E	303	6/6	0.89	0.27	72,90,95,96	0
6	GOL	A	305	6/6	0.90	0.22	56,74,76,81	0
6	GOL	A	304	6/6	0.91	0.20	72,75,75,76	0
6	GOL	A	303	6/6	0.92	0.17	70,81,88,92	0
6	GOL	B	101	6/6	0.92	0.36	51,67,72,78	0
7	SO4	E	304	5/5	0.92	0.16	71,81,92,101	0
6	GOL	A	302	6/6	0.93	0.15	57,79,80,83	0
6	GOL	A	301	6/6	0.95	0.25	49,51,60,60	0
7	SO4	D	202	5/5	0.97	0.09	56,66,75,89	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
6	GOL	C	101	6/6	0.97	0.19	50,52,55,64	0
7	SO4	E	305	5/5	0.97	0.09	69,79,84,93	0

6.5 Other polymers [i](#)

There are no such residues in this entry.