



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:35 PM UTC

PDB ID : 3PL6 / pdb_00003pl6
Title : Structure of Autoimmune TCR Hy.1B11 in complex with HLA-DQ1 and MBP 85-99
Authors : Sethi, D.K.; Wucherpennig, K.W.
Deposited on : 2010-11-13
Resolution : 2.55 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

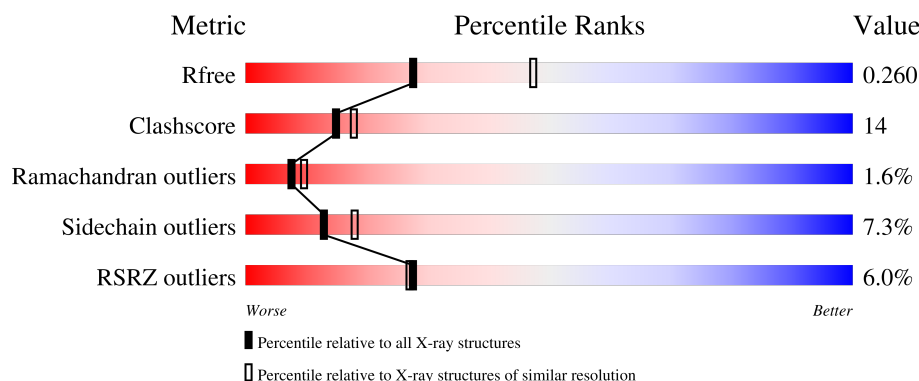
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	194	8% 61% 27% 5% • 6%
2	B	202	2% 67% 19% • 10%
3	C	206	8% 64% 23% • • 7%
4	D	268	4% 65% 24% 7% •

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6526 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 II HLA-DQ-alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	182	Total	C	N	O	S	0	0	0
			1438	922	237	274	5			

- Molecule 2 is a protein called MHC class II HLA-DQ-beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	181	Total	C	N	O	S	0	0	0
			1482	935	267	274	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	199	VAL	-	expression tag	UNP Q67AJ6
B	200	ASP	-	expression tag	UNP Q67AJ6

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	191	Total	C	N	O	S	0	0	0
			1486	923	251	305	7			

- Molecule 4 is a protein called MBP peptide / T-cell receptor beta chain chimera.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	257	Total	C	N	O	S	0	0	0
			1999	1263	348	383	5			

There are 9 discrepancies between the modelled and reference sequences:

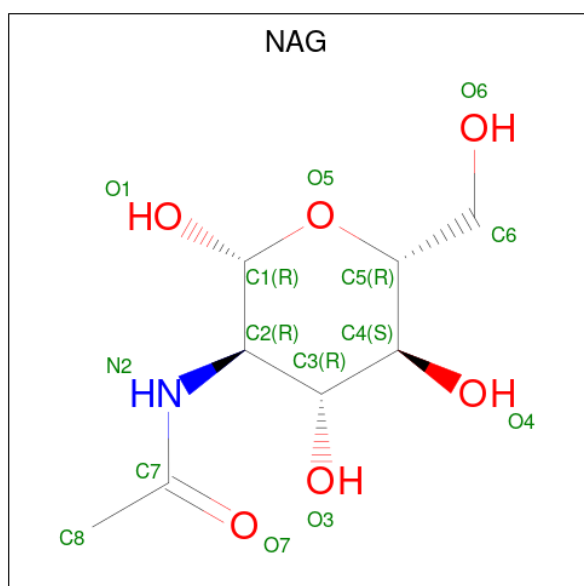
Chain	Residue	Modelled	Actual	Comment	Reference
D	1	MET	-	expression tag	UNP D3YTB3

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Chain	Residue	Modelled	Actual	Comment	Reference
D	2	LYS	-	expression tag	UNP D3YTB3
D	18	GLY	-	expression tag	PDB 3PL6
D	19	GLY	-	expression tag	PDB 3PL6
D	20	SER	-	expression tag	PDB 3PL6
D	21	GLY	-	expression tag	PDB 3PL6
D	22	GLY	-	expression tag	PDB 3PL6
D	23	GLY	-	expression tag	PDB 3PL6
D	24	GLY	-	expression tag	PDB 3PL6

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	21	Total	O	0	0
			21	21		
6	B	13	Total	O	0	0
			13	13		
6	C	30	Total	O	0	0
			30	30		

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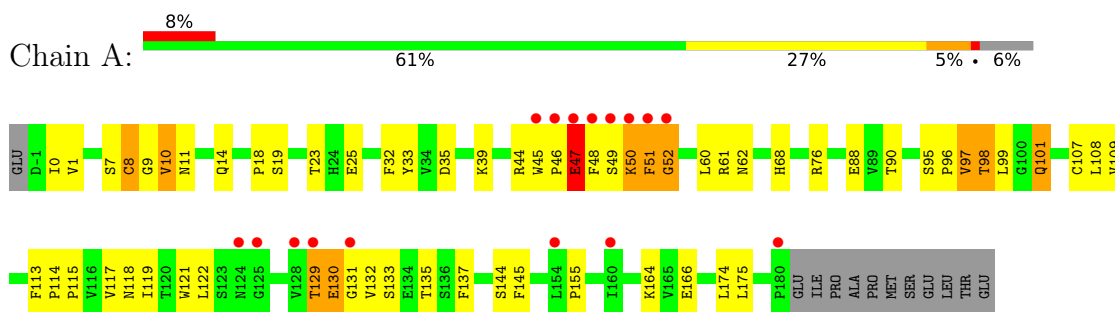
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	29	Total	O	0	0
			29	29		

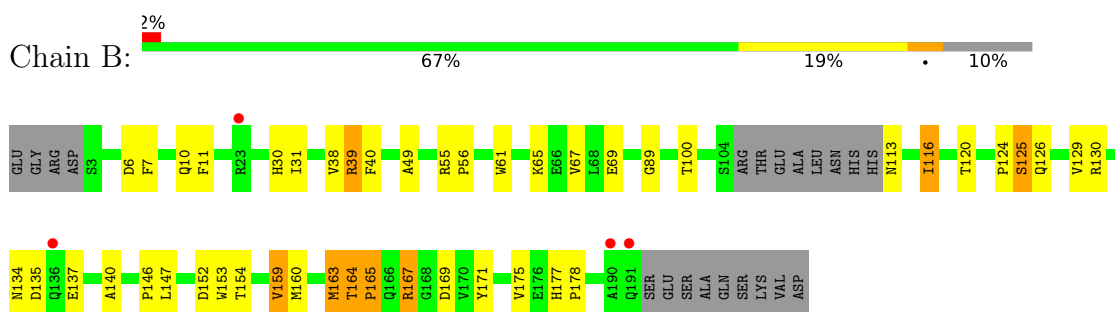
3 Residue-property plots [i](#)

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.

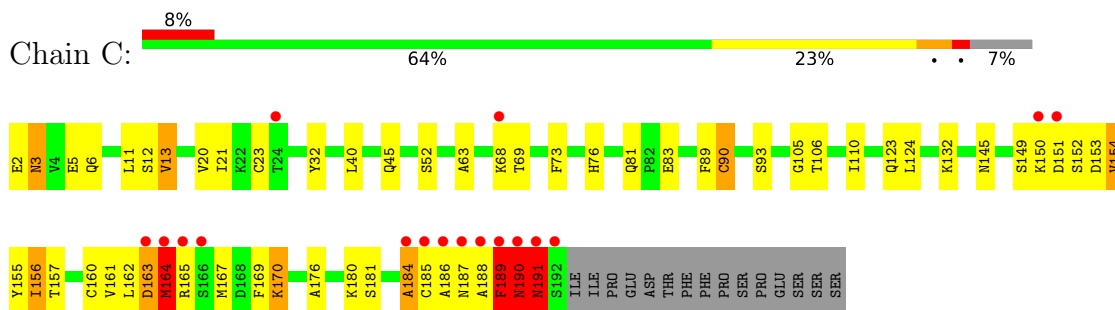
- Molecule 1: MHC class II HLA-DQ-alpha chain



- Molecule 2: MHC class II HLA-DQ-beta chain

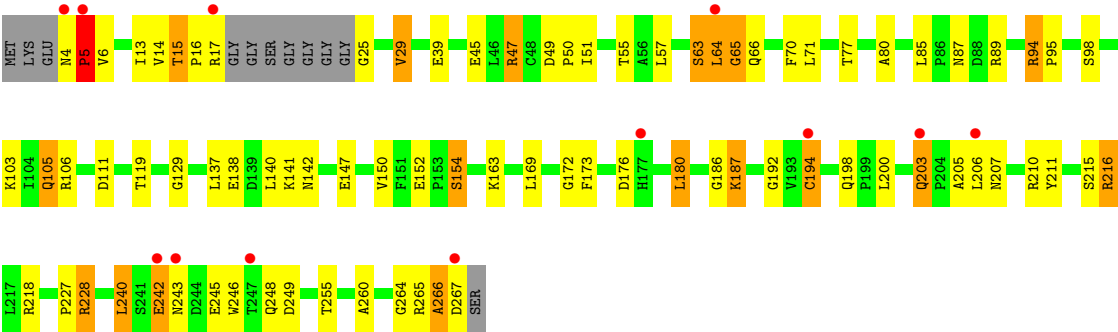


- Molecule 3: T-cell receptor alpha chain



- Molecule 4: MBP peptide / T-cell receptor beta chain chimera





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	73.12Å 124.81Å 134.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.00 – 2.55 42.00 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.6 (42.00-2.55) 99.5 (42.00-2.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 2.54Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.232 , 0.258 0.233 , 0.260	Depositor DCC
R_{free} test set	2045 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	55.9	Xtriage
Anisotropy	0.601	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 28.9	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	0.93	EDS
Total number of atoms	6526	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.73% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG

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	0.45	0/1482	0.98	10/2023 (0.5%)
2	B	0.47	0/1519	1.04	11/2067 (0.5%)
3	C	0.48	0/1515	1.04	7/2052 (0.3%)
4	D	0.47	0/2054	1.09	21/2803 (0.7%)
All	All	0.47	0/6570	1.04	49/8945 (0.5%)

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	191	ASN	N-CA-C	14.92	130.94	110.35
4	D	65	GLY	N-CA-C	-9.97	102.82	114.69
4	D	152	GLU	CA-C-N	8.81	128.89	119.90
4	D	152	GLU	C-N-CA	8.81	128.89	119.90
2	B	140	ALA	N-CA-C	-7.98	99.50	110.35
1	A	52	GLY	N-CA-C	-7.71	94.92	113.18
4	D	203	GLN	CA-C-N	7.32	127.96	119.47
4	D	203	GLN	C-N-CA	7.32	127.96	119.47
2	B	167	ARG	N-CA-C	7.13	119.91	111.71
1	A	131	GLY	N-CA-C	-6.88	105.59	115.64
1	A	8	CYS	N-CA-C	-6.71	102.19	110.65
2	B	164	THR	CA-C-N	6.62	126.85	119.90
2	B	164	THR	C-N-CA	6.62	126.85	119.90
4	D	63	SER	N-CA-C	-6.53	98.59	109.24
4	D	15	THR	CA-C-N	6.42	126.37	120.21
4	D	15	THR	C-N-CA	6.42	126.37	120.21
3	C	189	PHE	N-CA-C	6.41	124.45	110.80
2	B	120	THR	N-CA-C	6.37	119.39	109.52
4	D	94	ARG	CA-C-N	6.21	127.60	119.84
4	D	94	ARG	C-N-CA	6.21	127.60	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	85	LEU	CA-C-N	6.15	126.33	119.32
4	D	85	LEU	C-N-CA	6.15	126.33	119.32
3	C	181	SER	N-CA-C	6.14	118.76	111.33
4	D	70	PHE	N-CA-C	6.12	118.87	110.24
2	B	175	VAL	N-CA-C	6.11	116.66	108.11
2	B	116	ILE	N-CA-C	6.02	116.54	107.75
4	D	49	ASP	CA-C-N	5.88	125.85	120.21
4	D	49	ASP	C-N-CA	5.88	125.85	120.21
1	A	144	SER	N-CA-C	-5.79	101.42	110.17
4	D	173	PHE	N-CA-C	5.75	119.03	110.52
3	C	188	ALA	N-CA-C	-5.75	98.55	110.80
2	B	125	SER	N-CA-C	5.65	118.38	111.82
1	A	47	GLU	N-CA-C	-5.39	107.07	113.97
2	B	146	PRO	N-CA-C	-5.36	103.15	111.13
2	B	169	ASP	N-CA-C	5.34	118.18	110.28
2	B	165	PRO	N-CA-C	5.33	118.84	110.80
1	A	135	THR	N-CA-C	-5.30	102.90	110.59
4	D	255	THR	N-CA-C	-5.30	102.45	110.24
3	C	184	ALA	N-CA-C	5.26	122.21	113.89
4	D	207	ASN	N-CA-C	-5.17	106.98	113.28
3	C	90	CYS	N-CA-C	-5.13	102.47	110.42
1	A	10	VAL	N-CA-C	-5.12	99.09	106.88
1	A	35	ASP	N-CA-C	-5.12	101.78	109.15
4	D	51	ILE	N-CA-C	-5.11	102.91	109.30
1	A	48	PHE	N-CA-C	-5.08	106.68	112.92
4	D	172	GLY	N-CA-C	5.06	122.09	115.36
4	D	147	GLU	N-CA-C	-5.06	100.93	109.07
3	C	132	LYS	N-CA-C	5.04	117.35	110.55
1	A	97	VAL	N-CA-C	5.01	115.90	108.23

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	1438	0	1347	44	0
2	B	1482	0	1425	31	0
3	C	1486	0	1411	56	0
4	D	1999	0	1875	61	0
5	A	14	0	13	0	0
5	B	14	0	13	0	0
6	A	21	0	0	1	0
6	B	13	0	0	0	0
6	C	30	0	0	2	0
6	D	29	0	0	1	0
All	All	6526	0	6084	173	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 14.

All (173) 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)
3:C:160:CYS:HB3	4:D:216:ARG:HH12	1.16	1.10
3:C:160:CYS:HB3	4:D:216:ARG:NH1	1.79	0.97
3:C:160:CYS:CB	4:D:216:ARG:HH12	1.78	0.96
2:B:129:VAL:HG11	2:B:159:VAL:HG11	1.53	0.90
1:A:129:THR:OG1	1:A:130:GLU:HG2	1.74	0.86
4:D:228:ARG:HH11	4:D:228:ARG:HB2	1.42	0.85
4:D:64:LEU:HD22	4:D:65:GLY:H	1.41	0.84
4:D:140:LEU:HD13	4:D:240:LEU:HD22	1.58	0.84
2:B:126:GLN:H	2:B:126:GLN:CD	1.85	0.84
4:D:210:ARG:HG3	6:D:296:HOH:O	1.80	0.81
3:C:163:ASP:HB2	4:D:192:GLY:H	1.47	0.79
3:C:6:GLN:NE2	3:C:106:THR:HG22	1.99	0.78
4:D:94:ARG:HD2	4:D:98:SER:O	1.83	0.78
4:D:228:ARG:HB2	4:D:228:ARG:NH1	1.98	0.78
1:A:68:HIS:CD2	4:D:14:VAL:HG13	2.19	0.78
4:D:205:ALA:C	4:D:206:LEU:HD22	2.14	0.73
3:C:21:ILE:HG12	3:C:106:THR:HG21	1.71	0.73
4:D:4:ASN:CG	4:D:5:PRO:HD2	2.13	0.72
2:B:125:SER:HB2	2:B:147:LEU:HD21	1.72	0.70
3:C:145:ASN:HB3	3:C:190:ASN:OD1	1.92	0.69
3:C:164:MET:HG2	3:C:170:LYS:HD3	1.75	0.67
3:C:190:ASN:HB2	6:C:211:HOH:O	1.93	0.67
2:B:125:SER:HB2	2:B:147:LEU:CD2	2.25	0.66
4:D:227:PRO:HA	4:D:264:GLY:O	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:40:LEU:HD22	3:C:40:LEU:H	1.58	0.66
3:C:163:ASP:CB	4:D:192:GLY:H	2.08	0.66
1:A:49:SER:O	1:A:51:PHE:N	2.29	0.66
2:B:113:ASN:HB3	2:B:163:MET:O	1.95	0.65
3:C:184:ALA:C	3:C:186:ALA:H	2.04	0.65
3:C:189:PHE:O	3:C:190:ASN:ND2	2.32	0.62
3:C:191:ASN:N	3:C:191:ASN:OD1	2.31	0.62
1:A:101:GLN:O	1:A:155:PRO:HD2	2.00	0.62
4:D:200:LEU:H	4:D:200:LEU:HD23	1.64	0.61
2:B:163:MET:HE3	2:B:165:PRO:HD3	1.82	0.60
4:D:63:SER:HB2	4:D:66:GLN:OE1	2.02	0.59
4:D:141:LYS:O	4:D:141:LYS:HD3	2.02	0.59
3:C:149:SER:HB3	3:C:156:ILE:HG23	1.85	0.59
3:C:156:ILE:HD12	3:C:157:THR:O	2.03	0.59
1:A:11:ASN:HB2	2:B:11:PHE:HB3	1.85	0.59
2:B:124:PRO:HG2	2:B:126:GLN:OE1	2.03	0.59
1:A:7:SER:HB3	1:A:10:VAL:HG23	1.84	0.58
4:D:228:ARG:HH11	4:D:228:ARG:CB	2.14	0.58
4:D:180:LEU:C	4:D:180:LEU:HD23	2.28	0.58
4:D:246:TRP:CE2	4:D:248:GLN:HB2	2.39	0.57
1:A:44:ARG:HB3	1:A:45:TRP:CE3	2.39	0.57
3:C:52:SER:HB2	3:C:68:LYS:HD2	1.86	0.57
2:B:55:ARG:HB2	2:B:56:PRO:HD3	1.86	0.57
4:D:25:GLY:HA3	4:D:50:PRO:O	2.03	0.57
3:C:163:ASP:C	3:C:165:ARG:H	2.13	0.56
1:A:174:LEU:C	1:A:174:LEU:HD23	2.30	0.56
3:C:2:GLU:O	3:C:3:ASN:HB2	2.04	0.56
3:C:167:MET:HE1	4:D:163:LYS:HE3	1.87	0.56
3:C:162:LEU:HD11	4:D:218:ARG:HB2	1.88	0.56
1:A:98:THR:O	1:A:98:THR:HG22	2.06	0.55
4:D:29:VAL:HG22	4:D:129:GLY:HA3	1.88	0.55
1:A:115:PRO:HG3	1:A:145:PHE:CD2	2.41	0.55
1:A:121:TRP:O	1:A:122:LEU:HD12	2.06	0.55
4:D:176:ASP:HB3	4:D:211:TYR:CE2	2.41	0.55
1:A:98:THR:O	1:A:101:GLN:HB2	2.07	0.55
4:D:89:ARG:NH2	4:D:111:ASP:OD2	2.40	0.55
3:C:163:ASP:O	3:C:165:ARG:HG2	2.06	0.54
3:C:189:PHE:CD1	3:C:189:PHE:C	2.84	0.54
1:A:132:VAL:HG12	1:A:133:SER:N	2.22	0.54
1:A:164:LYS:HG3	1:A:175:LEU:CD2	2.37	0.54
3:C:20:VAL:C	3:C:21:ILE:HD12	2.34	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:163:ASP:HB2	4:D:192:GLY:N	2.19	0.53
1:A:14:GLN:NE2	2:B:6:ASP:OD1	2.40	0.53
1:A:68:HIS:NE2	4:D:14:VAL:HG13	2.24	0.52
2:B:130:ARG:HD3	2:B:137:GLU:OE2	2.10	0.52
4:D:265:ARG:O	4:D:266:ALA:O	2.27	0.52
2:B:177:HIS:CG	2:B:178:PRO:HD2	2.45	0.51
2:B:30:HIS:C	2:B:31:ILE:HG13	2.33	0.51
3:C:63:ALA:HB3	3:C:76:HIS:HB2	1.93	0.51
1:A:99:LEU:HA	1:A:155:PRO:HG2	1.92	0.51
1:A:39:LYS:HG2	1:A:60:LEU:HD21	1.91	0.51
1:A:129:THR:HG1	1:A:130:GLU:HG2	1.74	0.51
3:C:124:LEU:HD12	3:C:124:LEU:N	2.26	0.51
3:C:184:ALA:O	3:C:186:ALA:N	2.43	0.51
1:A:164:LYS:HG3	1:A:175:LEU:HD21	1.93	0.50
4:D:245:GLU:HA	4:D:245:GLU:OE1	2.10	0.50
4:D:150:VAL:HG12	4:D:260:ALA:HB1	1.93	0.50
1:A:113:PHE:HA	1:A:114:PRO:C	2.37	0.49
4:D:13:ILE:HG22	4:D:14:VAL:O	2.11	0.49
4:D:55:THR:O	4:D:94:ARG:NH2	2.45	0.49
2:B:129:VAL:HG11	2:B:159:VAL:CG1	2.36	0.49
4:D:186:GLY:O	4:D:187:LYS:HB2	2.10	0.49
1:A:107:CYS:HB2	1:A:121:TRP:CH2	2.47	0.49
4:D:50:PRO:CB	4:D:119:THR:HG21	2.42	0.49
1:A:90:THR:HG23	1:A:108:LEU:HB3	1.93	0.49
4:D:47:ARG:HH11	4:D:47:ARG:HB3	1.78	0.48
1:A:115:PRO:HG3	1:A:145:PHE:CE2	2.47	0.48
4:D:242:GLU:HG2	4:D:243:ASN:N	2.20	0.48
4:D:29:VAL:HG22	4:D:129:GLY:CA	2.44	0.48
3:C:6:GLN:HE21	3:C:106:THR:HG22	1.77	0.48
3:C:163:ASP:O	3:C:165:ARG:N	2.46	0.48
3:C:184:ALA:C	3:C:186:ALA:N	2.70	0.48
2:B:10:GLN:HB2	2:B:31:ILE:HB	1.96	0.47
1:A:109:VAL:HG21	1:A:119:ILE:CD1	2.45	0.47
3:C:5:GLU:O	3:C:23:CYS:HA	2.15	0.47
3:C:160:CYS:CB	4:D:216:ARG:NH1	2.56	0.47
1:A:97:VAL:O	1:A:97:VAL:HG23	2.15	0.47
4:D:57:LEU:C	4:D:57:LEU:HD23	2.39	0.47
1:A:118:ASN:HB2	1:A:166:GLU:HB2	1.97	0.47
2:B:125:SER:CB	2:B:147:LEU:HD21	2.44	0.47
3:C:164:MET:HE2	3:C:170:LYS:HD3	1.96	0.47
2:B:30:HIS:HB2	2:B:38:VAL:HG12	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:ARG:HH21	4:D:17:ARG:HB2	1.80	0.46
2:B:126:GLN:CD	2:B:126:GLN:N	2.62	0.46
1:A:45:TRP:HZ3	6:A:208:HOH:O	1.99	0.46
3:C:155:TYR:O	3:C:176:ALA:HA	2.15	0.46
3:C:21:ILE:CG1	3:C:106:THR:HG21	2.42	0.46
1:A:52:GLY:HA2	4:D:5:PRO:O	2.16	0.46
4:D:4:ASN:O	4:D:5:PRO:C	2.59	0.46
1:A:99:LEU:H	1:A:99:LEU:HD12	1.81	0.46
3:C:13:VAL:HG22	3:C:110:ILE:CD1	2.46	0.46
3:C:156:ILE:HD12	3:C:156:ILE:C	2.41	0.45
2:B:163:MET:HG3	2:B:171:TYR:CZ	2.51	0.45
3:C:164:MET:HA	3:C:169:PHE:O	2.16	0.45
3:C:123:GLN:O	4:D:154:SER:HB2	2.17	0.45
2:B:6:ASP:OD2	2:B:7:PHE:N	2.46	0.45
4:D:45:GLU:HG2	4:D:103:LYS:HG2	1.98	0.45
1:A:50:LYS:HE2	2:B:89:GLY:HA3	1.99	0.44
4:D:142:ASN:CG	4:D:142:ASN:O	2.59	0.44
4:D:176:ASP:HB3	4:D:211:TYR:CD2	2.53	0.44
2:B:116:ILE:HG12	2:B:160:MET:HG3	1.98	0.44
3:C:40:LEU:H	3:C:40:LEU:CD2	2.27	0.44
1:A:32:PHE:C	1:A:32:PHE:CD1	2.95	0.44
2:B:134:ASN:O	2:B:135:ASP:HB2	2.17	0.44
3:C:2:GLU:O	3:C:3:ASN:CB	2.66	0.44
3:C:189:PHE:O	3:C:189:PHE:CG	2.70	0.44
1:A:23:THR:HG22	1:A:33:TYR:CB	2.48	0.44
2:B:49:ALA:HB2	2:B:55:ARG:HA	2.00	0.44
3:C:163:ASP:C	3:C:165:ARG:N	2.75	0.44
1:A:132:VAL:CG1	1:A:133:SER:N	2.81	0.43
4:D:50:PRO:HB3	4:D:119:THR:HG21	2.01	0.43
4:D:64:LEU:HD22	4:D:65:GLY:N	2.21	0.43
3:C:153:ASP:HB2	3:C:180:LYS:HE2	2.00	0.43
1:A:7:SER:HB2	1:A:25:GLU:HB2	2.00	0.43
3:C:52:SER:O	3:C:68:LYS:HD2	2.19	0.43
1:A:95:SER:HB2	1:A:96:PRO:CD	2.49	0.43
3:C:81:GLN:HB3	3:C:83:GLU:OE1	2.18	0.43
3:C:151:ASP:HB2	3:C:154:VAL:HG12	2.00	0.43
2:B:167:ARG:HH11	2:B:167:ARG:HG3	1.84	0.42
3:C:89:PHE:CE1	3:C:105:GLY:HA3	2.54	0.42
2:B:39:ARG:HG2	2:B:40:PHE:N	2.32	0.42
4:D:39:GLU:HA	4:D:137:LEU:O	2.19	0.42
2:B:124:PRO:CG	2:B:126:GLN:OE1	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:152:ASP:O	2:B:153:TRP:HB2	2.20	0.42
1:A:45:TRP:C	1:A:47:GLU:H	2.28	0.41
1:A:61:ARG:HA	4:D:80:ALA:HB1	2.01	0.41
2:B:65:LYS:O	2:B:69:GLU:HG3	2.20	0.41
3:C:151:ASP:HB2	3:C:154:VAL:CG1	2.49	0.41
3:C:40:LEU:HD22	3:C:40:LEU:N	2.32	0.41
4:D:4:ASN:OD1	4:D:5:PRO:HD2	2.18	0.41
2:B:61:TRP:HB3	2:B:67:VAL:HG11	2.02	0.41
1:A:18:PRO:O	1:A:19:SER:HB2	2.21	0.41
4:D:15:THR:OG1	4:D:16:PRO:HD2	2.21	0.41
3:C:73:PHE:CE1	3:C:90:CYS:HB2	2.55	0.41
3:C:169:PHE:CD2	4:D:163:LYS:HE2	2.56	0.41
1:A:98:THR:O	1:A:99:LEU:C	2.63	0.41
3:C:45:GLN:O	3:C:45:GLN:HG3	2.21	0.41
3:C:21:ILE:HD12	3:C:21:ILE:N	2.36	0.41
4:D:64:LEU:HA	4:D:64:LEU:HD23	1.63	0.41
4:D:105:GLN:O	4:D:106:ARG:C	2.62	0.41
4:D:194:CYS:O	4:D:215:SER:HA	2.21	0.41
1:A:98:THR:HG22	1:A:101:GLN:HB2	2.02	0.41
4:D:87:ASN:HD21	4:D:89:ARG:NH1	2.18	0.41
2:B:152:ASP:OD1	2:B:154:THR:OG1	2.38	0.40
6:C:231:HOH:O	4:D:64:LEU:HD21	2.20	0.40
3:C:32:TYR:HB3	3:C:93:SER:HB3	2.03	0.40
1:A:117:VAL:HG12	1:A:137:PHE:CZ	2.57	0.40
1:A:7:SER:HB3	1:A:10:VAL:CG2	2.51	0.40
4:D:246:TRP:CZ2	4:D:248:GLN:HB2	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	180/194 (93%)	166 (92%)	10 (6%)	4 (2%)	5	5
2	B	177/202 (88%)	170 (96%)	7 (4%)	0	100	100
3	C	189/206 (92%)	172 (91%)	12 (6%)	5 (3%)	4	3
4	D	253/268 (94%)	238 (94%)	11 (4%)	4 (2%)	7	9
All	All	799/870 (92%)	746 (93%)	40 (5%)	13 (2%)	7	9

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	50	LYS
4	D	5	PRO
4	D	266	ALA
3	C	3	ASN
3	C	164	MET
1	A	9	GLY
3	C	152	SER
3	C	185	CYS
3	C	190	ASN
1	A	46	PRO
1	A	51	PHE
4	D	187	LYS
4	D	95	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	157/173 (91%)	147 (94%)	10 (6%)	16	23
2	B	163/182 (90%)	158 (97%)	5 (3%)	35	53
3	C	170/187 (91%)	155 (91%)	15 (9%)	9	12
4	D	212/224 (95%)	191 (90%)	21 (10%)	7	9
All	All	702/766 (92%)	651 (93%)	51 (7%)	13	18

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	0	ILE
1	A	1	VAL
1	A	8	CYS
1	A	47	GLU
1	A	62	ASN
1	A	88	GLU
1	A	98	THR
1	A	101	GLN
1	A	129	THR
1	A	130	GLU
2	B	39	ARG
2	B	100	THR
2	B	159	VAL
2	B	163	MET
2	B	164	THR
3	C	11	LEU
3	C	12	SER
3	C	13	VAL
3	C	69	THR
3	C	150	LYS
3	C	154	VAL
3	C	156	ILE
3	C	161	VAL
3	C	163	ASP
3	C	164	MET
3	C	170	LYS
3	C	187	ASN
3	C	189	PHE
3	C	190	ASN
3	C	191	ASN
4	D	5	PRO
4	D	6	VAL
4	D	29	VAL
4	D	47	ARG
4	D	64	LEU
4	D	71	LEU
4	D	77	THR
4	D	105	GLN
4	D	138	GLU
4	D	154	SER
4	D	169	LEU
4	D	180	LEU
4	D	194	CYS

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Mol	Chain	Res	Type
4	D	198	GLN
4	D	203	GLN
4	D	216	ARG
4	D	228	ARG
4	D	240	LEU
4	D	242	GLU
4	D	249	ASP
4	D	267	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	GLN
2	B	113	ASN
2	B	166	GLN
3	C	6	GLN
3	C	7	HIS
3	C	14	GLN
3	C	81	GLN
3	C	97	ASN
3	C	123	GLN
3	C	145	ASN
4	D	177	HIS

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](#)

2 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	$\# Z > 2$
5	NAG	B	5467	2	14,14,15	0.65	0	17,19,21	0.69	0
5	NAG	A	1181	1	14,14,15	0.89	1 (7%)	17,19,21	0.70	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
5	NAG	B	5467	2	-	0/6/23/26	0/1/1/1
5	NAG	A	1181	1	-	4/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1181	NAG	C1-C2	2.63	1.55	1.52

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1181	NAG	C8-C7-N2-C2
5	A	1181	NAG	O7-C7-N2-C2
5	A	1181	NAG	O5-C5-C6-O6
5	A	1181	NAG	C4-C5-C6-O6

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	182/194 (93%)	0.70	16 (8%) 15 15	39, 59, 92, 99	0
2	B	181/202 (89%)	0.27	4 (2%) 62 63	40, 56, 69, 89	0
3	C	191/206 (92%)	0.56	17 (8%) 15 14	32, 55, 80, 92	0
4	D	257/268 (95%)	0.38	12 (4%) 36 36	38, 54, 75, 91	0
All	All	811/870 (93%)	0.47	49 (6%) 27 27	32, 56, 83, 99	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	188	ALA	7.2
1	A	46	PRO	6.0
1	A	48	PHE	5.0
1	A	51	PHE	5.0
3	C	187	ASN	4.8
3	C	186	ALA	4.7
3	C	190	ASN	4.7
3	C	189	PHE	4.6
3	C	192	SER	4.4
3	C	164	MET	4.3
3	C	191	ASN	4.3
3	C	163	ASP	4.2
4	D	64	LEU	4.0
4	D	17	ARG	3.9
1	A	154	LEU	3.9
2	B	191	GLN	3.8
1	A	50	LYS	3.5
1	A	160	ILE	3.5
1	A	45	TRP	3.5
1	A	131	GLY	3.4
1	A	129	THR	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	49	SER	3.2
1	A	128	VAL	3.1
1	A	125	GLY	3.1
3	C	166	SER	3.0
2	B	190	ALA	3.0
1	A	52	GLY	2.9
3	C	151	ASP	2.9
3	C	24	THR	2.8
4	D	247	THR	2.7
1	A	180	PRO	2.7
3	C	184	ALA	2.7
1	A	47	GLU	2.5
4	D	243	ASN	2.5
4	D	177	HIS	2.5
4	D	242	GLU	2.4
3	C	185	CYS	2.4
4	D	194	CYS	2.4
3	C	150	LYS	2.2
4	D	267	ASP	2.2
3	C	68	LYS	2.2
2	B	136	GLN	2.2
4	D	203	GLN	2.1
2	B	23	ARG	2.0
3	C	165	ARG	2.0
4	D	206	LEU	2.0
4	D	5	PRO	2.0
1	A	124	ASN	2.0
4	D	4	ASN	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($\text{\AA}^2$)	Q<0.9
5	NAG	B	5467	14/15	0.67	0.17	90,93,95,95	0
5	NAG	A	1181	14/15	0.81	0.14	81,82,83,83	0

6.5 Other polymers [i](#)

There are no such residues in this entry.