



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:10 PM UTC

PDB ID : 2X70 / pdb_00002x70
Title : Crystal structure of MHC CClass I HLA-A2.1 bound to a photocleavable peptide
Authors : Celie, P.H.N.; Toebe, M.; Rodenko, B.; Ovaa, H.; Perrakis, A.; Schumacher, T.N.M.
Deposited on : 2010-02-22
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

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:

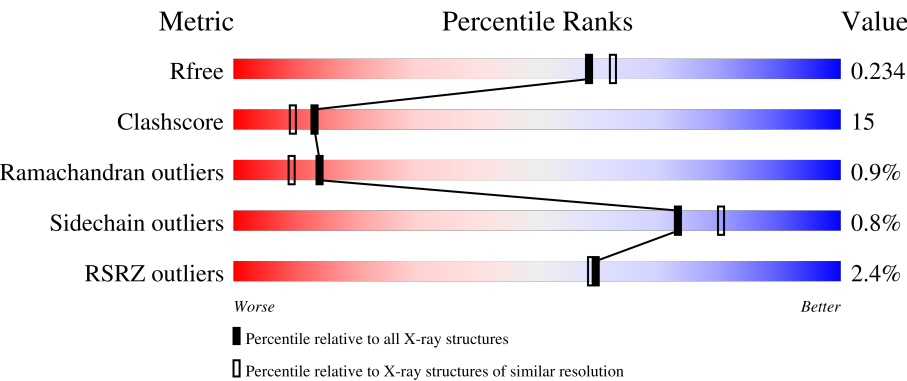
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (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 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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 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	
1	D	275	
2	B	100	
2	E	100	
3	C	9	

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Mol	Chain	Length	Quality of chain
3	F	9	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	GOL	D	1278[A]	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7337 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-2.1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	7	0
			2306	1441	422	433	10			
1	D	275	Total	C	N	O	S	0	13	0
			2319	1452	420	437	10			

- 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			
2	E	100	Total	C	N	O	S	0	3	0
			850	542	141	162	5			

There are 2 discrepancies between the modelled and reference sequences:

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

- Molecule 3 is a protein called HLA-A2.1-RESTRICTED INFLUENZA A MATRIX EPI-TOPE.

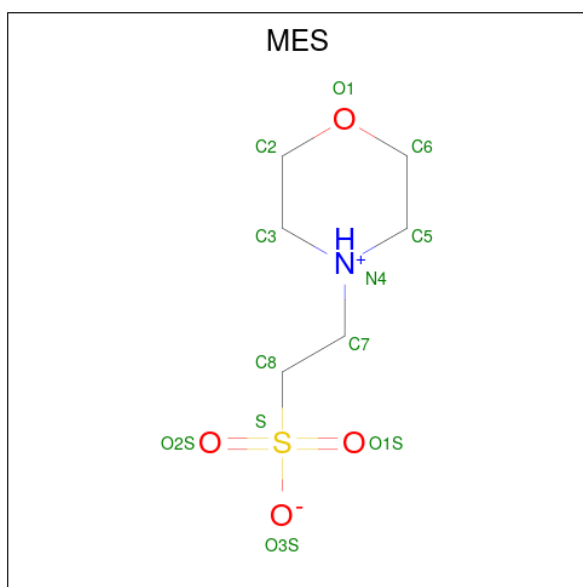
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			82	56	12	14			
3	F	9	Total	C	N	O	0	0	0
			82	56	12	14			

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	1
			12	6	6		
4	D	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	1
			12	6	6		
4	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
5	D	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
5	D	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
5	E	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

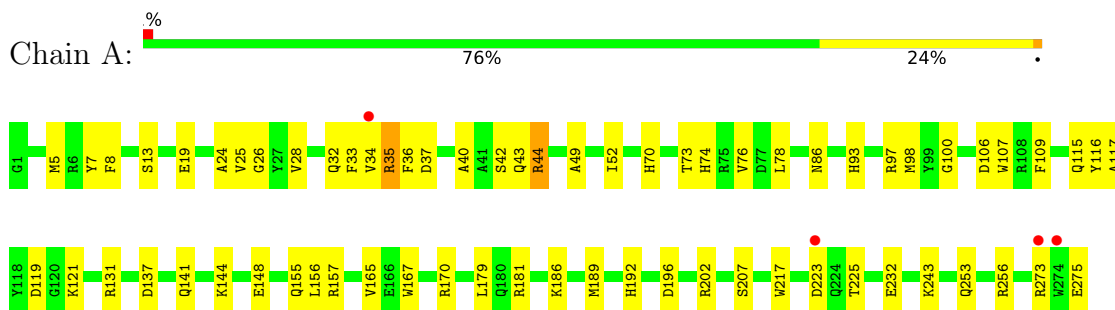
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	230	Total	O	0	0
			230	230		
6	B	98	Total	O	0	0
			98	98		
6	C	7	Total	O	0	0
			7	7		
6	D	302	Total	O	0	0
			302	302		
6	E	125	Total	O	0	0
			125	125		
6	F	9	Total	O	0	0
			9	9		

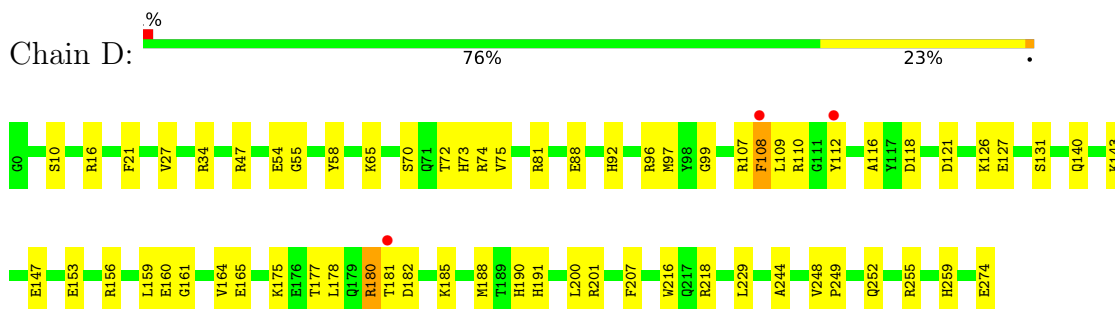
3 Residue-property plots [i](#)

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.

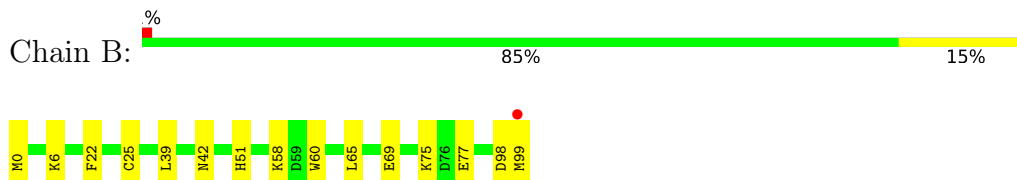
- Molecule 1: HLA CLASS I HISTOCOMPATIBILITY ANTIGEN, A-2.1



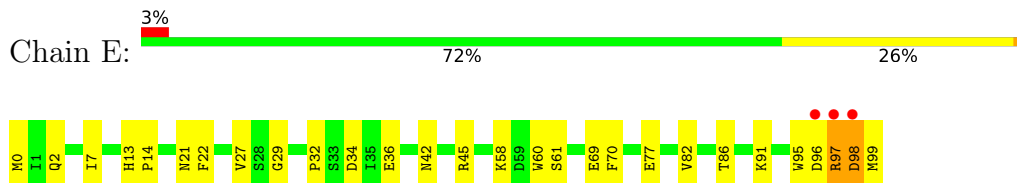
- Molecule 1: HLA CLASS I HISTOCOMPATIBILITY ANTIGEN, A-2.1



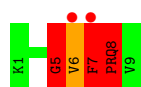
- Molecule 2: BETA-2-MICROGLOBULIN



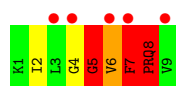
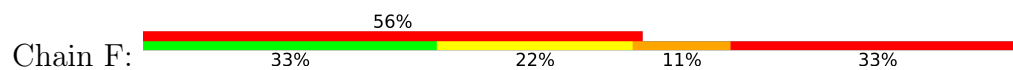
- Molecule 2: BETA-2-MICROGLOBULIN



- Molecule 3: HLA-A2.1-RESTRICTED INFLUENZA A MATRIX EPITOPE



● Molecule 3: HLA-A2.1-RESTRICTED INFLUENZA A MATRIX EPITOPE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.98Å 81.80Å 79.71Å 90.00° 90.77° 90.00°	Depositor
Resolution (Å)	19.93 – 2.00 19.93 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.1 (19.93-2.00) 97.0 (19.93-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.10 (at 2.01Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.159 , 0.225 0.177 , 0.234	Depositor DCC
R_{free} test set	2662 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	25.5	Xtriage
Anisotropy	0.139	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 60.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.020 for -h,-l,-k 0.009 for -h,l,k 0.054 for h,-k,-l	Xtriage
Reported twinning fraction	0.043 for -h,-k,l	Depositor
Outliers	0 of 52222 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7337	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.33% 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: PRQ, GOL, PRV, MES

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	0.33	1/2384 (0.0%)	0.67	0/3231
1	D	0.31	0/2421	0.69	2/3281 (0.1%)
2	B	0.30	0/860	0.67	0/1162
2	E	0.34	0/882	0.64	0/1191
3	C	0.27	0/53	0.68	0/65
3	F	0.26	0/53	0.57	0/65
All	All	0.32	1/6653 (0.0%)	0.67	2/8995 (0.0%)

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
3	C	0	5
3	F	0	6
All	All	0	11

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	121	LYS	C-N	-5.85	1.25	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	55	GLY	CA-C-N	5.09	124.88	119.28
1	D	55	GLY	C-N-CA	5.09	124.88	119.28

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	5	PRV	Mainchain,Peptide
3	C	7	PHE	Mainchain
3	C	8	PRQ	Mainchain,Peptide
3	F	4	GLY	Mainchain,Peptide
3	F	5	PRV	Mainchain,Peptide
3	F	7	PHE	Mainchain
3	F	8	PRQ	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	2306	0	2164	70	0
1	D	2319	0	2196	66	0
2	B	837	0	803	13	0
2	E	850	0	823	36	0
3	C	82	0	76	12	0
3	F	82	0	77	20	0
4	A	18	0	24	2	0
4	D	24	0	32	6	0
5	B	12	0	12	0	0
5	D	24	0	24	0	0
5	E	12	0	12	0	0
6	A	230	0	0	10	0
6	B	98	0	0	1	0
6	C	7	0	0	0	0
6	D	302	0	0	14	0
6	E	125	0	0	5	0
6	F	9	0	0	1	0
All	All	7337	0	6243	190	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 15.

All (190) 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:D:96[B]:ARG:NH2	3:F:7:PHE:CE1	2.10	1.18
1:D:191:HIS:HE1	2:E:98:ASP:OD1	1.30	1.14
2:E:96:ASP:HB3	2:E:99:MET:HE3	1.12	1.12
4:A:1277[B]:GOL:H32	2:B:58:LYS:HA	1.33	1.10
2:E:96:ASP:CB	2:E:99:MET:HE3	1.83	1.08
1:A:26:GLY:HA3	1:A:34:VAL:CG2	1.87	1.04
3:F:6:VAL:HG13	3:F:7:PHE:H	1.21	1.04
1:D:96[B]:ARG:NH2	3:F:7:PHE:CD1	2.26	1.03
4:D:1278[A]:GOL:H11	2:E:58:LYS:HA	1.42	1.00
1:A:44[A]:ARG:HH11	1:A:44[A]:ARG:HG3	1.26	0.99
1:A:26:GLY:N	1:A:34:VAL:HG23	1.77	0.98
1:D:191:HIS:CE1	2:E:98:ASP:OD1	2.18	0.95
3:F:6:VAL:O	3:F:7:PHE:HB2	1.72	0.89
1:D:96[B]:ARG:HH21	3:F:7:PHE:HE1	1.20	0.87
3:C:6:VAL:HG13	3:C:7:PHE:H	1.40	0.86
1:A:26:GLY:CA	1:A:34:VAL:HG23	2.04	0.85
1:A:223:ASP:CG	1:A:225:THR:HG23	2.01	0.84
3:F:6:VAL:HG13	3:F:7:PHE:N	1.94	0.82
1:A:26:GLY:CA	1:A:34:VAL:CG2	2.57	0.82
1:A:44[A]:ARG:HH11	1:A:44[A]:ARG:CG	1.93	0.81
1:A:35:ARG:HD2	1:A:35:ARG:C	2.11	0.75
1:A:26:GLY:H	1:A:34:VAL:HG23	1.49	0.74
2:E:96:ASP:HB3	2:E:99:MET:CE	2.07	0.74
1:A:33:PHE:CD2	1:A:34:VAL:HG13	2.24	0.72
1:D:110:ARG:NH1	6:D:2146:HOH:O	2.23	0.72
2:E:97:ARG:HG3	6:E:2119:HOH:O	1.88	0.72
1:A:223:ASP:OD2	1:A:225:THR:HG23	1.89	0.71
1:A:97[B]:ARG:CZ	1:A:116[B]:TYR:HE1	2.04	0.71
1:A:189:MET:HE3	1:A:217:TRP:HH2	1.57	0.69
2:B:98:ASP:HA	2:B:99:MET:OXT	1.91	0.69
1:A:167:TRP:CZ3	1:A:170[A]:ARG:HD2	2.27	0.69
1:D:191:HIS:HD2	6:D:2214:HOH:O	1.76	0.68
1:A:156:LEU:HD21	3:C:7:PHE:HZ	1.57	0.67
3:C:6:VAL:O	3:C:7:PHE:HB2	1.93	0.67
1:D:96[B]:ARG:HH22	3:F:7:PHE:HD1	1.35	0.67
1:D:177:THR:O	1:D:180:ARG:HD3	1.93	0.67
1:D:108:PHE:HD1	1:D:164:VAL:CG1	2.09	0.66
2:E:95:TRP:CH2	2:E:97:ARG:HG2	2.30	0.66
1:D:108:PHE:HD1	1:D:164:VAL:HG11	1.60	0.65
1:A:24:ALA:O	1:A:35:ARG:HA	1.97	0.65
1:A:35:ARG:NH2	6:A:2039:HOH:O	2.30	0.65
1:D:252:GLN:NE2	1:D:255:ARG:HH11	1.96	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:THR:HG23	3:C:8:PRQ:HAH	1.79	0.64
1:A:33:PHE:CE2	1:A:34:VAL:HG13	2.33	0.64
1:D:116:ALA:HB2	2:E:60:TRP:CE2	2.32	0.64
1:D:188[A]:MET:HE3	1:D:216:TRP:HH2	1.63	0.64
2:E:97:ARG:O	2:E:99:MET:N	2.30	0.64
1:D:191:HIS:HE1	2:E:98:ASP:CG	2.05	0.63
1:A:93:HIS:HD2	1:A:119:ASP:OD2	1.81	0.63
1:A:170[B]:ARG:HD2	6:A:2148:HOH:O	1.98	0.63
2:B:98:ASP:HA	2:B:99:MET:C	2.25	0.62
1:D:72:THR:HA	3:F:8:PRQ:CAF	2.29	0.61
2:E:77:GLU:HG2	6:E:2100:HOH:O	2.01	0.60
3:F:5:PRV:HA	3:F:5:PRV:O1	2.01	0.60
1:A:223:ASP:OD2	1:A:225:THR:CG2	2.49	0.60
1:D:107:ARG:O	1:D:108:PHE:C	2.45	0.59
3:F:5:PRV:O	6:F:2009:HOH:O	2.17	0.59
1:A:232[A]:GLU:CD	2:B:6:LYS:HE3	2.28	0.59
1:D:191:HIS:CE1	2:E:98:ASP:CG	2.79	0.59
1:A:223:ASP:OD1	1:A:225:THR:HG23	2.02	0.59
2:E:42:ASN:ND2	2:E:77:GLU:H	2.00	0.59
1:A:97[B]:ARG:CZ	1:A:116[B]:TYR:CE1	2.86	0.58
1:A:7:TYR:HE1	1:A:34:VAL:HG21	1.68	0.58
1:A:192:HIS:HD2	6:A:2171:HOH:O	1.88	0.57
1:D:259:HIS:HB3	6:D:2288:HOH:O	2.04	0.57
1:D:126:LYS:HD2	1:D:131:SER:OG	2.05	0.57
4:D:1278[A]:GOL:H32	2:E:58:LYS:HA	1.87	0.57
1:D:99:GLY:C	1:D:112:TYR:CZ	2.83	0.56
1:A:144:LYS:O	1:A:148:GLU:HG3	2.06	0.55
1:A:156:LEU:HD21	3:C:7:PHE:CZ	2.39	0.55
3:F:6:VAL:CG1	3:F:7:PHE:N	2.66	0.54
1:D:65:LYS:HB2	3:F:2:ILE:HD11	1.90	0.54
2:E:96:ASP:CG	2:E:99:MET:HE3	2.33	0.54
1:D:54:GLU:HG2	1:D:58:TYR:CG	2.43	0.54
1:D:180:ARG:O	1:D:180:ARG:HG2	2.05	0.54
1:A:137:ASP:HB3	6:A:2127:HOH:O	2.07	0.53
1:D:92:HIS:HE1	6:D:2125:HOH:O	1.90	0.53
2:B:25:CYS:HB2	2:B:39:LEU:HD21	1.91	0.53
2:E:22:PHE:CZ	2:E:69:GLU:HG2	2.44	0.53
1:D:161:GLY:O	1:D:165:GLU:HG3	2.09	0.52
1:A:7:TYR:CE1	1:A:34:VAL:HG21	2.44	0.52
1:A:26:GLY:HA3	1:A:34:VAL:HG21	1.83	0.52
2:E:42:ASN:HD21	2:E:77:GLU:H	1.56	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:TRP:CE3	1:A:170[A]:ARG:HD2	2.44	0.52
1:A:40:ALA:HB3	6:A:2040:HOH:O	2.09	0.52
1:A:131:ARG:HE	1:A:157:ARG:NH1	2.08	0.52
1:A:19:GLU:HG3	6:A:2019:HOH:O	2.10	0.52
1:A:74:HIS:CE1	1:A:97[B]:ARG:HD2	2.44	0.52
1:D:185:LYS:HD2	6:D:2230:HOH:O	2.10	0.52
1:A:93:HIS:HE1	6:A:2099:HOH:O	1.93	0.51
4:D:1278[A]:GOL:H11	2:E:58:LYS:CA	2.27	0.51
1:A:253:GLN:NE2	1:A:256:ARG:HH11	2.08	0.51
1:D:73:HIS:CE1	1:D:96[B]:ARG:HD2	2.44	0.51
1:D:99:GLY:CA	1:D:112:TYR:CE1	2.93	0.51
1:D:72:THR:HA	3:F:8:PRQ:HAF	1.93	0.51
1:A:26:GLY:HA3	1:A:34:VAL:HG22	1.89	0.51
1:A:44[A]:ARG:HG3	1:A:44[A]:ARG:NH1	2.05	0.51
1:D:72:THR:HG23	3:F:8:PRQ:CAH	2.40	0.51
1:D:65:LYS:CB	3:F:2:ILE:HD11	2.42	0.50
3:F:5:PRV:CD1	3:F:6:VAL:HB	2.42	0.50
1:A:76:VAL:HG11	3:C:8:PRQ:CAG	2.43	0.49
2:E:34[B]:ASP:HB2	6:E:2055:HOH:O	2.13	0.49
1:D:218:ARG:HA	1:D:255:ARG:O	2.12	0.49
1:D:10[B]:SER:OG	1:D:21:PHE:HD1	1.96	0.49
1:D:109:LEU:O	1:D:110:ARG:HB3	2.13	0.49
3:C:6:VAL:HG22	3:C:7:PHE:N	2.28	0.49
1:D:121:ASP:OD1	4:D:1276:GOL:H12	2.13	0.48
1:A:32:GLN:HE21	1:A:35:ARG:HB3	1.79	0.48
1:A:37:ASP:HB3	1:A:40:ALA:HB2	1.96	0.48
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.48	0.48
3:C:5:PRV:C	3:C:5:PRV:N1	2.77	0.48
1:D:34:ARG:NH1	1:D:47:ARG:CZ	2.77	0.47
1:D:75:VAL:HB	3:F:8:PRQ:HAF	1.96	0.47
1:D:108:PHE:CD1	1:D:164:VAL:HG11	2.46	0.47
2:E:96:ASP:OD2	2:E:99:MET:CE	2.63	0.47
3:F:5:PRV:CE1	3:F:6:VAL:HB	2.44	0.47
1:D:97:MET:HE2	4:D:1278[A]:GOL:H2	1.96	0.47
3:C:6:VAL:HG23	6:D:2074:HOH:O	2.14	0.47
3:C:6:VAL:HG13	3:C:7:PHE:N	2.20	0.46
1:A:5:MET:O	1:A:100:GLY:HA3	2.16	0.46
1:D:27:VAL:HG11	1:D:178:LEU:HD13	1.97	0.46
1:D:92:HIS:HD2	1:D:118:ASP:OD2	1.98	0.46
1:A:273:ARG:HD2	1:A:275:GLU:CD	2.40	0.46
1:D:92:HIS:CD2	2:E:0[B]:MET:HE1	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98[A]:MET:HE2	4:A:1277[A]:GOL:H12	1.97	0.46
2:B:75:LYS:HE3	6:B:2083:HOH:O	2.16	0.45
1:D:181:THR:HG21	6:D:2201:HOH:O	2.16	0.45
1:A:70:HIS:O	1:A:74:HIS:HD2	2.00	0.45
1:A:106:ASP:O	1:A:107:TRP:HB2	2.16	0.45
1:D:140[B]:GLN:OE1	1:D:143:LYS:HD3	2.16	0.45
1:D:274:GLU:HB3	6:D:2299:HOH:O	2.16	0.45
1:A:13:SER:HB3	1:A:78:LEU:HD13	1.98	0.45
1:D:153:GLU:HG3	6:D:2181:HOH:O	2.16	0.45
1:D:81:ARG:NH1	1:D:88:GLU:HG2	2.32	0.45
1:D:108:PHE:CE1	1:D:159:LEU:O	2.70	0.44
1:A:44[A]:ARG:CG	1:A:44[A]:ARG:NH1	2.62	0.44
1:D:156:ARG:NH1	6:D:2183:HOH:O	2.50	0.44
1:A:28:VAL:HG11	1:A:179:LEU:HD13	1.98	0.44
1:D:127[A]:GLU:HG2	6:D:2160:HOH:O	2.16	0.44
1:A:202:ARG:NH1	2:B:99:MET:HA	2.33	0.44
1:D:156:ARG:HD2	6:D:2184:HOH:O	2.16	0.44
1:A:186:LYS:HE2	1:A:207:SER:HB3	2.00	0.44
2:B:0:MET:HE3	2:B:0:MET:HB2	1.90	0.44
2:B:42:ASN:ND2	2:B:77:GLU:H	2.16	0.44
1:A:109:PHE:HD1	1:A:165:VAL:HG11	1.83	0.43
1:D:201:ARG:NH1	2:E:99:MET:OXT	2.48	0.43
1:D:229:LEU:CD1	1:D:244:ALA:HB2	2.48	0.43
2:E:29:GLY:HA2	2:E:61:SER:OG	2.17	0.43
2:E:97:ARG:CG	6:E:2119:HOH:O	2.56	0.43
6:D:2044:HOH:O	2:E:32:PRO:HG3	2.17	0.43
1:A:34:VAL:HG11	6:A:2008:HOH:O	2.18	0.43
1:D:16[B]:ARG:HD2	1:D:16[B]:ARG:HA	1.85	0.43
1:D:99:GLY:C	1:D:112:TYR:CE1	2.97	0.43
2:E:95:TRP:CZ2	2:E:97:ARG:HG2	2.53	0.43
1:A:49:ALA:O	1:A:52:ILE:HG22	2.19	0.42
1:D:160:GLU:O	1:D:164:VAL:HG22	2.18	0.42
2:E:96:ASP:OD2	2:E:99:MET:HE2	2.18	0.42
1:D:143:LYS:O	1:D:147:GLU:HG3	2.18	0.42
1:D:180:ARG:HG3	1:D:182:ASP:OD1	2.20	0.42
2:E:27:VAL:HG23	2:E:27:VAL:O	2.19	0.42
2:E:2:GLN:HB3	2:E:86:THR:HG22	2.02	0.42
1:D:99:GLY:HA3	1:D:112:TYR:CE1	2.55	0.42
1:A:32:GLN:NE2	1:A:35:ARG:HB3	2.35	0.42
2:E:45:ARG:NH2	6:E:2066:HOH:O	2.53	0.41
1:A:137:ASP:O	1:A:141:GLN:HG2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:22:PHE:CE1	2:B:69:GLU:HG2	2.55	0.41
1:A:24:ALA:O	1:A:36:PHE:N	2.50	0.41
1:A:196:ASP:OD1	1:A:196:ASP:C	2.63	0.41
1:D:190:HIS:HD2	1:D:200[B]:LEU:HD21	1.85	0.41
1:D:248:VAL:HA	1:D:249:PRO:HD3	1.98	0.41
3:F:7:PHE:C	3:F:8:PRQ:OAD	2.63	0.41
1:A:97[B]:ARG:NH2	3:C:7:PHE:CD2	2.88	0.41
2:B:51:HIS:HA	2:B:65:LEU:O	2.20	0.41
4:D:1278[B]:GOL:H12	6:D:2149:HOH:O	2.20	0.41
1:A:86:ASN:HB2	6:A:2097:HOH:O	2.20	0.41
1:D:70:SER:O	1:D:74:ARG:HG3	2.20	0.41
2:E:7:ILE:HD12	2:E:91:LYS:HD3	2.03	0.41
2:E:21:ASN:HB3	2:E:70:PHE:CE2	2.56	0.41
1:A:155:GLN:NE2	3:C:5:PRV:HA	2.35	0.41
1:A:243:LYS:HE2	6:A:2177:HOH:O	2.21	0.41
2:E:13:HIS:O	2:E:14:PRO:C	2.62	0.41
1:A:42:SER:O	1:A:43:GLN:HB2	2.21	0.41
1:A:74:HIS:HE1	1:A:97[B]:ARG:HB2	1.86	0.40
1:A:93:HIS:CD2	2:B:0:MET:HE1	2.57	0.40
2:E:95:TRP:CZ3	2:E:97:ARG:HD2	2.57	0.40
1:D:182:ASP:O	1:D:207:PHE:HA	2.21	0.40
1:A:8:PHE:HB2	1:A:25:VAL:HG22	2.02	0.40
2:E:36:GLU:O	2:E:82:VAL:HA	2.21	0.40
1:D:65:LYS:HE3	3:F:2:ILE:HG13	2.03	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	280/275 (102%)	272 (97%)	8 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	286/275 (104%)	278 (97%)	7 (2%)	1 (0%)	36	35
2	B	98/100 (98%)	94 (96%)	4 (4%)	0	100	100
2	E	100/100 (100%)	97 (97%)	1 (1%)	2 (2%)	6	2
3	C	5/9 (56%)	3 (60%)	0	2 (40%)	0	0
3	F	5/9 (56%)	1 (20%)	2 (40%)	2 (40%)	0	0
All	All	774/768 (101%)	745 (96%)	22 (3%)	7 (1%)	14	9

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	98	ASP
3	F	7	PHE
3	C	7	PHE
2	E	97	ARG
1	D	108	PHE
3	F	6	VAL
3	C	6	VAL

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	238/231 (103%)	233 (98%)	5 (2%)	47	52
1	D	244/231 (106%)	243 (100%)	1 (0%)	84	89
2	B	95/95 (100%)	95 (100%)	0	100	100
2	E	98/95 (103%)	98 (100%)	0	100	100
3	C	6/6 (100%)	6 (100%)	0	100	100
3	F	6/6 (100%)	6 (100%)	0	100	100
All	All	687/664 (104%)	681 (99%)	6 (1%)	73	78

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	ARG
1	A	44[A]	ARG
1	A	44[B]	ARG
1	A	115	GLN
1	A	181	ARG
1	D	180	ARG

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	32	GLN
1	A	74	HIS
1	A	87	GLN
1	A	93	HIS
1	A	96	GLN
1	A	155	GLN
1	A	192	HIS
1	A	197	HIS
1	A	224	GLN
1	A	226	GLN
1	A	253	GLN
2	B	42	ASN
1	D	53	GLN
1	D	92	HIS
1	D	95	GLN
1	D	154	GLN
1	D	191	HIS
1	D	223	GLN
1	D	252	GLN
2	E	42	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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
3	PRQ	F	8	3	14,14,15	1.04	1 (7%)	13,18,20	1.28	2 (15%)
3	PRV	C	5	3	11,13,14	1.03	1 (9%)	9,17,19	1.64	1 (11%)
3	PRV	F	5	3	11,13,14	0.97	1 (9%)	9,17,19	1.57	2 (22%)
3	PRQ	C	8	3	14,14,15	1.06	2 (14%)	13,18,20	1.50	4 (30%)

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
3	PRQ	F	8	3	-	5/9/11/12	0/1/1/1
3	PRV	C	5	3	-	4/6/10/12	0/1/1/1
3	PRV	F	5	3	-	3/6/10/12	0/1/1/1
3	PRQ	C	8	3	-	4/9/11/12	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	8	PRQ	CAL-NAN	-2.63	1.40	1.45
3	C	5	PRV	CD2-N1	-2.58	1.41	1.45
3	C	8	PRQ	CAL-NAN	-2.47	1.41	1.45
3	F	5	PRV	CD2-N1	-2.35	1.41	1.45
3	C	8	PRQ	CAK-CA	-2.13	1.48	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	5	PRV	CD1-CG-CA	-3.69	115.82	120.61
3	F	5	PRV	CD1-CG-CD2	2.80	119.99	117.08
3	F	5	PRV	CE2-CD2-N1	2.77	119.43	116.47
3	C	8	PRQ	CAH-CAK-CAL	2.71	119.89	117.08
3	C	8	PRQ	CAL-CAK-CA	-2.57	121.64	124.68
3	F	8	PRQ	CAH-CAK-CAL	2.20	119.37	117.08
3	F	8	PRQ	O-C-CAJ	-2.14	119.16	125.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	8	PRQ	CAI-CAL-NAN	2.13	118.75	116.47
3	C	8	PRQ	CAI-CAL-CAK	-2.05	119.32	121.32

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	5	PRV	C-CA-CG-CD2
3	C	5	PRV	N-CA-CG-CD2
3	F	5	PRV	C-CA-CG-CD2
3	C	8	PRQ	CAK-CA-CAJ-C
3	C	8	PRQ	CAK-CAL-NAN-OAC
3	F	8	PRQ	N-CA-CAK-CAH
3	F	8	PRQ	N-CA-CAK-CAL
3	F	8	PRQ	CAK-CAL-NAN-OAC
3	F	8	PRQ	CAI-CAL-NAN-OAC
3	C	8	PRQ	N-CA-CAJ-C
3	C	5	PRV	CE2-CD2-N1-O2
3	F	5	PRV	CE2-CD2-N1-O2
3	C	8	PRQ	CAI-CAL-NAN-OAC
3	C	5	PRV	CG-CD2-N1-O2
3	F	5	PRV	CG-CD2-N1-O2
3	F	8	PRQ	CAJ-CA-CAK-CAH

There are no ring outliers.

4 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	8	PRQ	5	0
3	C	5	PRV	2	0
3	F	5	PRV	4	0
3	C	8	PRQ	2	0

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$
4	GOL	A	1276	-	5,5,5	0.37	0	5,5,5	0.41	0
4	GOL	A	1277[A]	-	5,5,5	0.39	0	5,5,5	0.34	0
4	GOL	D	1278[B]	-	5,5,5	0.41	0	5,5,5	0.24	0
5	MES	E	1100	-	12,12,12	2.06	1 (8%)	15,16,16	5.96	8 (53%)
4	GOL	D	1279	-	5,5,5	0.41	0	5,5,5	0.51	0
5	MES	D	1277	-	12,12,12	2.22	1 (8%)	15,16,16	6.08	10 (66%)
4	GOL	A	1277[B]	-	5,5,5	0.42	0	5,5,5	0.28	0
4	GOL	D	1278[A]	-	5,5,5	0.39	0	5,5,5	0.23	0
5	MES	D	1275	-	12,12,12	2.21	1 (8%)	15,16,16	6.06	10 (66%)
5	MES	B	1100	-	12,12,12	2.30	1 (8%)	15,16,16	6.20	10 (66%)
4	GOL	D	1276	-	5,5,5	0.39	0	5,5,5	0.16	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
4	GOL	A	1276	-	-	4/4/4/4	-
4	GOL	A	1277[A]	-	-	1/4/4/4	-
4	GOL	D	1278[B]	-	-	2/4/4/4	-
5	MES	E	1100	-	-	2/6/14/14	0/1/1/1
4	GOL	D	1279	-	-	2/4/4/4	-
5	MES	D	1277	-	-	1/6/14/14	0/1/1/1
4	GOL	A	1277[B]	-	-	0/4/4/4	-
4	GOL	D	1278[A]	-	-	3/4/4/4	-
5	MES	D	1275	-	-	1/6/14/14	0/1/1/1
5	MES	B	1100	-	-	1/6/14/14	0/1/1/1
4	GOL	D	1276	-	-	1/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1100	MES	C8-S	-7.60	1.66	1.77
5	D	1277	MES	C8-S	-7.25	1.67	1.77
5	D	1275	MES	C8-S	-7.24	1.67	1.77
5	E	1100	MES	C8-S	-6.76	1.68	1.77

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	1275	MES	O1S-S-C8	-14.46	84.88	106.73
5	B	1100	MES	O1S-S-C8	-13.99	85.59	106.73
5	E	1100	MES	O1S-S-C8	-13.75	85.95	106.73
5	D	1277	MES	O2S-S-C8	-13.66	86.08	106.73
5	D	1277	MES	O1S-S-C8	-13.54	86.26	106.73
5	D	1275	MES	O2S-S-C8	-13.46	86.39	106.73
5	B	1100	MES	O2S-S-C8	-13.28	86.66	106.73
5	E	1100	MES	O2S-S-C8	-12.09	88.47	106.73
5	E	1100	MES	O3S-S-C8	-10.67	85.14	106.00
5	B	1100	MES	O3S-S-C8	-9.92	86.59	106.00
5	D	1277	MES	O3S-S-C8	-9.44	87.53	106.00
5	D	1275	MES	O3S-S-C8	-8.48	89.42	106.00
5	E	1100	MES	C5-N4-C3	5.78	121.29	108.84
5	D	1275	MES	C5-N4-C3	5.31	120.28	108.84
5	D	1277	MES	C5-N4-C3	5.02	119.65	108.84
5	B	1100	MES	C5-N4-C3	4.93	119.45	108.84
5	D	1277	MES	C6-C5-N4	-4.14	103.83	110.12
5	B	1100	MES	C2-C3-N4	-4.13	103.85	110.12
5	E	1100	MES	O3S-S-O1S	3.85	121.02	111.40
5	B	1100	MES	C7-N4-C5	3.50	120.56	111.24
5	D	1277	MES	C7-N4-C3	3.41	120.32	111.24
5	B	1100	MES	C6-C5-N4	-3.38	104.99	110.12
5	D	1275	MES	O3S-S-O1S	3.37	119.83	111.40
5	E	1100	MES	C7-N4-C5	3.35	120.16	111.24
5	E	1100	MES	O3S-S-O2S	3.34	119.76	111.40
5	D	1275	MES	C7-N4-C5	3.32	120.09	111.24
5	D	1277	MES	O3S-S-O2S	3.31	119.68	111.40
5	D	1277	MES	C7-N4-C5	3.29	120.00	111.24
5	B	1100	MES	C7-N4-C3	3.28	119.99	111.24
5	B	1100	MES	O3S-S-O2S	3.28	119.60	111.40
5	B	1100	MES	O3S-S-O1S	3.22	119.46	111.40
5	D	1275	MES	C7-N4-C3	3.15	119.63	111.24
5	D	1275	MES	O3S-S-O2S	3.00	118.91	111.40
5	D	1277	MES	O3S-S-O1S	2.91	118.69	111.40
5	D	1275	MES	C6-C5-N4	-2.85	105.79	110.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	1100	MES	C7-N4-C3	2.73	118.51	111.24
5	D	1277	MES	O2S-S-O1S	2.08	120.60	113.82
5	D	1275	MES	O2S-S-O1S	2.02	120.38	113.82

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1276	GOL	C1-C2-C3-O3
4	D	1278[A]	GOL	C1-C2-C3-O3
4	D	1278[B]	GOL	C1-C2-C3-O3
4	D	1278[B]	GOL	O2-C2-C3-O3
4	D	1279	GOL	C1-C2-C3-O3
5	B	1100	MES	C8-C7-N4-C5
5	D	1275	MES	C8-C7-N4-C3
4	D	1276	GOL	C1-C2-C3-O3
4	D	1279	GOL	O2-C2-C3-O3
4	A	1276	GOL	O2-C2-C3-O3
5	D	1277	MES	C8-C7-N4-C5
5	E	1100	MES	C8-C7-N4-C3
4	D	1278[A]	GOL	O1-C1-C2-O2
4	A	1276	GOL	O1-C1-C2-O2
4	A	1276	GOL	O1-C1-C2-C3
4	D	1278[A]	GOL	O2-C2-C3-O3
5	E	1100	MES	C7-C8-S-O3S
4	A	1277[A]	GOL	O2-C2-C3-O3

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1277[A]	GOL	1	0
4	D	1278[B]	GOL	1	0
4	A	1277[B]	GOL	1	0
4	D	1278[A]	GOL	4	0
4	D	1276	GOL	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	4 (1%)	72 71	12, 33, 62, 91	7 (2%)
1	D	275/275 (100%)	-0.34	3 (1%)	78 77	12, 26, 49, 63	13 (4%)
2	B	100/100 (100%)	-0.45	1 (1%)	79 79	18, 28, 59, 91	0
2	E	100/100 (100%)	-0.51	3 (3%)	52 51	13, 23, 46, 55	3 (3%)
3	C	7/9 (77%)	1.14	2 (28%)	1 1	19, 26, 88, 92	0
3	F	7/9 (77%)	2.85	5 (71%)	0 0	33, 43, 77, 103	0
All	All	764/768 (99%)	-0.28	18 (2%)	59 59	12, 28, 57, 103	23 (3%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	7	PHE	4.9
3	F	7	PHE	4.5
3	C	6	VAL	4.4
1	A	34	VAL	4.4
1	D	112	TYR	4.3
3	F	4	GLY	3.5
3	F	6	VAL	3.3
3	F	3	LEU	3.2
1	D	108	PHE	3.0
3	F	9	VAL	2.8
2	E	98	ASP	2.8
1	D	181	THR	2.8
1	A	274	TRP	2.8
2	E	97	ARG	2.7
2	E	96	ASP	2.4
2	B	99	MET	2.2
1	A	223	ASP	2.1
1	A	273	ARG	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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
3	PRV	F	5	13/14	0.50	0.19	65,70,75,76	0
3	PRV	C	5	13/14	0.74	0.15	51,72,85,85	0
3	PRQ	F	8	14/15	0.76	0.25	46,94,106,106	0
3	PRQ	C	8	14/15	0.78	0.15	42,96,105,107	0

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
4	GOL	D	1276	6/6	0.72	0.12	61,66,68,71	0
5	MES	B	1100	12/12	0.80	0.13	77,79,82,84	0
4	GOL	A	1277[A]	6/6	0.82	0.15	21,27,33,34	6
4	GOL	A	1277[B]	6/6	0.82	0.15	25,26,28,30	6
5	MES	D	1277	12/12	0.86	0.15	44,71,75,77	0
4	GOL	D	1278[A]	6/6	0.88	0.17	25,26,28,29	6
4	GOL	D	1278[B]	6/6	0.88	0.17	28,29,30,33	6
5	MES	D	1275	12/12	0.89	0.13	37,55,59,59	0
5	MES	E	1100	12/12	0.89	0.11	43,45,53,53	0
4	GOL	D	1279	6/6	0.90	0.12	42,43,44,45	0
4	GOL	A	1276	6/6	0.91	0.10	34,39,42,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.