



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2020 – 04:51 PM BST

PDB ID : 2SEB
Title : X-RAY CRYSTAL STRUCTURE OF HLA-DR4 COMPLEXED WITH A PEPTIDE FROM HUMAN COLLAGEN II
Authors : Dessen, A.; Lawrence, C.M.; Cupo, S.; Zaller, D.M.; Wiley, D.C.
Deposited on : 1997-10-16
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

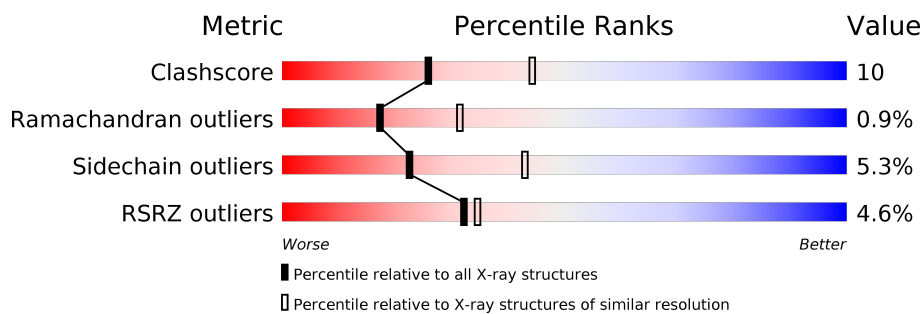
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on 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	181	<div> <div>2%</div> <div>75%</div> <div>23%</div> <div>..</div> </div>
2	B	192	<div> <div>4%</div> <div>61%</div> <div>27%</div> <div>8%</div> </div>
3	D	239	<div> <div>6%</div> <div>71%</div> <div>18%</div> <div>9%</div> </div>
4	E	12	<div> <div>8%</div> <div>92%</div> <div>8%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4857 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 II HISTOCOMPATIBILITY ANTIGEN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	180	Total	C	N	O	S	0	0	0
			1469	954	238	272	5			

- Molecule 2 is a protein called HLA CLASS II HISTOCOMPATIBILITY ANTIGEN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	177	Total	C	N	O	S	19	0	0
			1451	920	253	273	5			

- Molecule 3 is a protein called ENTEROTOXIN TYPE B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	217	Total	C	N	O	S	0	0	0
			1784	1150	287	337	10			

- Molecule 4 is a protein called PEPTIDE FROM COLLAGEN II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	12	Total	C	N	O	S	0	0	0
			77	46	15	15	1			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	1174	ALA	GLN	conflict	UNP P02458

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	15	Total	O	0	0
			15	15		
6	B	19	Total	O	0	0
			19	19		
6	D	20	Total	O	0	0
			20	20		
6	E	8	Total	O	0	0
			8	8		



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.53 Å 100.28 Å 100.87 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	12.00 – 2.50 11.94 – 2.50	Depositor EDS
% Data completeness (in resolution range)	87.7 (12.00-2.50) 97.6 (11.94-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.74 (at 2.49 Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.229 , 0.280 0.239 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	24.9	Xtriage
Anisotropy	0.245	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.021 for -h,l,k	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	4857	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% 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	0.54	2/1514 (0.1%)	0.70	0/2065
2	B	0.45	0/1489	0.73	1/2025 (0.0%)
3	D	0.46	0/1826	0.68	1/2461 (0.0%)
4	E	0.47	0/77	0.81	0/102
All	All	0.48	2/4906 (0.0%)	0.70	2/6653 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	127	PRO	N-CD	-12.40	1.30	1.47
1	A	127	PRO	N-CA	5.36	1.56	1.47

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	23	ARG	O-C-N	-7.24	111.12	122.70
3	D	37	ASN	N-CA-CB	6.46	122.24	110.60

There are no chirality outliers.

There are no planarity outliers.

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	1469	0	1404	28	0
2	B	1451	0	1339	42	0
3	D	1784	0	1694	36	0
4	E	77	0	70	1	0
5	A	14	0	13	0	0
6	A	15	0	0	1	0
6	B	19	0	0	1	0
6	D	20	0	0	0	0
6	E	8	0	0	0	0
All	All	4857	0	4520	97	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 10.

All (97) 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)
2:B:27:LEU:HD23	2:B:29:ARG:HD2	1.40	1.03
3:D:126:LEU:HD23	3:D:153:LYS:HE2	1.44	0.97
3:D:126:LEU:HD21	3:D:153:LYS:HD3	1.58	0.85
2:B:7:PHE:HA	2:B:33:HIS:HE1	1.46	0.81
3:D:126:LEU:HD23	3:D:153:LYS:CE	2.12	0.80
2:B:37:TYR:CD2	2:B:38:VAL:HG22	2.23	0.74
2:B:121:GLY:HA2	2:B:154:THR:HG23	1.71	0.73
3:D:126:LEU:CD2	3:D:153:LYS:CE	2.70	0.69
2:B:119:VAL:HB	2:B:157:THR:HG22	1.74	0.68
3:D:126:LEU:CD2	3:D:153:LYS:HE2	2.24	0.66
1:A:118:ASN:HB2	1:A:166:GLU:HB2	1.77	0.65
3:D:41:ILE:O	3:D:42:ASP:HB3	1.97	0.64
1:A:92:LEU:HD12	1:A:92:LEU:N	2.12	0.62
2:B:7:PHE:HA	2:B:33:HIS:CE1	2.32	0.62
2:B:31:PHE:O	2:B:32:TYR:O	2.18	0.61
3:D:36:ILE:O	3:D:36:ILE:HG22	1.99	0.61
2:B:93:ARG:HG2	2:B:123:TYR:CD1	2.35	0.61
2:B:18:PHE:HB2	2:B:23:ARG:HB3	1.81	0.61
3:D:95:PHE:O	3:D:111:LYS:NZ	2.22	0.61
3:D:176:GLU:HB2	3:D:179:ASN:O	2.01	0.59
2:B:113:ASN:O	2:B:163:THR:HB	2.02	0.59
2:B:138:GLU:HG2	2:B:161:LEU:HD11	1.84	0.58
1:A:92:LEU:HD12	1:A:92:LEU:H	1.69	0.58
3:D:126:LEU:HD21	3:D:153:LYS:CD	2.31	0.57
1:A:98:GLU:HB2	1:A:101:GLU:HG3	1.84	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:163:THR:HG23	2:B:171:TYR:OH	2.05	0.57
2:B:75:VAL:HG23	6:B:920:HOH:O	2.05	0.57
2:B:117:CYS:HB2	2:B:131:TRP:CZ2	2.39	0.56
1:A:9:GLN:HG3	1:A:24:PHE:CE1	2.41	0.55
2:B:128:GLU:HB2	2:B:176:GLU:HB2	1.88	0.55
2:B:10:GLN:HB3	2:B:31:PHE:HB2	1.87	0.55
2:B:27:LEU:HD23	2:B:29:ARG:CD	2.25	0.55
3:D:185:GLY:HA2	3:D:233:TYR:O	2.06	0.54
3:D:80:LYS:HD3	3:D:120:GLU:HG3	1.89	0.54
1:A:30:GLU:OE2	6:A:957:HOH:O	2.18	0.53
2:B:180:LEU:HD13	2:B:184:LEU:HG	1.90	0.53
1:A:129:THR:O	1:A:132:VAL:HG22	2.08	0.53
2:B:66:ASP:O	2:B:70:GLN:HG3	2.09	0.53
2:B:121:GLY:HA2	2:B:154:THR:CG2	2.39	0.52
2:B:145:THR:HG22	2:B:158:LEU:H	1.73	0.52
3:D:130:ARG:NH1	3:D:226:LYS:HB2	2.24	0.52
2:B:97:PRO:HB3	2:B:122:PHE:HB3	1.93	0.51
1:A:143:HIS:HD2	2:B:12:LYS:NZ	2.08	0.51
3:D:224:ASP:O	3:D:228:VAL:HG22	2.10	0.51
2:B:150:ASN:HD22	2:B:154:THR:HG22	1.75	0.50
1:A:95:SER:HB2	1:A:96:PRO:HD2	1.94	0.50
3:D:124:ASN:HD21	3:D:152:LYS:HB3	1.77	0.49
2:B:27:LEU:HD12	2:B:41:ASP:HA	1.93	0.49
3:D:148:VAL:HG13	3:D:163:LEU:HD13	1.94	0.49
1:A:111:LYS:HG2	1:A:140:ARG:CZ	2.42	0.49
1:A:57:GLN:HB2	3:D:92:GLN:CD	2.33	0.48
2:B:21:THR:O	2:B:80:ARG:NH1	2.46	0.48
3:D:95:PHE:CE1	3:D:111:LYS:HB3	2.49	0.48
1:A:16:PRO:HD2	2:B:6:ARG:HD3	1.96	0.47
3:D:165:ARG:HG2	3:D:200:MET:HE3	1.95	0.47
3:D:124:ASN:O	3:D:152:LYS:HA	2.16	0.46
3:D:126:LEU:CD2	3:D:153:LYS:NZ	2.79	0.46
2:B:177:HIS:CD2	2:B:178:PRO:HD2	2.51	0.46
2:B:90:THR:HG22	2:B:153:TRP:CH2	2.50	0.46
1:A:73:MET:SD	1:A:76:ARG:NH2	2.88	0.46
3:D:17:PHE:CZ	3:D:208:PHE:HB2	2.51	0.45
3:D:37:ASN:HD22	3:D:37:ASN:HA	1.57	0.45
2:B:30:TYR:HB2	2:B:38:VAL:HG23	1.98	0.45
3:D:93:CYS:SG	3:D:112:THR:C	2.95	0.45
2:B:28:ASP:HB3	2:B:40:PHE:HB3	1.97	0.45
1:A:14:LEU:HD11	2:B:6:ARG:HB3	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57:GLN:HE21	3:D:92:GLN:HA	1.82	0.45
2:B:37:TYR:CD2	2:B:38:VAL:CG2	2.98	0.44
2:B:128:GLU:O	2:B:175:VAL:HA	2.18	0.44
1:A:11:GLU:HA	1:A:21:GLU:O	2.17	0.44
2:B:28:ASP:O	2:B:39:ARG:HA	2.18	0.44
1:A:30:GLU:O	1:A:44:ARG:HB2	2.18	0.44
3:D:124:ASN:ND2	3:D:152:LYS:HB3	2.33	0.44
3:D:134:VAL:HA	3:D:230:ILE:O	2.18	0.43
3:D:134:VAL:HG22	3:D:230:ILE:HB	2.00	0.43
1:A:7:ILE:HG12	1:A:26:PHE:HD1	1.82	0.43
2:B:4:ARG:HA	2:B:5:PRO:HD3	1.81	0.43
1:A:107:CYS:HB2	1:A:121:TRP:CH2	2.53	0.42
1:A:86:PRO:HA	1:A:87:PRO:HD3	1.86	0.42
3:D:92:GLN:O	3:D:92:GLN:HG2	2.19	0.42
1:A:39:LYS:HD3	3:D:92:GLN:NE2	2.34	0.42
3:D:134:VAL:HG11	3:D:164:THR:HG23	2.01	0.42
1:A:34:VAL:HG21	1:A:59:ALA:HB2	2.01	0.42
3:D:136:VAL:O	3:D:142:ASN:HA	2.19	0.42
1:A:16:PRO:HD2	2:B:6:ARG:CD	2.50	0.41
1:A:94:ASN:HB2	1:A:104:VAL:HB	2.02	0.41
3:D:183:GLU:HB3	3:D:235:THR:HG22	2.02	0.41
1:A:136:VAL:HG13	1:A:137:PHE:N	2.35	0.41
2:B:85:VAL:HG21	4:E:1169:TYR:O	2.21	0.41
2:B:127:ILE:HD11	2:B:175:VAL:CG1	2.51	0.41
2:B:30:TYR:N	2:B:30:TYR:CD1	2.89	0.41
3:D:11:LEU:HD11	3:D:186:TYR:CD1	2.56	0.41
3:D:73:LEU:HA	3:D:73:LEU:HD12	1.89	0.41
1:A:71:GLU:O	1:A:75:LYS:HG3	2.22	0.40
3:D:80:LYS:HD3	3:D:120:GLU:CG	2.52	0.40
1:A:9:GLN:HB3	2:B:13:HIS:HB2	2.03	0.40
1:A:143:HIS:CE1	2:B:31:PHE:HE2	2.40	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	178/181 (98%)	173 (97%)	5 (3%)	0	100	100
2	B	171/192 (89%)	163 (95%)	6 (4%)	2 (1%)	13	24
3	D	211/239 (88%)	195 (92%)	13 (6%)	3 (1%)	11	20
4	E	10/12 (83%)	10 (100%)	0	0	100	100
All	All	570/624 (91%)	541 (95%)	24 (4%)	5 (1%)	17	31

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	32	TYR
3	D	37	ASN
2	B	87	GLU
3	D	91	TYR
3	D	42	ASP

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	162/166 (98%)	155 (96%)	7 (4%)	29	53
2	B	154/173 (89%)	146 (95%)	8 (5%)	23	44
3	D	193/225 (86%)	181 (94%)	12 (6%)	18	35
4	E	4/4 (100%)	4 (100%)	0	100	100
All	All	513/568 (90%)	486 (95%)	27 (5%)	22	43

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	92	LEU
1	A	113	THR
1	A	117	VAL

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Mol	Chain	Res	Type
1	A	128	VAL
1	A	130	THR
1	A	136	VAL
1	A	159	ASP
2	B	2	ASP
2	B	6	ARG
2	B	38	VAL
2	B	39	ARG
2	B	98	GLU
2	B	138	GLU
2	B	154	THR
2	B	172	THR
3	D	5	ASP
3	D	27	LEU
3	D	37	ASN
3	D	38	VAL
3	D	45	LEU
3	D	73	LEU
3	D	78	LYS
3	D	119	THR
3	D	120	GLU
3	D	131	SER
3	D	149	GLN
3	D	228	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	ASN
1	A	143	HIS
1	A	149	HIS
2	B	13	HIS
3	D	3	GLN
3	D	23	ASN
3	D	37	ASN
3	D	125	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	A	900	1	14,14,15	1.09	0	17,19,21	1.49	4 (23%)

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	A	900	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	900	NAG	C4-C3-C2	-4.08	105.04	111.02
5	A	900	NAG	O4-C4-C3	2.57	116.29	110.35
5	A	900	NAG	C2-N2-C7	-2.19	119.79	122.90
5	A	900	NAG	C6-C5-C4	-2.05	108.20	113.00

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	180/181 (99%)	-0.02	3 (1%) 70 72	4, 19, 39, 51	0
2	B	177/192 (92%)	0.21	8 (4%) 33 36	3, 19, 38, 55	4 (2%)
3	D	217/239 (90%)	0.21	15 (6%) 16 17	5, 19, 40, 53	0
4	E	12/12 (100%)	0.09	1 (8%) 11 11	13, 21, 37, 67	0
All	All	586/624 (93%)	0.14	27 (4%) 32 34	3, 19, 40, 67	4 (0%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	190	ALA	6.8
3	D	110	ARG	6.5
2	B	3	THR	4.3
3	D	9	ASP	4.0
2	B	66	ASP	3.9
1	A	1	ILE	3.4
3	D	62	ASP	3.2
2	B	59	GLU	3.2
3	D	79	ASP	3.1
4	E	1179	ALA	3.1
3	D	206	ASP	3.0
3	D	31	ASN	3.0
2	B	140	THR	2.8
3	D	81	TYR	2.5
3	D	61	TYR	2.5
3	D	37	ASN	2.5
2	B	164	VAL	2.4
3	D	36	ILE	2.2
1	A	126	LYS	2.2
3	D	141	LYS	2.1
3	D	8	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
2	B	134	ASN	2.1
3	D	30	ASP	2.1
3	D	219	ASP	2.1
3	D	2	SER	2.0
2	B	113	ASN	2.0
1	A	71	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	NAG	A	900	14/15	0.60	0.36	55,70,81,84	0

6.5 Other polymers [i](#)

There are no such residues in this entry.