



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:04 PM UTC

PDB ID : 7JYX / pdb_00007jyx
Title : Crystal Structure of HLA A*2402 in complex with TYQWIIRNWET, an 11-mer epitope from Influenza
Authors : Gras, S.; Nguyen, A.T.; Szeto, C.; Rossjohn, J.
Deposited on : 2020-09-01
Resolution : 2.95 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
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

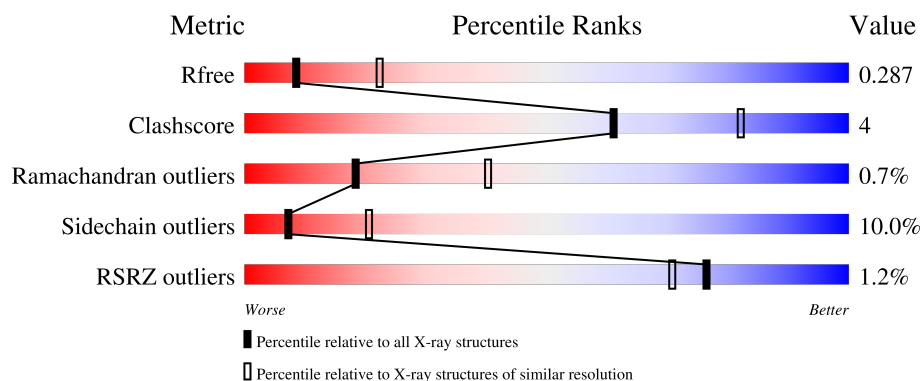
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	278	3% 79% 17% ..
1	D	278	75% 21% ..
2	B	100	76% 20% ..
2	E	100	% 71% 27% .
3	C	11	82% 18%

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Mol	Chain	Length	Quality of chain
3	F	11	 A horizontal bar chart showing the quality of chain F. The bar is divided into four segments: green (55%), yellow (27%), orange (9%), and red (9%).

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6373 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 MHC class I antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	272	Total	C	N	O	S	0	1	0
			2207	1371	401	425	10			
1	D	274	Total	C	N	O	S	0	0	0
			2221	1382	403	426	10			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	98	Total	C	N	O	S	0	0	0
			820	523	139	156	2			
2	E	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			

There are 2 discrepancies between the modelled and reference sequences:

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

- Molecule 3 is a protein called PB2 peptide from Influenza, TYQWIIRNWET.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	11	Total	C	N	O	0	0	0
			108	71	18	19			
3	F	11	Total	C	N	O	0	0	0
			108	71	18	19			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	28	Total	O	0	0
			28	28		

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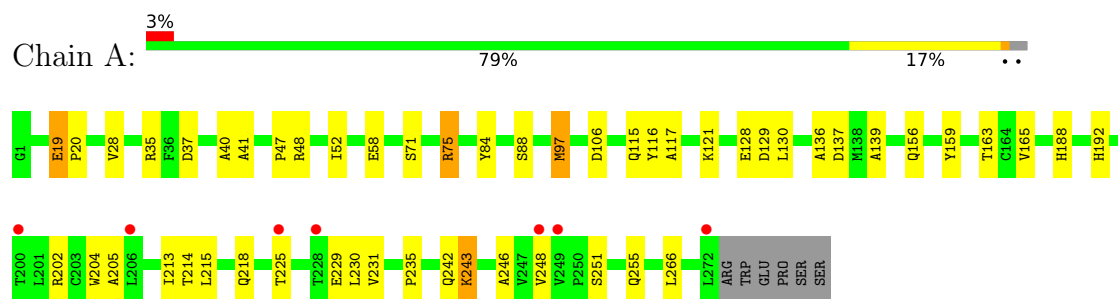
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	10	Total 10	O 10	0	0
4	C	2	Total 2	O 2	0	0
4	D	19	Total 19	O 19	0	0
4	E	12	Total 12	O 12	0	0
4	F	1	Total 1	O 1	0	0

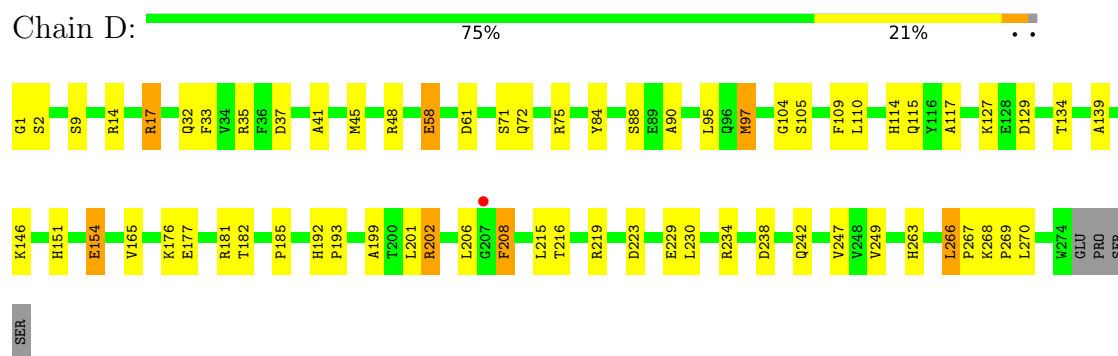
3 Residue-property plots [i](#)

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.

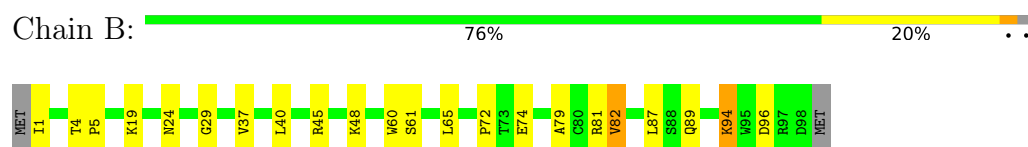
- Molecule 1: MHC class I antigen



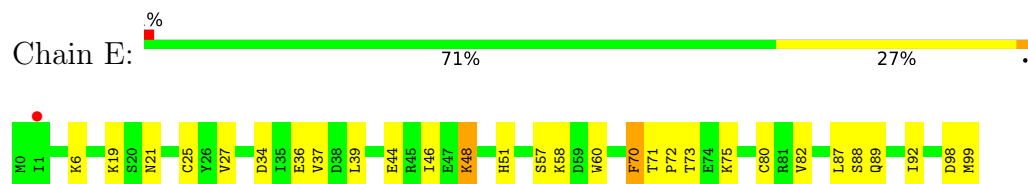
- Molecule 1: MHC class I antigen



- Molecule 2: Beta-2-microglobulin

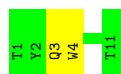


- Molecule 2: Beta-2-microglobulin



- Molecule 3: PB2 peptide from Influenza, TYQWIIRNWET

Chain C:  82% 18%



- Molecule 3: PB2 peptide from Influenza, TYQWIIRNWET

Chain F:  55% 27% 9% 9%



4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	89.68Å 43.65Å 236.49Å 90.00° 95.56° 90.00°	Depositor
Resolution (Å)	39.21 – 2.95 39.21 – 2.95	Depositor EDS
% Data completeness (in resolution range)	99.4 (39.21-2.95) 99.3 (39.21-2.95)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 2.95Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.208 , 0.275 0.217 , 0.287	Depositor DCC
R_{free} test set	989 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	59.3	Xtriage
Anisotropy	0.311	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 52.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6373	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.96% 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 ⓘ

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	0.83	0/2265	1.36	18/3069 (0.6%)
1	D	0.85	0/2281	1.34	17/3092 (0.5%)
2	B	0.85	0/843	1.30	3/1142 (0.3%)
2	E	0.81	0/860	1.35	9/1162 (0.8%)
3	C	0.77	0/112	1.06	0/152
3	F	0.73	0/112	1.11	1/152 (0.7%)
All	All	0.83	0/6473	1.33	48/8769 (0.5%)

There are no bond length outliers.

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	71	THR	CB-CA-C	7.03	117.50	110.33
1	D	33	PHE	CA-CB-CG	7.00	120.80	113.80
1	A	40	ALA	CA-C-N	6.80	129.72	120.54
1	A	40	ALA	C-N-CA	6.80	129.72	120.54
1	D	208	PHE	CA-CB-CG	6.74	120.54	113.80
1	A	75	ARG	CA-C-N	6.50	129.31	120.54
1	A	75	ARG	C-N-CA	6.50	129.31	120.54
2	B	96	ASP	CA-CB-CG	6.22	118.82	112.60
1	D	45	MET	N-CA-C	-6.17	99.11	108.67
2	E	70	PHE	CA-CB-CG	6.12	119.92	113.80
2	E	87	LEU	CA-C-N	6.00	128.23	120.44
2	E	87	LEU	C-N-CA	6.00	128.23	120.44
3	F	8	ASN	CA-CB-CG	5.91	118.51	112.60
1	A	41	ALA	CA-C-N	5.89	128.49	120.54
1	A	41	ALA	C-N-CA	5.89	128.49	120.54
1	A	28	VAL	N-CA-C	-5.83	98.01	107.28
1	A	75	ARG	N-CA-C	-5.81	104.86	111.14
1	D	238	ASP	CA-CB-CG	5.71	118.31	112.60
2	E	21	ASN	CA-CB-CG	5.69	118.29	112.60
1	D	269	PRO	CA-C-N	5.68	128.90	120.95
1	D	269	PRO	C-N-CA	5.68	128.90	120.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	75	ARG	N-CA-C	-5.57	105.21	112.23
1	D	37	ASP	CA-CB-CG	5.57	118.17	112.60
1	D	41	ALA	CA-C-N	5.57	132.18	121.54
1	D	41	ALA	C-N-CA	5.57	132.18	121.54
1	D	249	VAL	N-CA-C	5.56	114.22	107.61
2	E	73	THR	CA-C-N	5.56	127.99	120.38
2	E	73	THR	C-N-CA	5.56	127.99	120.38
1	A	19	GLU	CB-CG-CD	5.54	122.03	112.60
1	A	205	ALA	N-CA-C	-5.54	99.86	108.90
1	A	137	ASP	CA-C-N	5.50	127.65	120.28
1	A	137	ASP	C-N-CA	5.50	127.65	120.28
1	A	106	ASP	CA-CB-CG	5.46	118.06	112.60
1	D	61	ASP	CA-CB-CG	5.36	117.96	112.60
2	B	87	LEU	CA-C-N	5.32	127.36	120.44
2	B	87	LEU	C-N-CA	5.32	127.36	120.44
1	A	115	GLN	N-CA-C	5.26	117.07	109.07
1	A	37	ASP	CA-CB-CG	5.17	117.77	112.60
1	D	71	SER	CA-C-N	5.16	127.15	120.44
1	D	71	SER	C-N-CA	5.16	127.15	120.44
1	D	58	GLU	CA-C-N	5.11	127.08	120.44
1	D	58	GLU	C-N-CA	5.11	127.08	120.44
2	E	75	LYS	CA-C-N	5.10	127.96	120.71
2	E	75	LYS	C-N-CA	5.10	127.96	120.71
1	D	129	ASP	CA-CB-CG	5.09	117.69	112.60
1	A	231	VAL	CA-C-N	5.03	128.01	120.87
1	A	231	VAL	C-N-CA	5.03	128.01	120.87
1	A	129	ASP	CA-CB-CG	5.02	117.62	112.60

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	2207	0	2073	16	0
1	D	2221	0	2082	20	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	820	0	785	10	0
2	E	837	0	805	6	0
3	C	108	0	100	4	0
3	F	108	0	100	3	0
4	A	28	0	0	1	0
4	B	10	0	0	0	0
4	C	2	0	0	0	0
4	D	19	0	0	0	0
4	E	12	0	0	0	0
4	F	1	0	0	0	0
All	All	6373	0	5945	51	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 4.

All (51) 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:A:121:LYS:HD2	2:B:1:ILE:HG12	1.66	0.77
2:E:25:CYS:HB2	2:E:39:LEU:HD21	1.79	0.65
1:A:35:ARG:HD2	1:A:48[B]:ARG:HE	1.64	0.63
2:B:24:ASN:HB3	2:B:65:LEU:HD11	1.85	0.58
1:D:185:PRO:HB3	1:D:208:PHE:HB3	1.87	0.55
3:F:8:ASN:HD22	3:F:10:GLU:H	1.54	0.55
2:B:19:LYS:O	2:B:72:PRO:HD2	2.06	0.55
1:D:234:ARG:HG3	1:D:242:GLN:HG3	1.89	0.55
3:F:3:GLN:HG3	3:F:4:TRP:N	2.22	0.54
2:E:19:LYS:O	2:E:72:PRO:HD2	2.08	0.54
1:A:97:MET:HG3	1:A:116:TYR:CE1	2.41	0.53
2:B:79:ALA:HB2	2:B:94:LYS:HE3	1.90	0.52
1:D:35:ARG:HG3	1:D:48:ARG:CG	2.40	0.52
2:B:40:LEU:HD11	2:B:81:ARG:HB2	1.92	0.52
1:D:177:GLU:O	1:D:181:ARG:HB3	2.10	0.52
1:A:20:PRO:HG2	1:A:75:ARG:HG3	1.91	0.51
1:A:235:PRO:HG2	2:B:65:LEU:HD22	1.91	0.51
1:D:35:ARG:HG3	1:D:48:ARG:HG3	1.94	0.50
2:E:46:ILE:HG22	2:E:48:LYS:H	1.76	0.49
1:D:151:HIS:O	1:D:154:GLU:HG2	2.13	0.49
1:D:117:ALA:HB2	2:E:60:TRP:CD2	2.48	0.48
1:A:84:TYR:HB3	1:A:139:ALA:HB1	1.95	0.48
1:A:188:HIS:HB3	1:A:204:TRP:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:GLU:O	1:A:130:LEU:HD12	2.14	0.47
1:D:1:GLY:HA2	1:D:105:SER:HB3	1.95	0.47
2:E:37:VAL:HG22	2:E:82:VAL:HG22	1.97	0.46
1:A:47:PRO:HB3	1:A:52:ILE:HG23	1.97	0.46
1:A:156:GLN:HE21	3:C:3:GLN:HE22	1.63	0.46
1:D:14:ARG:HB2	1:D:17:ARG:HG3	1.98	0.45
1:D:185:PRO:HD2	1:D:266:LEU:HD13	1.97	0.45
1:D:88:SER:C	1:D:90:ALA:H	2.24	0.45
1:D:193:PRO:HA	1:D:199:ALA:HA	1.98	0.45
2:B:37:VAL:HG22	2:B:82:VAL:HG13	1.98	0.44
1:D:263:HIS:HB3	1:D:266:LEU:HB2	1.99	0.44
1:A:202:ARG:HB3	1:A:246:ALA:HB2	2.00	0.43
1:D:109:PHE:HB2	1:D:165:VAL:HG11	2.00	0.43
1:A:156:GLN:HE21	3:C:3:GLN:NE2	2.17	0.43
1:A:58:GLU:HB2	4:A:310:HOH:O	2.19	0.43
3:C:3:GLN:HG2	3:C:4:TRP:N	2.34	0.43
1:D:146:LYS:HD3	3:F:10:GLU:O	2.19	0.42
1:D:35:ARG:CG	1:D:48:ARG:HG3	2.49	0.42
1:D:2:SER:HA	1:D:104:GLY:HA2	2.01	0.41
2:B:29:GLY:HA2	2:B:61:SER:OG	2.19	0.41
1:D:192:HIS:CD2	1:D:202:ARG:HH21	2.38	0.41
2:E:25:CYS:SG	2:E:80:CYS:HB2	2.61	0.41
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.55	0.41
1:A:215:LEU:HD23	1:A:243:LYS:HD2	2.02	0.41
1:D:84:TYR:HB3	1:D:139:ALA:HB1	2.03	0.41
2:B:4:THR:HA	2:B:5:PRO:HD3	1.96	0.41
1:D:97:MET:HE1	1:D:114:HIS:CD2	2.56	0.40
1:A:159:TYR:CG	3:C:3:GLN:HG3	2.55	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	271/278 (98%)	255 (94%)	14 (5%)	2 (1%)	18	40
1	D	272/278 (98%)	235 (86%)	34 (12%)	3 (1%)	11	29
2	B	96/100 (96%)	89 (93%)	7 (7%)	0	100	100
2	E	98/100 (98%)	92 (94%)	6 (6%)	0	100	100
3	C	9/11 (82%)	8 (89%)	1 (11%)	0	100	100
3	F	9/11 (82%)	9 (100%)	0	0	100	100
All	All	755/778 (97%)	688 (91%)	62 (8%)	5 (1%)	18	40

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	176	LYS
1	D	223	ASP
1	A	136	ALA
1	A	251	SER
1	D	267	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	229/234 (98%)	211 (92%)	18 (8%)	11	29
1	D	230/234 (98%)	205 (89%)	25 (11%)	6	18
2	B	93/95 (98%)	87 (94%)	6 (6%)	15	37
2	E	95/95 (100%)	80 (84%)	15 (16%)	2	8
3	C	11/11 (100%)	11 (100%)	0	100	100
3	F	11/11 (100%)	8 (73%)	3 (27%)	0	0
All	All	669/680 (98%)	602 (90%)	67 (10%)	7	20

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	GLU
1	A	71	SER
1	A	88	SER
1	A	97	MET
1	A	163	THR
1	A	165	VAL
1	A	192	HIS
1	A	213	ILE
1	A	214	THR
1	A	218	GLN
1	A	225	THR
1	A	229	GLU
1	A	230	LEU
1	A	242	GLN
1	A	243	LYS
1	A	248	VAL
1	A	255	GLN
1	A	266	LEU
2	B	45	ARG
2	B	48	LYS
2	B	74	GLU
2	B	82	VAL
2	B	89	GLN
2	B	94	LYS
1	D	9	SER
1	D	17	ARG
1	D	32	GLN
1	D	58	GLU
1	D	72	GLN
1	D	95	LEU
1	D	97	MET
1	D	110	LEU
1	D	115	GLN
1	D	127	LYS
1	D	134	THR
1	D	154	GLU
1	D	182	THR
1	D	201	LEU
1	D	202	ARG
1	D	206	LEU
1	D	215	LEU
1	D	216	THR
1	D	219	ARG

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Mol	Chain	Res	Type
1	D	229	GLU
1	D	230	LEU
1	D	247	VAL
1	D	266	LEU
1	D	268	LYS
1	D	270	LEU
2	E	6	LYS
2	E	27	VAL
2	E	34	ASP
2	E	36	GLU
2	E	44	GLU
2	E	48	LYS
2	E	51	HIS
2	E	57	SER
2	E	58	LYS
2	E	70	PHE
2	E	88	SER
2	E	89	GLN
2	E	92	ILE
2	E	98	ASP
2	E	99	MET
3	F	3	GLN
3	F	6	ILE
3	F	8	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	72	GLN
1	A	87	GLN
1	A	93	HIS
1	A	114	HIS
1	A	155	GLN
1	A	180	GLN
1	A	218	GLN
1	A	226	GLN
1	A	242	GLN
1	A	255	GLN
2	B	17	ASN
2	B	83	ASN
3	C	3	GLN

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Mol	Chain	Res	Type
1	D	32	GLN
1	D	87	GLN
1	D	93	HIS
1	D	174	ASN
1	D	191	HIS
1	D	218	GLN
1	D	226	GLN
2	E	31	HIS
2	E	42	ASN
3	F	3	GLN
3	F	8	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	272/278 (97%)	0.17	7 (2%) 57 49	20, 52, 124, 148	1 (0%)
1	D	274/278 (98%)	0.11	1 (0%) 88 86	32, 59, 97, 109	0
2	B	98/100 (98%)	0.36	0 100 100	43, 79, 101, 112	0
2	E	100/100 (100%)	0.22	1 (1%) 79 74	31, 64, 89, 100	0
3	C	11/11 (100%)	-0.35	0 100 100	31, 39, 46, 58	0
3	F	11/11 (100%)	-0.21	0 100 100	31, 41, 50, 51	0
All	All	766/778 (98%)	0.17	9 (1%) 76 71	20, 60, 108, 148	1 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	1	ILE	3.0
1	A	249	VAL	2.8
1	A	248	VAL	2.6
1	A	206	LEU	2.5
1	A	228	THR	2.4
1	A	200	THR	2.3
1	A	225	THR	2.2
1	D	207	GLY	2.2
1	A	272	LEU	2.2

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.