



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:40 AM GMT

PDB ID : 3BQK  
Title : Structure of a chondroitin sulphate binding DBL3X from a var2csa encoded PfEMP1 protein in complex with sulphate  
Authors : Higgins, M.K.  
Deposited on : 2007-12-20  
Resolution : 1.80 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

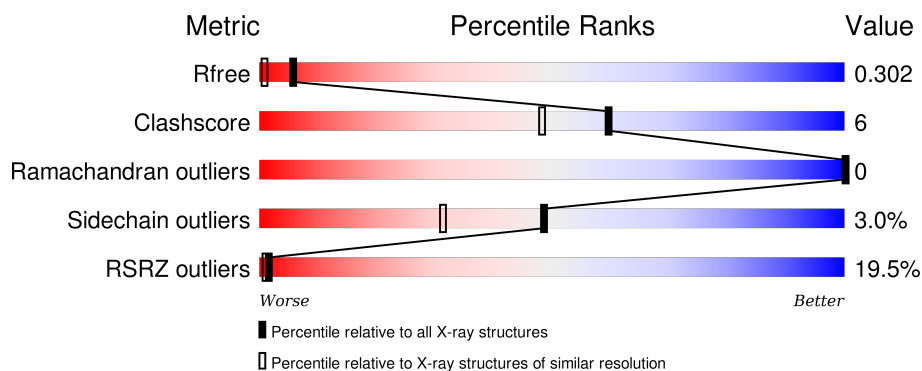
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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))
$R_{free}$	91344	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	360	<div> <div>18%</div> <div>82%</div> <div>11%</div> <div>6%</div> </div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2975 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 Erythrocyte membrane protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	338	2732	1713	471	530	18	0	0	0

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

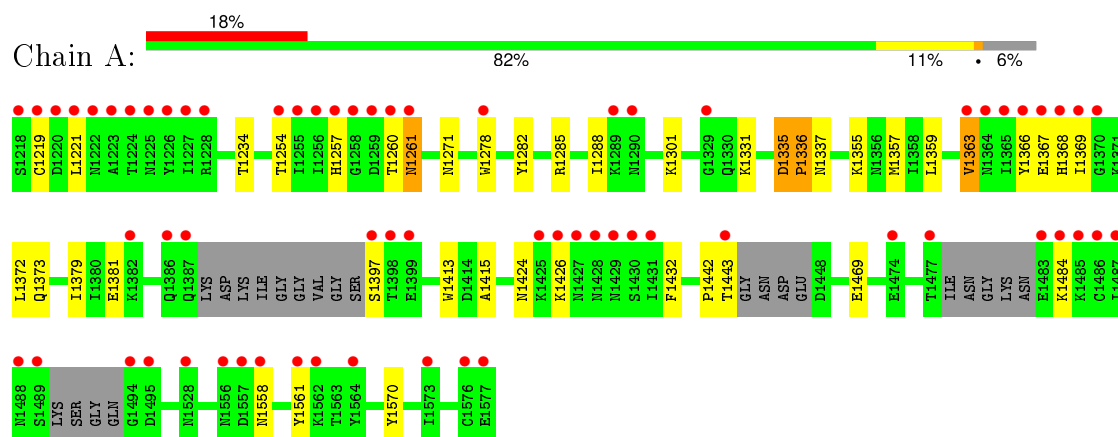
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	232	Total	O	0	0
			232	232		

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Erythrocyte membrane protein 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	41.26 Å 86.56 Å 92.33 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.18 – 1.80 46.16 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (46.18-1.80) 99.8 (46.16-1.80)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.39 (at 1.79 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.231 , 0.286 0.250 , 0.302	Depositor DCC
$R_{free}$ test set	1575 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.2	Xtriage
Anisotropy	0.123	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 52.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 31359 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	2975	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.27% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SO4

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	0.50	1/2785 (0.0%)	0.55	1/3733 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1413	TRP	CG-CD1	-5.38	1.29	1.36

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1336	PRO	N-CA-C	-7.14	93.54	112.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1335	ASP	Peptide

## 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	2732	0	2649	31	0
2	A	5	0	0	0	0
3	A	6	0	8	0	0
4	A	232	0	0	1	0
All	All	2975	0	2657	31	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 6.

All (31) 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:A:1357:MET:HA	1:A:1373:GLN:HE21	1.28	0.97
1:A:1442:PRO:HA	1:A:1443:THR:HB	1.51	0.92
1:A:1424:ASN:HD21	1:A:1432:PHE:H	1.18	0.85
1:A:1357:MET:HA	1:A:1373:GLN:NE2	1.91	0.85
1:A:1442:PRO:CA	1:A:1443:THR:HB	2.15	0.77
1:A:1234:THR:H	1:A:1271:ASN:HD21	1.37	0.70
1:A:1278:TRP:CH2	1:A:1368:HIS:HB2	2.30	0.67
1:A:1369:ILE:HD13	1:A:1372:LEU:HD12	1.76	0.65
1:A:1357:MET:CA	1:A:1373:GLN:HE21	2.09	0.63
1:A:1335:ASP:OD1	1:A:1335:ASP:C	2.36	0.63
1:A:1336:PRO:CD	1:A:1337:ASN:H	2.11	0.61
1:A:1278:TRP:CZ2	1:A:1368:HIS:HB2	2.38	0.59
1:A:1234:THR:H	1:A:1271:ASN:ND2	2.02	0.57
1:A:1558:ASN:HB2	1:A:1561:TYR:CD2	2.42	0.55
1:A:1355:LYS:O	1:A:1359:LEU:HD13	2.06	0.55
1:A:1381:GLU:HG2	1:A:1397:SER:N	2.23	0.54
1:A:1257:HIS:CD2	1:A:1363:VAL:HG23	2.43	0.54
1:A:1335:ASP:CB	1:A:1336:PRO:HD3	2.39	0.53
1:A:1288:ILE:O	1:A:1379:ILE:HD11	2.09	0.52
1:A:1366:TYR:HB2	1:A:1369:ILE:HB	1.91	0.52
1:A:1336:PRO:CD	1:A:1337:ASN:N	2.68	0.50
1:A:1442:PRO:HA	1:A:1443:THR:CB	2.31	0.49
1:A:1331:LYS:HE3	4:A:1806:HOH:O	2.12	0.48

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1335:ASP:HB2	1:A:1336:PRO:HD3	1.94	0.47
1:A:1336:PRO:HD2	1:A:1337:ASN:N	2.31	0.46
1:A:1301:LYS:HG3	1:A:1415:ALA:HB2	1.98	0.45
1:A:1261:ASN:N	1:A:1261:ASN:OD1	2.49	0.45
1:A:1469:GLU:HB2	1:A:1570:TYR:CZ	2.54	0.43
1:A:1336:PRO:HD2	1:A:1337:ASN:H	1.81	0.42
1:A:1282:TYR:HB2	1:A:1285:ARG:HH12	1.85	0.41
1:A:1335:ASP:CG	1:A:1336:PRO:HD3	2.40	0.41

There are no symmetry-related clashes.

## 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	328/360 (91%)	317 (97%)	11 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	300/316 (95%)	291 (97%)	9 (3%)	48	31

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1219	CYS
1	A	1221	LEU
1	A	1254	THR
1	A	1260	THR
1	A	1261	ASN
1	A	1363	VAL
1	A	1367	GLU
1	A	1426	LYS
1	A	1484	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1271	ASN
1	A	1305	HIS
1	A	1356	ASN
1	A	1373	GLN
1	A	1424	ASN

### 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 carbohydrates 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	$\# Z  > 2$
2	SO4	A	1	-	4,4,4	0.19	0	6,6,6	0.23	0
3	GOL	A	2	-	5,5,5	0.39	0	5,5,5	0.20	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
2	SO4	A	1	-	-	0/0/0/0	0/0/0/0
3	GOL	A	2	-	-	0/4/4/4	0/0/0/0

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	338/360 (93%)	1.19	66 (19%) ⓘ ⓘ	2, 18, 35, 52	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1494	GLY	9.9
1	A	1219	CYS	9.1
1	A	1222	ASN	8.7
1	A	1221	LEU	8.3
1	A	1226	TYR	8.1
1	A	1224	THR	8.0
1	A	1254	THR	7.7
1	A	1484	LYS	7.4
1	A	1487	ILE	6.7
1	A	1397	SER	6.3
1	A	1218	SER	5.8
1	A	1561	TYR	5.6
1	A	1364	ASN	5.6
1	A	1223	ALA	5.5
1	A	1577	GLU	5.5
1	A	1365	ILE	5.5
1	A	1443	THR	5.3
1	A	1228	ARG	5.0
1	A	1429	ASN	5.0
1	A	1258	GLY	4.9
1	A	1363	VAL	4.8
1	A	1398	THR	4.8
1	A	1220	ASP	4.6
1	A	1256	ILE	4.6
1	A	1486	CYS	4.5
1	A	1488	ASN	4.5
1	A	1387	GLN	4.4

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	1259	ASP	4.4
1	A	1485	LYS	4.3
1	A	1428	ASN	4.2
1	A	1225	ASN	4.1
1	A	1368	HIS	4.1
1	A	1257	HIS	4.0
1	A	1255	ILE	3.9
1	A	1557	ASP	3.7
1	A	1477	THR	3.7
1	A	1227	ILE	3.6
1	A	1367	GLU	3.5
1	A	1556	ASN	3.5
1	A	1430	SER	3.4
1	A	1369	ILE	3.3
1	A	1564	TYR	3.3
1	A	1382	LYS	3.2
1	A	1495	ASP	3.1
1	A	1260	THR	3.0
1	A	1489	SER	3.0
1	A	1289	LYS	2.9
1	A	1290	ASN	2.8
1	A	1386	GLN	2.7
1	A	1576	CYS	2.7
1	A	1528	ASN	2.7
1	A	1431	ILE	2.7
1	A	1427	ASN	2.7
1	A	1425	LYS	2.7
1	A	1562	LYS	2.7
1	A	1261	ASN	2.6
1	A	1558	ASN	2.6
1	A	1399	GLU	2.6
1	A	1366	TYR	2.5
1	A	1329	GLY	2.4
1	A	1278	TRP	2.4
1	A	1426	LYS	2.4
1	A	1483	GLU	2.2
1	A	1370	GLY	2.1
1	A	1573	ILE	2.1
1	A	1474	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	GOL	A	2	6/6	0.78	0.29	1.74	36,38,39,40	0
2	SO4	A	1	5/5	0.93	0.12	-1.07	42,42,43,43	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.