



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:48 AM UTC

PDB ID : 6Q86 / pdb_00006q86
Title : Structure of Fucosylated D-antimicrobial peptide SB4 in complex with the
Fucose-binding lectin PA-IIL at 2.008 Angstrom resolution
Authors : Baeriswyl, S.; Stocker, A.; Reymond, J.L.
Deposited on : 2018-12-14
Resolution : 2.01 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

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 2009 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 Fucose-binding lectin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	114	Total	C	N	O	0	0	0
			818	504	140	174			
1	B	114	Total	C	N	O	0	0	0
			818	504	140	174			

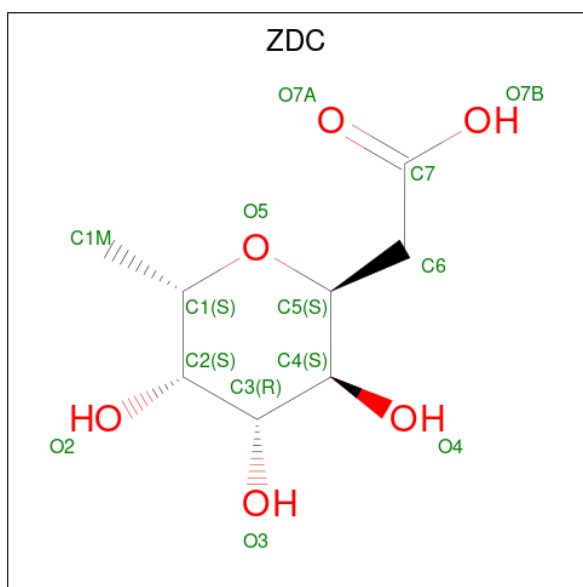
- Molecule 2 is a protein (with D amino acids) called SB4.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	13	Total	C	N	O	0	0	0
			98	68	16	14			
2	D	13	Total	C	N	O	0	0	0
			108	75	19	14			

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

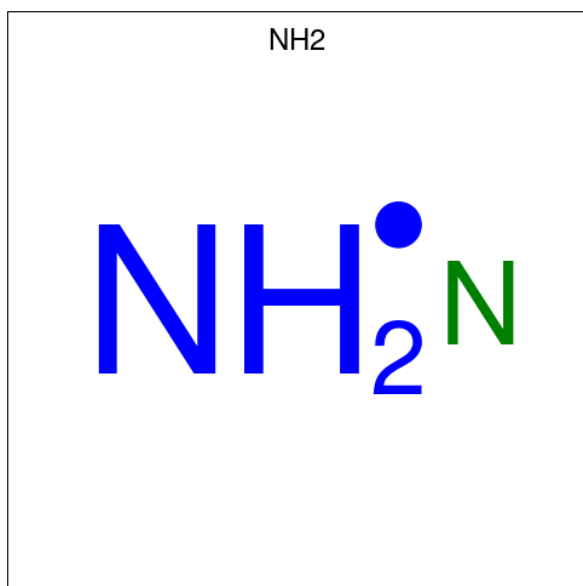
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Ca	0	0
			2	2		
3	B	2	Total	Ca	0	0
			2	2		

- Molecule 4 is 3,7-anhydro-2,8-dideoxy-L-glycero-D-gluco-octonic acid (CCD ID: ZDC) (formula: C₈H₁₄O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			13	8	5		
4	D	1	Total	C	O	0	0
			13	8	5		

- Molecule 5 is AMINO GROUP (CCD ID: NH2) (formula: H₂N).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	1	Total	N	0	0
			1	1		
5	D	1	Total	N	0	0
			1	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	71	Total 71	O 71	0	0
6	C	7	Total 7	O 7	0	0
6	B	48	Total 48	O 48	0	0
6	D	9	Total 9	O 9	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: Fucose-binding lectin

Chain A:  88% 12%



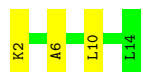
- Molecule 1: Fucose-binding lectin

Chain B:  82% 18%



- Molecule 2: SB4

Chain C:  77% 23%



- Molecule 2: SB4

Chain D:  92% 8%



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	55.60Å 55.60Å 150.62Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.15 – 2.01 48.15 – 2.01	Depositor EDS
% Data completeness (in resolution range)	99.3 (48.15-2.01) 99.7 (48.15-2.01)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.03	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.61 (at 2.01Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.157 , 0.190 0.160 , 0.183	Depositor DCC
R_{free} test set	937 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	39.4	Xtriage
Anisotropy	0.327	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 35.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.254 for -h,-k,l	Xtriage
Reported twinning fraction	0.260 for -h,-k,l	Depositor
Outliers	0 of 18701 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	2009	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.50% 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: DLY, DTY, DAL, ZDC, DLE, NH2, CA

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	0/828	0.46	0/1135
1	B	0.32	0/828	0.47	0/1135
All	All	0.32	0/1656	0.46	0/2270

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	818	0	787	11	0
1	B	818	0	787	17	0
2	C	98	0	101	2	0
2	D	108	0	128	1	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	C	13	0	9	1	0
4	D	13	0	10	1	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
6	A	71	0	0	4	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	48	0	0	7	0
6	C	7	0	0	0	0
6	D	9	0	0	1	0
All	All	2009	0	1822	29	1

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 8.

All (29) 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:26:GLN:NE2	6:B:301:HOH:O	1.91	0.98
1:A:33:ASN:OD1	6:A:301:HOH:O	1.91	0.88
1:B:58:GLY:N	6:B:303:HOH:O	2.21	0.72
1:A:90:ALA:HB3	1:A:109:ILE:HB	1.72	0.71
1:B:72:ARG:O	6:B:302:HOH:O	2.10	0.69
1:A:2:THR:OG1	6:A:302:HOH:O	2.10	0.69
1:A:12:THR:H	1:A:57:SER:HB3	1.59	0.66
1:B:25:THR:HG22	1:B:43:GLN:HB2	1.81	0.63
2:D:2:DLY:NZ	6:D:201:HOH:O	2.25	0.60
1:B:12:THR:O	1:B:57:SER:OG	2.26	0.53
1:B:34:ASN:OD1	1:B:62:LYS:HE2	2.12	0.49
1:B:58:GLY:C	6:B:303:HOH:O	2.54	0.49
1:A:33:ASN:HA	6:A:301:HOH:O	2.13	0.48
2:C:6:DAL:O	2:C:10:DLE:HD12	2.15	0.47
1:B:1:ALA:N	6:B:308:HOH:O	2.47	0.46
1:B:99:ASP:OD2	4:D:1:ZDC:O3	2.34	0.45
1:B:90:ALA:HB3	1:B:109:ILE:HB	1.99	0.45
1:A:57:SER:OG	1:A:61:GLY:HA2	2.17	0.45
1:A:96:ASP:OD1	4:C:1:ZDC:O4	2.35	0.45
1:A:102:TYR:CZ	1:B:1:ALA:HB2	2.52	0.44
1:B:92:VAL:HB	1:B:107:VAL:HB	1.98	0.44
1:B:22:SER:OG	1:B:23:SER:N	2.52	0.43
6:A:337:HOH:O	2:C:2:DLY:HD2	2.19	0.43
1:B:58:GLY:CA	6:B:303:HOH:O	2.64	0.43
1:B:26:GLN:O	1:B:41:SER:HA	2.19	0.43
1:A:40:PHE:CG	1:A:51:GLY:HA3	2.54	0.43
1:B:58:GLY:O	6:B:303:HOH:O	2.21	0.42
1:A:1:ALA:HB2	1:B:102:TYR:CZ	2.56	0.41
1:A:26:GLN:O	1:A:41:SER:HA	2.20	0.41

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:368:HOH:O	6:A:371:HOH:O[4_555]	2.10	0.10

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	112/114 (98%)	108 (96%)	4 (4%)	0	100	100
1	B	112/114 (98%)	108 (96%)	3 (3%)	1 (1%)	14	9
All	All	224/228 (98%)	216 (96%)	7 (3%)	1 (0%)	30	27

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	84	THR

5.3.2 Protein sidechains [i](#)

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	91/93 (98%)	91 (100%)	0	100	100
1	B	91/93 (98%)	91 (100%)	0	100	100
All	All	182/186 (98%)	182 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

Mol	Chain	Res	Type
1	A	26	GLN
1	A	29	ASN
1	A	66	GLN
1	B	26	GLN
1	B	46	ASN
1	B	66	GLN
1	B	80	GLN
1	B	88	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

26 non-standard protein/DNA/RNA residues are modelled in this entry.

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.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 4 are monoatomic and 2 are modelled with single atom - leaving 2 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
4	ZDC	C	1	3	13,13,14	0.78	0	18,18,20	2.45	5 (27%)
4	ZDC	D	1	3	13,13,14	0.78	0	18,18,20	2.45	5 (27%)

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
4	ZDC	C	1	3	-	3/3/23/24	0/1/1/1
4	ZDC	D	1	3	-	2/3/23/24	0/1/1/1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1	ZDC	C6-C5-C4	7.20	123.04	114.16
4	C	1	ZDC	C6-C5-C4	6.44	122.10	114.16
4	C	1	ZDC	O5-C5-C6	4.33	125.32	109.98
4	D	1	ZDC	O5-C5-C6	3.66	122.94	109.98
4	C	1	ZDC	O7A-C7-C6	-3.64	114.79	125.38
4	D	1	ZDC	O7A-C7-C6	-3.52	115.14	125.38
4	C	1	ZDC	C3-C4-C5	-3.23	104.37	110.23
4	D	1	ZDC	C3-C4-C5	-2.87	105.03	110.23
4	C	1	ZDC	C5-C6-C7	2.46	117.01	112.72
4	D	1	ZDC	O2-C2-C1	-2.11	105.08	109.74

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	1	ZDC	O5-C5-C6-C7
4	D	1	ZDC	O5-C5-C6-C7
4	C	1	ZDC	C5-C6-C7-O7A
4	C	1	ZDC	C4-C5-C6-C7

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Mol	Chain	Res	Type	Atoms
4	D	1	ZDC	C5-C6-C7-O7A

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	1	ZDC	1	0
4	D	1	ZDC	1	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 [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	114/114 (100%)	-1.12	0 100 100	26, 42, 59, 93	0
1	B	114/114 (100%)	-0.66	0 100 100	29, 52, 70, 84	0
2	C	0/13	-	-	-	-
2	D	0/13	-	-	-	-
All	All	228/254 (89%)	-0.89	0 100 100	26, 46, 70, 93	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
2	DLY	D	12	9/10	0.96	0.10	48,58,76,77	0
2	DLY	D	2	9/10	0.97	0.05	42,48,56,59	0
2	DLY	C	4	9/10	0.98	0.07	38,46,64,65	0
2	DLY	D	4	9/10	0.98	0.06	38,44,60,64	0
2	DLY	C	5	6/10	0.98	0.06	39,46,51,54	0
2	DLY	D	5	9/10	0.98	0.07	40,44,67,68	0
2	DAL	C	6	5/6	0.98	0.04	35,35,41,42	0
2	DLY	C	8	5/10	0.98	0.05	35,35,37,43	0
2	DLY	D	8	9/10	0.98	0.07	34,39,60,61	0
2	DLY	D	9	9/10	0.98	0.04	37,45,50,60	0
2	DLY	C	12	6/10	0.98	0.05	41,45,49,54	0
2	DLY	C	2	9/10	0.98	0.06	45,50,70,71	0
2	DLE	D	13	8/9	0.98	0.06	39,48,58,59	0
2	DLE	D	14	8/9	0.98	0.05	42,44,55,55	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	DLY	C	9	9/10	0.99	0.04	37,38,59,66	0
2	DAL	D	6	5/6	0.99	0.03	40,40,45,48	0
2	DLE	D	10	8/9	0.99	0.05	40,44,49,53	0
2	DAL	C	11	5/6	0.99	0.05	40,42,48,52	0
2	DAL	D	11	5/6	0.99	0.09	40,40,47,48	0
2	DLE	C	7	8/9	0.99	0.04	35,39,42,47	0
2	DLE	D	7	8/9	0.99	0.03	31,33,39,39	0
2	DLE	C	13	8/9	0.99	0.04	40,42,43,47	0
2	DTY	D	3	12/13	0.99	0.04	41,46,54,58	0
2	DLE	C	14	8/9	0.99	0.04	39,42,42,46	0
2	DTY	C	3	12/13	0.99	0.05	45,70,74,76	0
2	DLE	C	10	8/9	1.00	0.03	37,43,47,49	0

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	NH2	D	101	1/1	0.96	0.15	66,66,66,66	0
4	ZDC	D	1	13/14	0.99	0.05	44,46,47,53	0
5	NH2	C	101	1/1	0.99	0.09	39,39,39,39	0
4	ZDC	C	1	13/14	0.99	0.03	33,39,51,53	0
3	CA	A	202	1/1	1.00	0.02	40,40,40,40	0
3	CA	A	203	1/1	1.00	0.02	42,42,42,42	0
3	CA	B	202	1/1	1.00	0.01	43,43,43,43	0
3	CA	B	203	1/1	1.00	0.04	47,47,47,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.