



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:51 AM UTC

PDB ID : 4FVS / pdb_00004fvs
Title : Crystal structure of a putative lipoprotein (BDI_3050) from Parabacteroides distasonis ATCC 8503 at 2.70 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2012-06-29
Resolution : 2.70 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

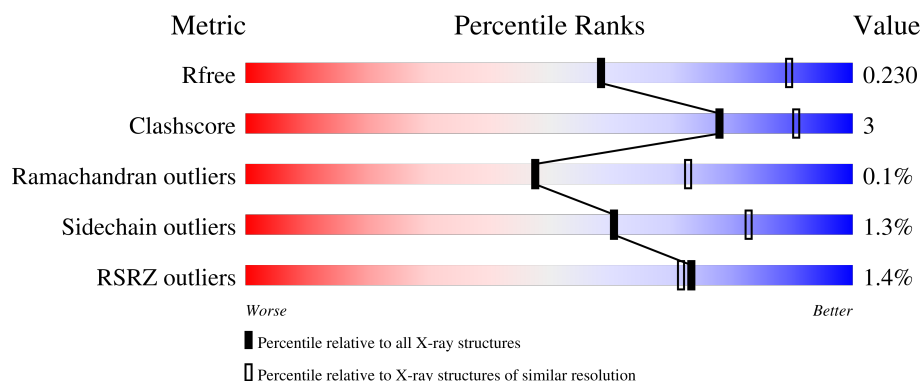
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	215	91% 9%
1	B	215	3% 88% 11%
1	C	215	87% 13%
1	D	215	91% 9%
1	E	215	88% 10%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	215	% 89% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10531 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 putative lipoprotein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	Se	0	1	0
			1719	1085	276	352	3	3			
1	B	214	Total	C	N	O	S	Se	0	1	0
			1715	1083	276	350	3	3			
1	C	214	Total	C	N	O	S	Se	0	1	0
			1711	1080	275	350	3	3			
1	D	214	Total	C	N	O	S	Se	0	0	0
			1722	1088	278	350	3	3			
1	E	213	Total	C	N	O	S	Se	0	0	0
			1705	1077	274	348	3	3			
1	F	214	Total	C	N	O	S	Se	0	1	0
			1716	1087	275	348	3	3			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP A6LGE3
B	0	GLY	-	expression tag	UNP A6LGE3
C	0	GLY	-	expression tag	UNP A6LGE3
D	0	GLY	-	expression tag	UNP A6LGE3
E	0	GLY	-	expression tag	UNP A6LGE3
F	0	GLY	-	expression tag	UNP A6LGE3

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	2	Total	Cl	0	0
			2	2		
2	D	1	Total	Cl	0	0
			1	1		
2	F	1	Total	Cl	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	34	Total 34	O 34	0	0
3	B	40	Total 40	O 40	0	0
3	C	40	Total 40	O 40	0	0
3	D	40	Total 40	O 40	0	0
3	E	52	Total 52	O 52	0	0
3	F	33	Total 33	O 33	0	0

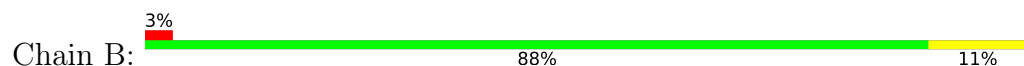
3 Residue-property plots

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.

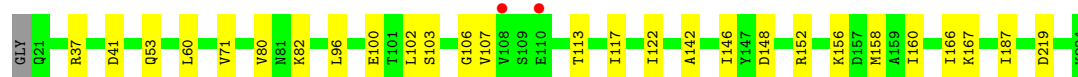
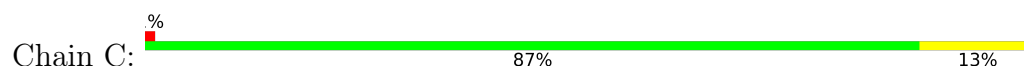
- Molecule 1: putative lipoprotein



- Molecule 1: putative lipoprotein



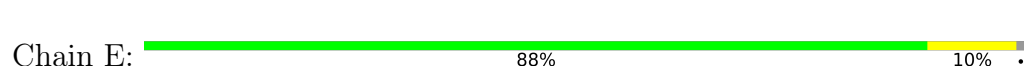
- Molecule 1: putative lipoprotein



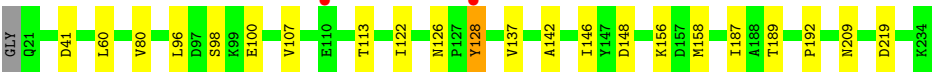
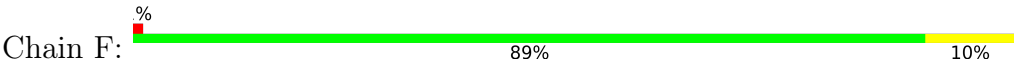
- Molecule 1: putative lipoprotein



- Molecule 1: putative lipoprotein



- Molecule 1: putative lipoprotein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	140.86Å 154.51Å 72.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.83 – 2.70 29.83 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.83-2.70) 99.6 (29.83-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.87 (at 2.68Å)	Xtriage
Refinement program	BUSTER-TNT 2.10.0, BUSTER 2.10.0	Depositor
R, R_{free}	0.187 , 0.224 0.197 , 0.230	Depositor DCC
R_{free} test set	2228 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	60.2	Xtriage
Anisotropy	0.526	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 49.6	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	0.95	EDS
Total number of atoms	10531	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.00% 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

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	0.82	1/1757 (0.1%)	1.15	6/2380 (0.3%)
1	B	0.86	1/1753 (0.1%)	1.13	7/2375 (0.3%)
1	C	0.82	0/1749	1.13	7/2371 (0.3%)
1	D	0.85	0/1757	1.11	4/2377 (0.2%)
1	E	0.87	0/1740	1.15	9/2357 (0.4%)
1	F	0.82	0/1755	1.13	8/2378 (0.3%)
All	All	0.84	2/10511 (0.0%)	1.13	41/14238 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	108	VAL	CA-C	6.87	1.61	1.52
1	A	90	GLN	CA-C	-5.25	1.46	1.52

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	148	ASP	CA-CB-CG	7.23	119.83	112.60
1	B	148	ASP	CA-CB-CG	7.17	119.77	112.60
1	E	148	ASP	CA-CB-CG	7.16	119.76	112.60
1	A	148	ASP	CA-CB-CG	6.99	119.59	112.60
1	C	148	ASP	CA-CB-CG	6.86	119.46	112.60
1	D	148	ASP	CA-CB-CG	6.71	119.31	112.60
1	B	140	ASP	CA-CB-CG	6.63	119.23	112.60
1	E	111	MSE	CA-C-N	6.56	132.86	122.36
1	E	111	MSE	C-N-CA	6.56	132.86	122.36
1	E	107	VAL	CA-C-N	6.14	128.88	120.46
1	E	107	VAL	C-N-CA	6.14	128.88	120.46
1	D	219	ASP	CA-CB-CG	5.99	118.58	112.60
1	B	111	MSE	N-CA-C	5.84	118.53	110.35

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	80	VAL	N-CA-C	-5.82	107.14	112.96
1	B	219	ASP	CA-CB-CG	5.77	118.37	112.60
1	A	219	ASP	CA-CB-CG	5.71	118.31	112.60
1	F	41	ASP	CA-CB-CG	5.51	118.11	112.60
1	B	41	ASP	CA-CB-CG	5.49	118.09	112.60
1	A	41	ASP	CA-CB-CG	5.42	118.02	112.60
1	F	219	ASP	CA-CB-CG	5.42	118.02	112.60
1	C	106	GLY	CA-C-N	5.41	127.38	120.56
1	C	106	GLY	C-N-CA	5.41	127.38	120.56
1	D	41	ASP	CA-CB-CG	5.40	118.00	112.60
1	C	219	ASP	CA-CB-CG	5.40	118.00	112.60
1	E	41	ASP	CA-CB-CG	5.31	117.91	112.60
1	C	41	ASP	CA-CB-CG	5.26	117.86	112.60
1	E	219	ASP	CA-CB-CG	5.25	117.86	112.60
1	E	140	ASP	CA-CB-CG	5.19	117.79	112.60
1	D	80	VAL	N-CA-C	-5.18	107.78	112.96
1	A	67	GLU	CB-CG-CD	5.11	121.29	112.60
1	B	161	LYS	CA-C-N	5.08	129.35	122.19
1	B	161	LYS	C-N-CA	5.08	129.35	122.19
1	A	130	ALA	CA-C-N	5.07	129.32	122.07
1	A	130	ALA	C-N-CA	5.07	129.32	122.07
1	C	107	VAL	CA-C-N	5.07	128.61	120.30
1	C	107	VAL	C-N-CA	5.07	128.61	120.30
1	F	192	PRO	CA-C-N	5.05	127.30	120.44
1	F	192	PRO	C-N-CA	5.05	127.30	120.44
1	F	107	VAL	CA-C-N	5.04	128.43	120.47
1	F	107	VAL	C-N-CA	5.04	128.43	120.47
1	E	133	ASP	CA-CB-CG	5.01	117.61	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	1719	0	1619	6	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1715	0	1615	13	0
1	C	1711	0	1604	14	0
1	D	1722	0	1637	10	0
1	E	1705	0	1607	11	0
1	F	1716	0	1616	10	0
2	C	2	0	0	1	0
2	D	1	0	0	0	0
2	F	1	0	0	0	0
3	A	34	0	0	0	0
3	B	40	0	0	0	0
3	C	40	0	0	0	0
3	D	40	0	0	0	0
3	E	52	0	0	0	0
3	F	33	0	0	0	0
All	All	10531	0	9698	55	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 3.

All (55) 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:B:158:MSE:HE2	1:B:187:ILE:HD11	1.56	0.88
1:E:158:MSE:HE2	1:E:187:ILE:HD11	1.78	0.64
1:D:158:MSE:HE2	1:D:187:ILE:HD11	1.82	0.59
1:A:158:MSE:HE2	1:A:187:ILE:HD11	1.83	0.59
1:B:158:MSE:HG2	1:B:189:THR:HG23	1.84	0.59
1:F:122:ILE:HG21	1:F:142:ALA:HB2	1.85	0.58
1:D:122:ILE:HG21	1:D:142:ALA:HB2	1.85	0.57
1:C:53:GLN:NE2	1:E:113:THR:OG1	2.34	0.56
1:B:98:SER:HB2	1:B:142:ALA:HB1	1.88	0.55
1:B:146:ILE:HG23	1:B:156:LYS:HB2	1.88	0.55
1:C:71:VAL:HG21	1:E:71:VAL:HG21	1.89	0.54
1:D:158:MSE:HG2	1:D:189:THR:HG23	1.88	0.54
1:E:96:LEU:HD21	1:E:100:GLU:HG2	1.90	0.54
1:F:158:MSE:HE2	1:F:187:ILE:HD11	1.90	0.54
1:B:96:LEU:HD21	1:B:100:GLU:HG2	1.90	0.54
1:A:96:LEU:HD21	1:A:100:GLU:HG2	1.89	0.54
1:D:223:VAL:CG2	1:F:209:ASN:ND2	2.70	0.54
1:F:96:LEU:HD21	1:F:100:GLU:HG2	1.91	0.53
1:B:122:ILE:HD13	1:B:142:ALA:HB2	1.90	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:142:ALA:HB3	1:C:160:ILE:HB	1.91	0.52
1:C:96:LEU:HD21	1:C:100:GLU:HG2	1.91	0.52
1:E:122:ILE:HG21	1:E:142:ALA:HB2	1.92	0.52
1:D:96:LEU:HD21	1:D:100:GLU:HG2	1.91	0.52
1:E:146:ILE:HG23	1:E:156:LYS:HB2	1.92	0.51
1:A:146:ILE:HG23	1:A:156:LYS:HB2	1.93	0.51
1:F:126:ASN:OD1	1:F:128[A]:TYR:HD2	1.93	0.51
1:C:146:ILE:HG23	1:C:156:LYS:HB2	1.93	0.50
1:B:227:TYR:OH	1:E:131:ASN:ND2	2.45	0.50
1:C:158:MSE:HE2	1:C:187:ILE:HD11	1.94	0.50
1:C:37:ARG:HH22	1:C:103:SER:HB2	1.77	0.49
1:A:108:VAL:O	1:A:111:MSE:HB2	2.12	0.49
1:B:108:VAL:HG22	1:B:111:MSE:HE2	1.95	0.48
1:F:126:ASN:OD1	1:F:128[A]:TYR:CD2	2.67	0.48
1:D:37:ARG:HH22	1:D:103:SER:HB2	1.78	0.48
1:F:146:ILE:HG23	1:F:156:LYS:HB2	1.95	0.48
1:C:122:ILE:HG21	1:C:142:ALA:HB2	1.96	0.48
1:B:133:ASP:OD2	1:B:136:ASN:HB3	2.14	0.47
1:B:122:ILE:HG21	1:B:142:ALA:HB2	1.96	0.47
1:D:102:LEU:HB3	1:D:107:VAL:HG21	1.98	0.45
1:A:111:MSE:HG2	1:A:115:VAL:HG11	1.99	0.44
1:B:43:LYS:CE	1:E:22:ASP:HB3	2.48	0.43
1:C:102:LEU:HD12	1:C:117:ILE:HD13	2.00	0.43
1:D:223:VAL:HG22	1:F:209:ASN:ND2	2.33	0.43
1:F:98:SER:HB2	1:F:142:ALA:HB1	2.01	0.43
1:B:102:LEU:HD21	1:B:160:ILE:HD11	1.99	0.43
1:C:152:ARG:NH1	2:C:301:CL:CL	2.89	0.43
1:C:82:LYS:HE3	1:E:69:ASP:OD2	2.19	0.43
1:D:102:LEU:HD12	1:D:117:ILE:HD13	2.01	0.43
1:D:142:ALA:HB3	1:D:160:ILE:HB	2.00	0.43
1:C:53:GLN:OE1	1:E:113:THR:HG21	2.19	0.42
1:C:166:ILE:HG22	1:C:167:LYS:HG2	2.01	0.42
1:A:145:LYS:HD3	1:A:155:ARG:HG3	2.02	0.42
1:F:158:MSE:HG2	1:F:189:THR:HG23	2.01	0.41
1:B:102:LEU:HB3	1:B:107:VAL:HG21	2.02	0.40
1:C:53:GLN:HE22	1:E:113:THR:CB	2.34	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	213/215 (99%)	205 (96%)	8 (4%)	0	100	100
1	B	213/215 (99%)	201 (94%)	12 (6%)	0	100	100
1	C	213/215 (99%)	204 (96%)	9 (4%)	0	100	100
1	D	212/215 (99%)	203 (96%)	9 (4%)	0	100	100
1	E	211/215 (98%)	205 (97%)	5 (2%)	1 (0%)	24	48
1	F	213/215 (99%)	203 (95%)	10 (5%)	0	100	100
All	All	1275/1290 (99%)	1221 (96%)	53 (4%)	1 (0%)	48	73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	23	CYS

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	190/189 (100%)	188 (99%)	2 (1%)	65	85
1	B	189/189 (100%)	187 (99%)	2 (1%)	65	85
1	C	188/189 (100%)	185 (98%)	3 (2%)	55	80
1	D	191/189 (101%)	190 (100%)	1 (0%)	81	92
1	E	188/189 (100%)	185 (98%)	3 (2%)	55	80
1	F	188/189 (100%)	183 (97%)	5 (3%)	39	69

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	1134/1134 (100%)	1118 (99%)	16 (1%)	61	82

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	LEU
1	A	113	THR
1	B	111	MSE
1	B	113	THR
1	C	60	LEU
1	C	80	VAL
1	C	113	THR
1	D	60	LEU
1	E	60	LEU
1	E	113	THR
1	E	145	LYS
1	F	60	LEU
1	F	113	THR
1	F	128[A]	TYR
1	F	128[B]	TYR
1	F	137	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	ASN
1	A	126	ASN
1	A	225	GLN
1	B	29	GLN
1	B	64	GLN
1	B	136	ASN
1	B	225	GLN
1	C	91	ASN
1	C	151	ASN
1	C	225	GLN
1	D	91	ASN
1	D	135	ASN
1	D	151	ASN
1	D	225	GLN
1	E	45	ASN
1	E	151	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	E	225	GLN
1	F	29	GLN
1	F	64	GLN
1	F	120	ASN
1	F	209	ASN
1	F	225	GLN

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](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

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 [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	211/215 (98%)	-0.18	3 (1%) 73 72	43, 61, 99, 115	1 (0%)
1	B	211/215 (98%)	0.03	7 (3%) 49 45	45, 68, 108, 123	1 (0%)
1	C	211/215 (98%)	-0.23	2 (0%) 81 80	44, 60, 95, 113	1 (0%)
1	D	211/215 (98%)	-0.15	3 (1%) 73 72	39, 60, 93, 111	0
1	E	210/215 (97%)	-0.18	1 (0%) 87 86	44, 60, 84, 104	0
1	F	211/215 (98%)	0.00	2 (0%) 81 80	37, 71, 97, 121	1 (0%)
All	All	1265/1290 (98%)	-0.12	18 (1%) 73 72	37, 63, 97, 123	4 (0%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	88	TYR	4.0
1	D	103	SER	3.6
1	C	110	GLU	3.2
1	B	142	ALA	3.1
1	C	108	VAL	3.0
1	F	110	GLU	2.7
1	B	103	SER	2.5
1	B	108	VAL	2.5
1	B	101	THR	2.5
1	B	104	TYR	2.5
1	E	143	SER	2.4
1	B	107	VAL	2.4
1	B	106	GLY	2.4
1	F	128[A]	TYR	2.3
1	A	108	VAL	2.2
1	D	106	GLY	2.1
1	A	109	SER	2.1
1	D	104	TYR	2.1

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
2	CL	C	301	1/1	0.77	0.19	88,88,88,88	0
2	CL	C	302	1/1	0.81	0.14	95,95,95,95	0
2	CL	D	301	1/1	0.89	0.13	73,73,73,73	0
2	CL	F	301	1/1	0.95	0.06	70,70,70,70	0

6.5 Other polymers [i](#)

There are no such residues in this entry.