



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 5, 2022 – 04:18 PM EST

PDB ID : 7K7R
Title : EBNA1 peptide AA386-405 with Fab MS39p2w174
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Deposited on : 2020-09-23
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 : 2.25
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) : 2.25

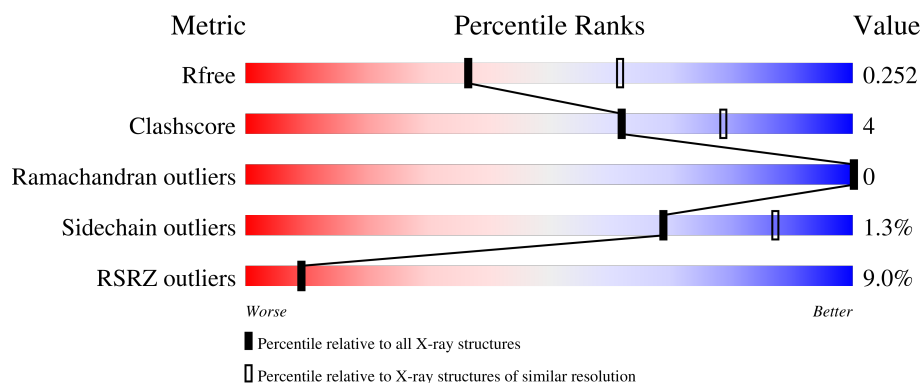
1 Overall quality at a glance

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))
R_{free}	130704	4661 (2.50-2.50)
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 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	219	<div> <div>17%</div> <div>91% 8%</div> </div>
1	D	219	<div> <div>14%</div> <div>82% 17%</div> </div>
2	B	224	<div> <div>14%</div> <div>87% 8% 5%</div> </div>
2	E	224	<div> <div>5%</div> <div>60% 10% 30%</div> </div>
3	C	20	

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Mol	Chain	Length	Quality of chain
3	F	20	<div><div></div><div></div><div></div><div></div></div>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6827 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 Fab LC MS39p2w174.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	0	0
			1686	1056	291	334	5			
1	D	217	Total	C	N	O	S	0	0	0
			1672	1047	286	334	5			

- Molecule 2 is a protein called Fab HC MS39p2w174.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	212	Total	C	N	O	S	0	0	0
			1596	1009	268	314	5			
2	E	219	Total	C	N	O	S	0	0	0
			1645	1036	277	326	6			

- Molecule 3 is a protein called EBNA1 peptide AA386-405.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	14	Total	C	N	O	0	0	0
			94	56	20	18			
3	F	12	Total	C	N	O	0	0	0
			85	51	18	16			

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	1	Total	Cl	0	0
			1	1		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	E	1	Total	C	O	0	0
			6	3	3		
5	E	1	Total	C	O	0	0
			6	3	3		
5	E	1	Total	C	O	0	0
			6	3	3		
5	E	1	Total	C	O	0	0
			6	3	3		

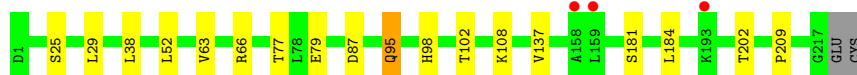
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	9	Total	O	0	0
			9	9		
6	B	6	Total	O	0	0
			6	6		
6	D	4	Total	O	0	0
			4	4		
6	E	4	Total	O	0	0
			4	4		
6	F	1	Total	O	0	0
			1	1		

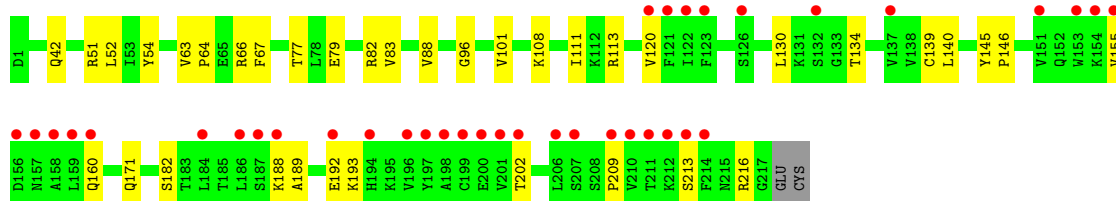
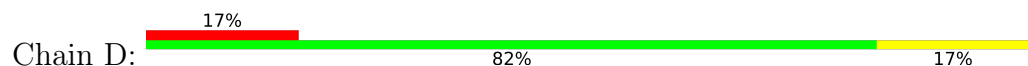
3 Residue-property plots [i](#)

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.

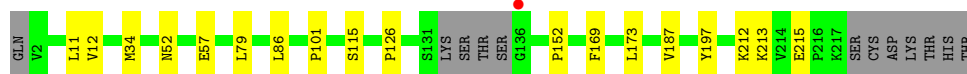
- Molecule 1: Fab LC MS39p2w174



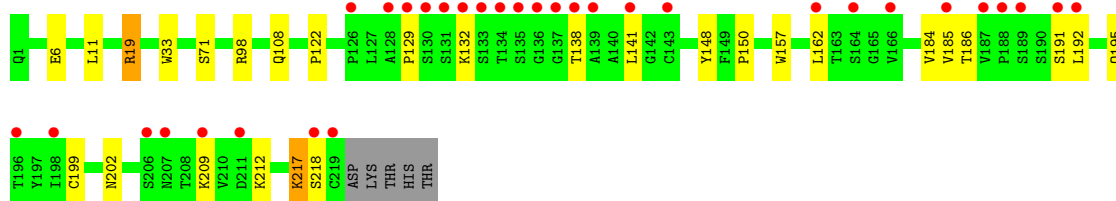
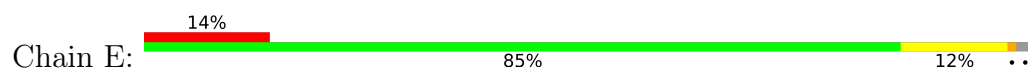
- Molecule 1: Fab LC MS39p2w174



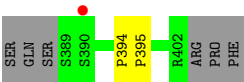
- Molecule 2: Fab HC MS39p2w174



- Molecule 2: Fab HC MS39p2w174



- Molecule 3: EBNA1 peptide AA386-405



● Molecule 3: EBNA1 peptide AA386-405



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	119.66Å 137.56Å 179.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.14 – 2.50 45.14 – 2.50	Depositor EDS
% Data completeness (in resolution range)	71.3 (45.14-2.50) 65.1 (45.14-2.50)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.211 , 0.252 0.211 , 0.252	Depositor DCC
R_{free} test set	1917 reflections (5.22%)	wwPDB-VP
Wilson B-factor (Å ²)	32.2	Xtriage
Anisotropy	0.716	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 38.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	6827	wwPDB-VP
Average B, all atoms (Å ²)	56.0	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: CL, GOL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.33	0/1723	0.54	0/2337
1	D	0.31	0/1709	0.52	0/2322
2	B	0.35	0/1636	0.57	0/2229
2	E	0.34	0/1686	0.57	0/2297
3	C	0.35	0/98	0.50	0/134
3	F	0.36	0/89	0.50	0/122
All	All	0.33	0/6941	0.55	0/9441

There are no bond length outliers.

There are no bond angle outliers.

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	1686	0	1651	12	0
1	D	1672	0	1618	19	0
2	B	1596	0	1549	11	0
2	E	1645	0	1601	19	0
3	C	94	0	88	2	0
3	F	85	0	83	1	0
4	D	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	E	24	0	32	3	0
6	A	9	0	0	0	0
6	B	6	0	0	1	0
6	D	4	0	0	0	0
6	E	4	0	0	0	0
6	F	1	0	0	0	0
All	All	6827	0	6622	59	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 (59) 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:E:191:SER:HB2	2:E:195:GLN:HB3	1.58	0.83
1:D:202:THR:HG22	1:D:209:PRO:HB3	1.73	0.70
2:E:11:LEU:HB2	2:E:150:PRO:HG3	1.73	0.69
2:B:12:VAL:HG11	2:B:86:LEU:HD13	1.78	0.66
1:D:140:LEU:HD22	2:E:184:VAL:HG11	1.78	0.64
2:E:162:LEU:HD21	2:E:185:VAL:HG11	1.82	0.61
2:E:217:LYS:HD2	2:E:218:SER:H	1.68	0.59
1:A:66:ARG:NH1	1:A:87:ASP:OD2	2.30	0.59
1:A:137:VAL:HG13	1:A:184:LEU:HB3	1.85	0.58
1:D:52:LEU:HA	1:D:63:VAL:HG21	1.85	0.58
1:A:29:LEU:HD11	1:A:95:GLN:HG3	1.85	0.58
2:E:157:TRP:CH2	2:E:199:CYS:HB3	2.39	0.57
1:A:95:GLN:HE21	1:A:98:HIS:H	1.53	0.55
2:E:19:ARG:NH1	5:E:304:GOL:O3	2.40	0.55
2:E:122:PRO:HB3	2:E:148:TYR:HB3	1.89	0.53
2:E:129:PRO:HG3	2:E:141:LEU:HD12	1.91	0.52
2:E:98:ARG:HH12	5:E:303:GOL:H2	1.76	0.50
1:A:66:ARG:HH12	1:A:87:ASP:CG	2.14	0.50
1:D:213:SER:H	2:E:132:LYS:NZ	2.09	0.50
2:E:138:THR:HG21	2:E:186:THR:HB	1.92	0.50
1:A:77:THR:HG22	1:A:79:GLU:HG3	1.94	0.50
2:B:126:PRO:HD3	2:B:212:LYS:HE2	1.94	0.48
1:D:77:THR:HG22	1:D:79:GLU:HG3	1.94	0.48
2:B:52:ASN:HB2	2:B:57:GLU:HB2	1.95	0.48
1:D:64:PRO:HG2	1:D:67:PHE:CD2	2.48	0.48
2:B:11:LEU:HD11	2:B:115:SER:HB3	1.96	0.48
1:D:130:LEU:HD22	1:D:188:LYS:HG3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:96:GLY:HA2	1:D:101:VAL:HG22	1.97	0.47
2:E:6:GLU:H	2:E:108:GLN:HE22	1.61	0.47
1:A:25:SER:OG	1:A:29:LEU:HD13	2.16	0.46
2:E:71:SER:HB3	5:E:304:GOL:H2	1.98	0.46
2:E:33:TRP:CD2	3:F:396:ARG:HD3	2.51	0.46
1:D:51:ARG:HD2	1:D:54:TYR:HB3	1.97	0.45
1:D:120:VAL:HA	1:D:140:LEU:O	2.16	0.45
1:D:42:GLN:HB2	1:D:52:LEU:HD11	1.99	0.45
1:A:181:SER:HB3	2:B:169:PHE:CE2	2.52	0.45
1:D:192:GLU:OE2	1:D:216:ARG:NH1	2.51	0.44
1:A:95:GLN:OE1	1:A:102:THR:OG1	2.34	0.44
2:B:187:VAL:HG11	2:B:197:TYR:CE1	2.52	0.44
1:D:155:VAL:HG22	1:D:160:GLN:NE2	2.34	0.43
2:B:213:LYS:HE3	2:B:215:GLU:OE1	2.19	0.43
1:A:202:THR:HG22	1:A:209:PRO:HB3	2.00	0.43
2:E:191:SER:HB2	2:E:195:GLN:CB	2.40	0.43
3:C:394:PRO:HA	3:C:395:PRO:HD3	1.86	0.43
2:E:11:LEU:HD22	2:E:150:PRO:HD3	2.00	0.43
2:E:129:PRO:HG3	2:E:192:LEU:HD22	2.00	0.43
1:D:139:CYS:HB3	1:D:182:SER:OG	2.19	0.42
1:D:145:TYR:CG	1:D:146:PRO:HA	2.54	0.42
1:D:66:ARG:HD2	1:D:82:ARG:O	2.20	0.42
1:A:66:ARG:H	1:A:66:ARG:HG2	1.71	0.42
1:D:88:VAL:HG21	1:D:171:GLN:HB3	2.01	0.42
2:E:202:ASN:OD1	2:E:209:LYS:HE2	2.21	0.41
2:B:101:PRO:HG3	3:C:395:PRO:HD2	2.03	0.41
2:B:152:PRO:HA	6:B:301:HOH:O	2.21	0.41
1:A:52:LEU:HA	1:A:63:VAL:HG21	2.03	0.40
1:D:189:ALA:O	1:D:193:LYS:HG3	2.22	0.40
2:B:34:MET:HB3	2:B:79:LEU:HD22	2.04	0.40
2:B:173:LEU:HD12	2:B:173:LEU:HA	1.94	0.40
1:D:83:VAL:HG12	1:D:111:ILE:HD11	2.03	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	215/219 (98%)	212 (99%)	3 (1%)	0	100	100
1	D	215/219 (98%)	212 (99%)	3 (1%)	0	100	100
2	B	208/224 (93%)	202 (97%)	6 (3%)	0	100	100
2	E	217/224 (97%)	211 (97%)	6 (3%)	0	100	100
3	C	12/20 (60%)	10 (83%)	2 (17%)	0	100	100
3	F	10/20 (50%)	8 (80%)	2 (20%)	0	100	100
All	All	877/926 (95%)	855 (98%)	22 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	193/195 (99%)	190 (98%)	3 (2%)	62	84
1	D	190/195 (97%)	187 (98%)	3 (2%)	62	84
2	B	181/193 (94%)	181 (100%)	0	100	100
2	E	188/193 (97%)	185 (98%)	3 (2%)	62	84
3	C	11/18 (61%)	11 (100%)	0	100	100
3	F	11/18 (61%)	10 (91%)	1 (9%)	9	18
All	All	774/812 (95%)	764 (99%)	10 (1%)	69	87

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	LEU
1	A	95	GLN
1	A	108	LYS

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Mol	Chain	Res	Type
1	D	108	LYS
1	D	113	ARG
1	D	134	THR
2	E	19	ARG
2	E	212	LYS
2	E	217	LYS
3	F	389	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 for Mogul analysis.

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	GOL	E	301	-	5,5,5	1.11	0	5,5,5	0.74	0
5	GOL	E	303	-	5,5,5	0.98	0	5,5,5	0.86	0
5	GOL	E	302	-	5,5,5	0.90	0	5,5,5	1.20	0
5	GOL	E	304	-	5,5,5	0.85	0	5,5,5	1.19	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	GOL	E	301	-	-	2/4/4/4	-
5	GOL	E	303	-	-	2/4/4/4	-
5	GOL	E	302	-	-	0/4/4/4	-
5	GOL	E	304	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	E	304	GOL	C1-C2-C3-O3
5	E	304	GOL	O2-C2-C3-O3
5	E	301	GOL	O1-C1-C2-C3
5	E	303	GOL	O1-C1-C2-C3
5	E	303	GOL	O1-C1-C2-O2
5	E	301	GOL	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	303	GOL	1	0
5	E	304	GOL	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	217/219 (99%)	-0.12	3 (1%) 75 77	29, 46, 68, 83	0
1	D	217/219 (99%)	0.69	37 (17%) 1 1	24, 56, 137, 151	0
2	B	212/224 (94%)	-0.09	1 (0%) 91 91	24, 41, 66, 104	0
2	E	219/224 (97%)	0.61	32 (14%) 2 2	21, 55, 134, 163	0
3	C	14/20 (70%)	0.33	1 (7%) 16 16	39, 55, 91, 92	0
3	F	12/20 (60%)	1.95	6 (50%) 0 0	46, 75, 114, 117	0
All	All	891/926 (96%)	0.30	80 (8%) 9 9	21, 46, 122, 163	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	155	VAL	8.4
1	D	198	ALA	7.9
1	D	154	LYS	7.6
2	E	192	LEU	7.2
2	E	139	ALA	7.1
3	F	389	SER	7.0
1	D	153	TRP	6.7
2	E	136	GLY	6.1
2	E	187	VAL	5.8
1	D	206	LEU	5.8
2	E	198	ILE	5.7
2	E	131	SER	5.5
1	D	196	VAL	5.4
2	E	135	SER	5.3
1	D	156	ASP	5.3
3	F	390	SER	4.8
2	E	134	THR	4.8
2	E	218	SER	4.7
1	D	211	THR	4.7

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Mol	Chain	Res	Type	RSRZ
1	D	122	ILE	4.7
2	E	211	ASP	4.6
1	D	186	LEU	4.5
1	D	123	PHE	4.4
1	D	199	CYS	4.4
1	D	157	ASN	4.3
1	D	187	SER	4.2
2	B	136	GLY	4.2
2	E	129	PRO	4.2
1	D	159	LEU	4.1
1	D	158	ALA	4.1
2	E	188	PRO	3.8
3	F	400	PRO	3.8
2	E	137	GLY	3.7
1	D	121	PHE	3.6
1	D	151	VAL	3.5
3	F	391	SER	3.4
2	E	138	THR	3.4
2	E	141	LEU	3.4
2	E	207	ASN	3.4
1	D	213	SER	3.3
2	E	164	SER	3.3
1	D	214	PHE	3.2
1	D	184	LEU	3.2
2	E	219	CYS	3.2
1	D	207	SER	3.1
1	D	209	PRO	3.0
1	D	120	VAL	2.8
2	E	128	ALA	2.8
1	D	197	TYR	2.8
3	F	392	GLY	2.8
1	A	159	LEU	2.7
1	D	201	VAL	2.7
1	D	212	LYS	2.7
2	E	133	SER	2.7
3	F	399	PRO	2.7
2	E	196	THR	2.6
1	D	160	GLN	2.6
1	D	202	THR	2.5
2	E	162	LEU	2.5
2	E	130	SER	2.5
2	E	189	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	200	GLU	2.4
1	A	193	LYS	2.4
2	E	132	LYS	2.3
2	E	166	VAL	2.3
2	E	126	PRO	2.3
2	E	143	CYS	2.3
1	D	188	LYS	2.2
2	E	185	VAL	2.2
2	E	209	LYS	2.2
1	D	126	SER	2.2
2	E	206	SER	2.2
1	D	210	VAL	2.2
1	A	158	ALA	2.2
1	D	192	GLU	2.1
1	D	132	SER	2.1
2	E	191	SER	2.1
3	C	390	SER	2.1
1	D	194	HIS	2.0
1	D	137	VAL	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	GOL	E	301	6/6	0.67	0.32	98,106,108,118	0
5	GOL	E	302	6/6	0.85	0.16	53,70,73,82	0
5	GOL	E	303	6/6	0.91	0.15	49,60,66,75	0
5	GOL	E	304	6/6	0.92	0.21	46,57,62,82	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	CL	D	301	1/1	0.96	0.06	45,45,45,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.