



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 04:04 AM UTC

PDB ID : 3QFJ / pdb_00003qfj
Title : The complex between TCR A6 and human Class I MHC HLA-A2 with the modified TAX (Y5F) peptide
Authors : Borbulevych, O.Y.; Baker, B.M.
Deposited on : 2011-01-21
Resolution : 2.29 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

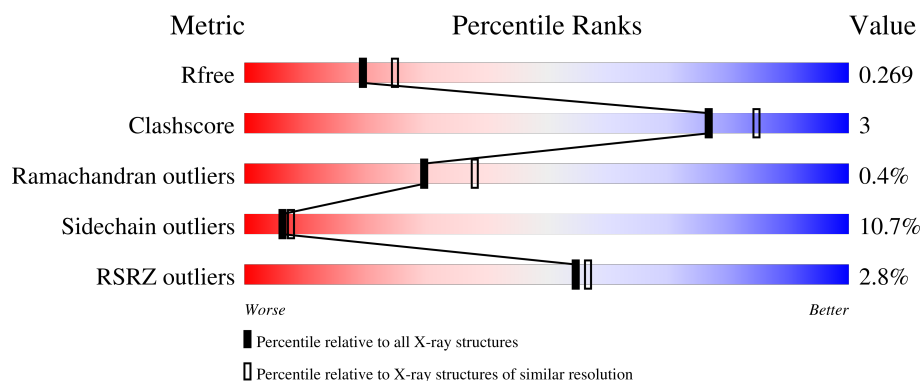
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.29 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	275	4% 83% 16% .
2	B	100	2% 86% 12% .
3	C	9	89% 11%
4	D	200	4% 81% 17% .
5	E	245	% 81% 16% .

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6898 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 HLA class I histocompatibility antigen, A-2 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	0	0
			2247	1403	409	426	9			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			

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 TAX(Y5F) peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			76	56	9	11			

- Molecule 4 is a protein called A6 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	200	Total	C	N	O	S	0	0	0
			1552	965	255	325	7			

- Molecule 5 is a protein called A6 beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	245	Total	C	N	O	S	0	0	0
			1927	1209	338	372	8			

- Molecule 6 is GLYCEROL (CCD ID: 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	B	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		
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		

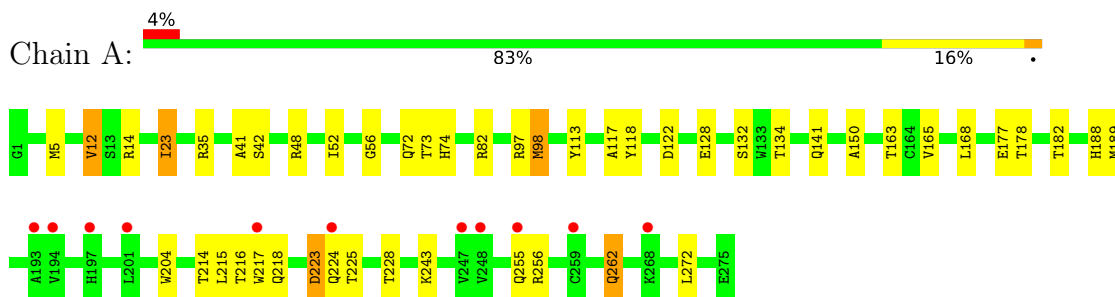
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	64	Total 64	O 64	0	0
7	B	31	Total 31	O 31	0	0
7	C	3	Total 3	O 3	0	0
7	D	34	Total 34	O 34	0	0
7	E	61	Total 61	O 61	0	0

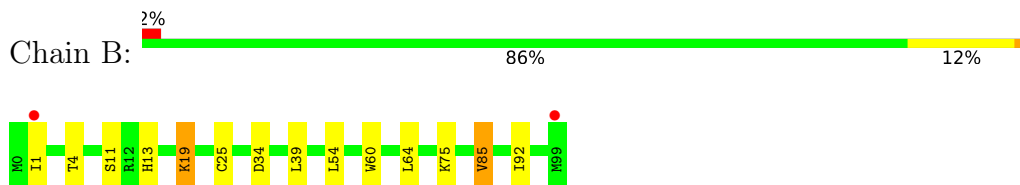
3 Residue-property plots

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.

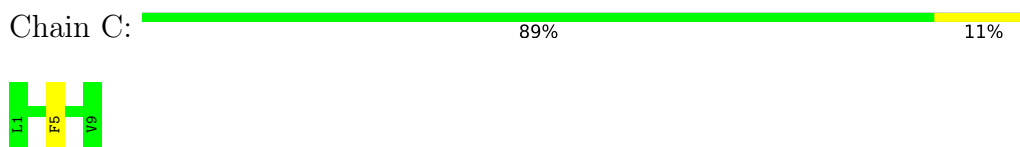
- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain



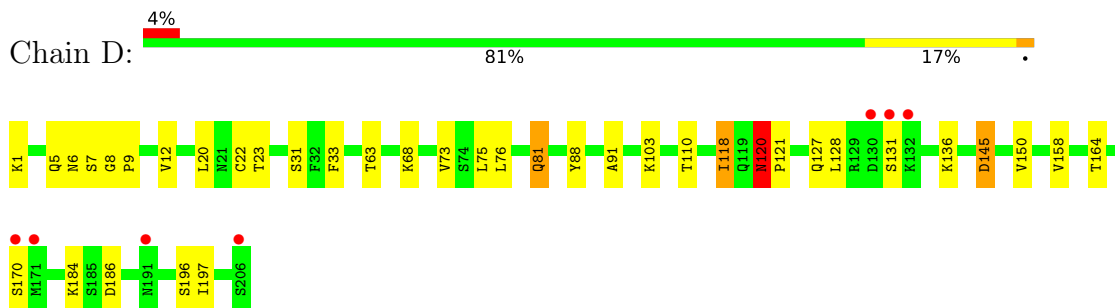
- Molecule 2: Beta-2-microglobulin



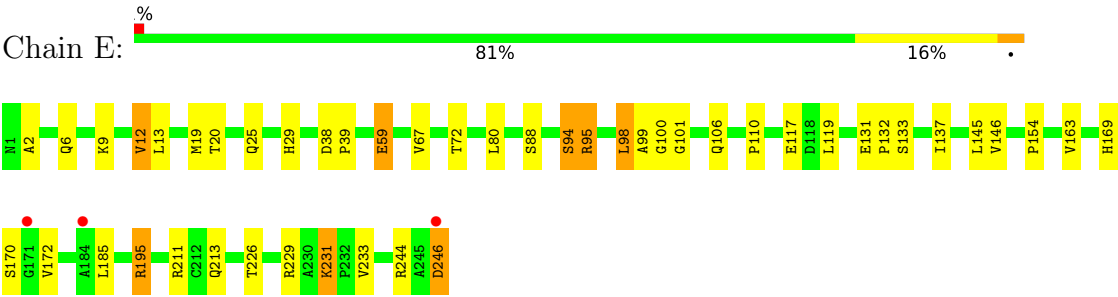
- Molecule 3: TAX(Y5F) peptide



- Molecule 4: A6 alpha chain



- Molecule 5: A6 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	224.01Å 48.10Å 92.96Å 90.00° 91.02° 90.00°	Depositor
Resolution (Å)	20.00 – 2.29 20.00 – 2.29	Depositor EDS
% Data completeness (in resolution range)	98.8 (20.00-2.29) 98.6 (20.00-2.29)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.41 (at 2.30Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.218 , 0.271 0.218 , 0.269	Depositor DCC
R_{free} test set	2244 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	45.4	Xtriage
Anisotropy	0.282	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 53.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.016 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6898	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.72% 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

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	0.73	3/2312 (0.1%)	1.06	8/3137 (0.3%)
2	B	0.67	0/860	0.98	1/1162 (0.1%)
3	C	0.92	0/79	1.29	1/106 (0.9%)
4	D	0.63	0/1585	0.98	5/2150 (0.2%)
5	E	0.79	0/1980	1.13	10/2699 (0.4%)
All	All	0.72	3/6816 (0.0%)	1.06	25/9254 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
4	D	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	12	VAL	CA-CB	5.45	1.61	1.54
1	A	117	ALA	C-O	5.33	1.30	1.23
1	A	118	TYR	C-O	5.12	1.29	1.24

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	99	ALA	CA-C-N	-10.36	114.18	122.16
5	E	99	ALA	C-N-CA	-10.36	114.18	122.16
5	E	38	ASP	CA-C-N	8.86	130.91	119.84
5	E	38	ASP	C-N-CA	8.86	130.91	119.84
5	E	100	GLY	N-CA-C	7.62	124.57	111.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	184	LYS	N-CA-C	7.21	121.72	112.34
1	A	98	MET	CG-SD-CE	-6.57	86.44	100.90
1	A	42	SER	N-CA-C	6.30	118.14	111.28
5	E	169	HIS	N-CA-C	-5.98	104.44	113.89
1	A	141	GLN	N-CA-C	-5.96	104.72	112.23
4	D	81	GLN	N-CA-C	-5.74	101.78	110.50
4	D	118	ILE	CB-CA-C	5.67	118.36	110.99
4	D	120	ASN	CA-C-N	5.67	125.62	119.78
4	D	120	ASN	C-N-CA	5.67	125.62	119.78
2	B	85	VAL	CB-CA-C	-5.58	101.13	111.79
5	E	131	GLU	CA-C-N	5.23	125.23	119.90
5	E	131	GLU	C-N-CA	5.23	125.23	119.90
1	A	41	ALA	CA-C-N	5.22	127.28	120.28
1	A	41	ALA	C-N-CA	5.22	127.28	120.28
1	A	56	GLY	CA-C-N	5.18	124.98	119.28
1	A	56	GLY	C-N-CA	5.18	124.98	119.28
5	E	231	LYS	CA-C-N	5.13	126.25	119.84
5	E	231	LYS	C-N-CA	5.13	126.25	119.84
1	A	74	HIS	N-CA-C	5.08	117.48	111.33
3	C	5	PHE	N-CA-C	5.04	117.85	109.79

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	D	131	SER	Peptide

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	2247	0	2096	18	0
2	B	837	0	803	5	0
3	C	76	0	79	0	0
4	D	1552	0	1461	9	0
5	E	1927	0	1830	12	0
6	A	12	0	16	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	18	0	24	0	0
6	C	6	0	8	0	0
6	D	6	0	8	0	0
6	E	24	0	32	0	0
7	A	64	0	0	1	0
7	B	31	0	0	0	0
7	C	3	0	0	0	0
7	D	34	0	0	0	0
7	E	61	0	0	0	0
All	All	6898	0	6357	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:164:THR:OG1	5:E:195:ARG:NH2	2.32	0.63
5:E:95:ARG:HG2	5:E:106:GLN:HB2	1.82	0.62
1:A:14:ARG:HA	6:A:277:GOL:H11	1.85	0.59
4:D:5:GLN:NE2	4:D:88:TYR:O	2.38	0.55
1:A:224:GLN:O	1:A:228:THR:OG1	2.24	0.54
1:A:14:ARG:HG2	6:A:277:GOL:H31	1.90	0.54
1:A:98:MET:HE1	1:A:113:TYR:CE1	2.43	0.53
1:A:223:ASP:OD1	1:A:223:ASP:N	2.34	0.52
1:A:214:THR:HB	1:A:262:GLN:HG3	1.91	0.52
1:A:73:THR:HG22	5:E:98:LEU:HG	1.92	0.52
4:D:33:PHE:HB2	4:D:91:ALA:HB3	1.93	0.51
5:E:231:LYS:HG2	5:E:233:VAL:HG13	1.92	0.50
2:B:11:SER:OG	2:B:13:HIS:O	2.27	0.50
1:A:98:MET:HE1	1:A:113:TYR:HE1	1.78	0.49
1:A:97:ARG:NH2	7:A:295:HOH:O	2.45	0.49
5:E:6:GLN:HB3	5:E:110:PRO:HD2	1.95	0.49
5:E:132:PRO:HD3	5:E:145:LEU:HG	1.97	0.47
1:A:122:ASP:OD2	2:B:60:TRP:NE1	2.42	0.47
4:D:118:ILE:HD11	4:D:145:ASP:HA	1.94	0.47
4:D:120:ASN:HA	4:D:121:PRO:HD2	1.74	0.46
4:D:63:THR:HB	4:D:76:LEU:HB2	1.98	0.46
5:E:12:VAL:HG13	5:E:154:PRO:HG3	1.97	0.46
1:A:188:HIS:HB2	1:A:204:TRP:HB2	1.98	0.46
1:A:189:MET:HE1	1:A:217:TRP:HH2	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:211:ARG:NH2	5:E:213:GLN:OE1	2.51	0.44
1:A:48:ARG:HD3	1:A:48:ARG:HA	1.74	0.44
2:B:19:LYS:HE3	2:B:19:LYS:HB2	1.78	0.43
1:A:163:THR:OG1	4:D:68:LYS:NZ	2.52	0.43
2:B:25:CYS:HB2	2:B:39:LEU:HD21	2.00	0.43
1:A:215:LEU:HD12	1:A:243:LYS:HD3	2.00	0.43
1:A:5:MET:HB2	1:A:168:LEU:HD13	2.00	0.42
4:D:8:GLY:HA2	4:D:9:PRO:HA	1.65	0.42
1:A:150:ALA:O	5:E:101:GLY:HA2	2.19	0.42
5:E:59:GLU:H	5:E:59:GLU:HG3	1.56	0.42
5:E:244:ARG:NH1	5:E:246:ASP:OD2	2.49	0.42
4:D:20:LEU:HD12	4:D:75:LEU:HD23	2.02	0.41
5:E:29:HIS:ND1	5:E:94:SER:OG	2.33	0.40
1:A:23:ILE:HD12	2:B:54:LEU:HB3	2.04	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	273/275 (99%)	261 (96%)	12 (4%)	0	100	100
2	B	98/100 (98%)	96 (98%)	2 (2%)	0	100	100
3	C	7/9 (78%)	7 (100%)	0	0	100	100
4	D	198/200 (99%)	186 (94%)	11 (6%)	1 (0%)	24	31
5	E	243/245 (99%)	229 (94%)	12 (5%)	2 (1%)	16	20
All	All	819/829 (99%)	779 (95%)	37 (4%)	3 (0%)	30	38

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
5	E	2	ALA
4	D	170	SER
5	E	39	PRO

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	231/231 (100%)	210 (91%)	21 (9%)	9	11
2	B	95/95 (100%)	87 (92%)	8 (8%)	10	14
3	C	8/8 (100%)	8 (100%)	0	100	100
4	D	178/178 (100%)	157 (88%)	21 (12%)	5	6
5	E	209/209 (100%)	182 (87%)	27 (13%)	4	4
All	All	721/721 (100%)	644 (89%)	77 (11%)	6	8

All (77) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	VAL
1	A	23	ILE
1	A	35	ARG
1	A	52	ILE
1	A	72	GLN
1	A	82	ARG
1	A	128	GLU
1	A	132	SER
1	A	134	THR
1	A	165	VAL
1	A	177	GLU
1	A	178	THR
1	A	182	THR
1	A	216	THR
1	A	218	GLN
1	A	223	ASP
1	A	225	THR

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Mol	Chain	Res	Type
1	A	255	GLN
1	A	256	ARG
1	A	262	GLN
1	A	272	LEU
2	B	1	ILE
2	B	4	THR
2	B	19	LYS
2	B	34	ASP
2	B	64	LEU
2	B	75	LYS
2	B	85	VAL
2	B	92	ILE
4	D	1	LYS
4	D	6	ASN
4	D	7	SER
4	D	12	VAL
4	D	22	CYS
4	D	23	THR
4	D	31	SER
4	D	73	VAL
4	D	81	GLN
4	D	103	LYS
4	D	110	THR
4	D	120	ASN
4	D	127	GLN
4	D	128	LEU
4	D	136	LYS
4	D	145	ASP
4	D	150	VAL
4	D	158	VAL
4	D	186	ASP
4	D	196	SER
4	D	197	ILE
5	E	9	LYS
5	E	12	VAL
5	E	13	LEU
5	E	19	MET
5	E	20	THR
5	E	25	GLN
5	E	59	GLU
5	E	67	VAL
5	E	72	THR

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Mol	Chain	Res	Type
5	E	80	LEU
5	E	88	SER
5	E	94	SER
5	E	95	ARG
5	E	98	LEU
5	E	117	GLU
5	E	119	LEU
5	E	133	SER
5	E	137	ILE
5	E	146	VAL
5	E	163	VAL
5	E	170	SER
5	E	172	VAL
5	E	185	LEU
5	E	195	ARG
5	E	226	THR
5	E	229	ARG
5	E	246	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	141	GLN
1	A	151	HIS
2	B	2	GLN
2	B	13	HIS
4	D	71	GLN
4	D	81	GLN
4	D	105	GLN
4	D	111	GLN
4	D	127	GLN
5	E	11	GLN
5	E	66	ASN
5	E	186	ASN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

11 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	B	100	-	5,5,5	0.35	0	5,5,5	0.49	0
6	GOL	E	247	-	5,5,5	0.55	0	5,5,5	0.60	0
6	GOL	E	250	-	5,5,5	0.43	0	5,5,5	0.55	0
6	GOL	B	102	-	5,5,5	0.31	0	5,5,5	0.68	0
6	GOL	A	277	-	5,5,5	0.20	0	5,5,5	0.77	0
6	GOL	E	248	-	5,5,5	0.41	0	5,5,5	0.42	0
6	GOL	D	207	-	5,5,5	0.38	0	5,5,5	0.56	0
6	GOL	A	276	-	5,5,5	0.36	0	5,5,5	0.30	0
6	GOL	C	10	-	5,5,5	0.44	0	5,5,5	0.47	0
6	GOL	E	249	-	5,5,5	0.79	0	5,5,5	0.70	0
6	GOL	B	101	-	5,5,5	0.56	0	5,5,5	0.96	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	100	-	-	4/4/4/4	-
6	GOL	E	247	-	-	4/4/4/4	-
6	GOL	E	250	-	-	4/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	GOL	B	102	-	-	2/4/4/4	-
6	GOL	A	277	-	-	0/4/4/4	-
6	GOL	E	248	-	-	4/4/4/4	-
6	GOL	D	207	-	-	4/4/4/4	-
6	GOL	A	276	-	-	1/4/4/4	-
6	GOL	C	10	-	-	0/4/4/4	-
6	GOL	E	249	-	-	2/4/4/4	-
6	GOL	B	101	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	100	GOL	O1-C1-C2-C3
6	B	100	GOL	C1-C2-C3-O3
6	B	101	GOL	C1-C2-C3-O3
6	B	102	GOL	C1-C2-C3-O3
6	D	207	GOL	O1-C1-C2-C3
6	E	247	GOL	O1-C1-C2-C3
6	E	247	GOL	O2-C2-C3-O3
6	E	248	GOL	O1-C1-C2-C3
6	E	248	GOL	C1-C2-C3-O3
6	E	249	GOL	O1-C1-C2-C3
6	E	250	GOL	O1-C1-C2-C3
6	E	250	GOL	C1-C2-C3-O3
6	B	100	GOL	O2-C2-C3-O3
6	E	249	GOL	O1-C1-C2-O2
6	E	250	GOL	O2-C2-C3-O3
6	E	247	GOL	C1-C2-C3-O3
6	B	100	GOL	O1-C1-C2-O2
6	B	101	GOL	O2-C2-C3-O3
6	B	102	GOL	O2-C2-C3-O3
6	D	207	GOL	O1-C1-C2-O2
6	D	207	GOL	O2-C2-C3-O3
6	E	247	GOL	O1-C1-C2-O2
6	E	248	GOL	O1-C1-C2-O2
6	E	248	GOL	O2-C2-C3-O3
6	E	250	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
6	A	276	GOL	O2-C2-C3-O3
6	D	207	GOL	C1-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	277	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, 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	275/275 (100%)	0.20	11 (4%) 42 44	23, 51, 155, 190	0
2	B	100/100 (100%)	-0.09	2 (2%) 65 66	23, 45, 86, 98	0
3	C	9/9 (100%)	-0.54	0 100 100	30, 31, 34, 35	0
4	D	200/200 (100%)	0.41	7 (3%) 47 49	31, 58, 110, 134	0
5	E	245/245 (100%)	0.03	3 (1%) 76 77	21, 50, 86, 111	0
All	All	829/829 (100%)	0.16	23 (2%) 55 57	21, 52, 122, 190	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	171	MET	4.5
4	D	206	SER	3.5
1	A	197	HIS	3.2
4	D	130	ASP	3.1
2	B	99	MET	3.0
1	A	193	ALA	3.0
2	B	1	ILE	3.0
4	D	131	SER	2.8
5	E	246	ASP	2.5
1	A	194	VAL	2.5
1	A	248	VAL	2.4
1	A	201	LEU	2.4
1	A	259	CYS	2.4
4	D	132	LYS	2.3
1	A	217	TRP	2.3
4	D	170	SER	2.3
1	A	255	GLN	2.2
4	D	191	ASN	2.2
5	E	184	ALA	2.2
5	E	171	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	224	GLN	2.1
1	A	247	VAL	2.0
1	A	268	LYS	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(Å ²)	Q<0.9
6	GOL	A	277	6/6	0.75	0.15	47,48,49,50	0
6	GOL	D	207	6/6	0.75	0.18	48,48,49,50	0
6	GOL	A	276	6/6	0.77	0.11	72,72,72,73	0
6	GOL	E	250	6/6	0.78	0.19	52,53,54,55	0
6	GOL	B	100	6/6	0.79	0.10	76,76,76,76	0
6	GOL	B	101	6/6	0.80	0.13	47,47,48,48	0
6	GOL	E	249	6/6	0.81	0.19	41,43,44,44	0
6	GOL	C	10	6/6	0.82	0.12	47,47,48,49	0
6	GOL	B	102	6/6	0.82	0.12	46,47,48,49	0
6	GOL	E	248	6/6	0.89	0.09	58,59,59,59	0
6	GOL	E	247	6/6	0.91	0.09	53,53,53,54	6

6.5 Other polymers [i](#)

There are no such residues in this entry.