

wwPDB X-ray Structure Validation Summary Report (i)

Oct 15, 2023 – 07:02 PM EDT

PDB ID	:	8DY2
Title	:	Crystal Structure of spFv GLK1
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Deposited on	:	2022-08-03
Resolution	:	1.65 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) 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 (1)) were used in the production of this report:

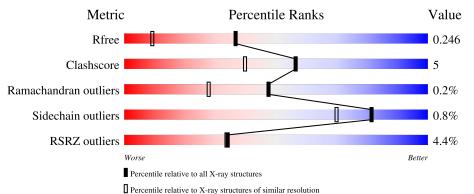
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 1.65 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	1827 (1.66-1.66)
Clashscore	141614	1931 (1.66-1.66)
Ramachandran outliers	138981	1891 (1.66-1.66)
Sidechain outliers	138945	1891 (1.66-1.66)
RSRZ outliers	127900	1791 (1.66-1.66)

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 <=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	А	252	2% 88%	8%	·
1	В	252	3%	10%	•
1	С	252	82%	15%	·
1	D	252	87%	10%	·

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard



residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	А	301	-	-	Х	-



2 Entry composition (i)

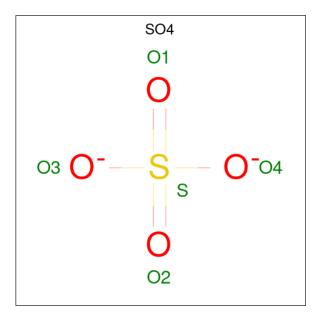
There are 3 unique types of molecules in this entry. The entry contains 7592 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	Δ	242	Total	С	Ν	0	\mathbf{S}	0	0	0
	А	242	1768	1104	294	359	11	0	0	0
1	В	245	Total	С	Ν	0	S	0	0	0
	D	240	1781	1110	297	363	11	0	0	0
1	С	244	Total	С	Ν	0	S	0	0	0
	U	244	1782	1111	297	363	11	0	0	0
1	Л	243	Total	С	Ν	0	S	0	0	0
		240	1783	1112	299	361	11	0	0	0

• Molecule 1 is a protein called spFv GLK1 LH.

• Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

• Molecule 3 is water.

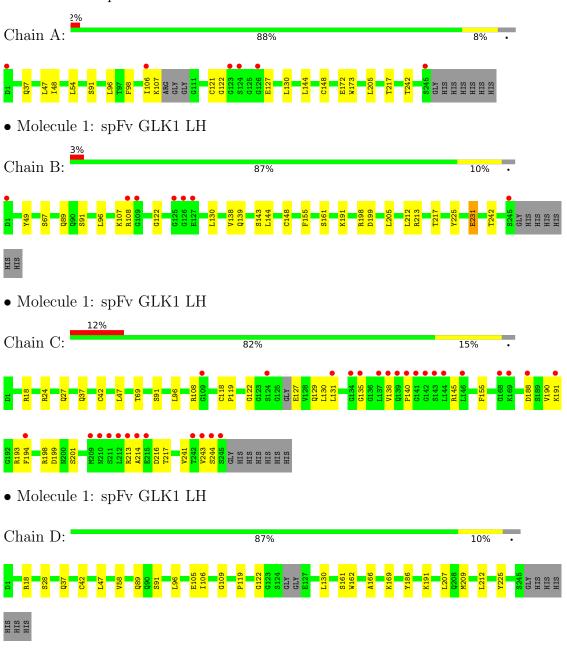


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	137	Total O 137 137	0	0
3	В	126	Total O 126 126	0	0
3	С	87	Total O 87 87	0	0
3	D	123	Total O 123 123	0	0



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.



 \bullet Molecule 1: spFv GLK1 LH



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	45.63Å 78.94Å 257.30Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.86 - 1.65	Depositor
	128.65 - 1.65	EDS
% Data completeness	47.8(35.86-1.65)	Depositor
(in resolution range)	44.7 (128.65 - 1.65)	EDS
R_{merge}	0.09	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.28 (at 1.65 \text{\AA})$	Xtriage
Refinement program	PHENIX 1.18_3855	Depositor
B B.	0.215 , 0.246	Depositor
R, R_{free}	0.215 , 0.246	DCC
R_{free} test set	2675 reflections $(4.93%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	20.5	Xtriage
Anisotropy	0.145	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35 , 36.0	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	0.94	EDS
Total number of atoms	7592	wwPDB-VP
Average B, all atoms $(Å^2)$	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.10% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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: 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		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.27	0/1806	0.50	0/2448	
1	В	0.27	0/1820	0.49	0/2469	
1	С	0.27	0/1820	0.49	0/2466	
1	D	0.28	0/1821	0.49	0/2467	
All	All	0.27	0/7267	0.49	0/9850	

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	А	1768	0	1675	14	0
1	В	1781	0	1679	16	0
1	С	1782	0	1689	23	0
1	D	1783	0	1694	16	0
2	А	5	0	0	2	0
3	А	137	0	0	3	1
3	В	126	0	0	4	1
3	С	87	0	0	5	0
3	D	123	0	0	3	0
All	All	7592	0	6737	69	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 5.

The worst 5 of 69 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:139:GLN:NE2	3:B:302:HOH:O	2.04	0.85
1:C:18:ARG:NH1	3:C:302:HOH:O	2.12	0.83
1:C:108:ARG:O	3:C:301:HOH:O	1.99	0.81
1:B:67:SER:O	3:B:301:HOH:O	1.99	0.79
1:C:217:THR:HG22	1:C:243:VAL:H	1.51	0.74

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)
3:A:456:HOH:O	3:B:401:HOH:O[4_445]	2.13	0.07

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	Perce	ntiles
1	А	238/252~(94%)	230~(97%)	7 (3%)	1 (0%)	34	16
1	В	243/252~(96%)	236~(97%)	7 (3%)	0	100	100
1	С	240/252~(95%)	228 (95%)	11 (5%)	1 (0%)	34	16
1	D	239/252~(95%)	233~(98%)	6 (2%)	0	100	100
All	All	960/1008~(95%)	927~(97%)	31 (3%)	2(0%)	47	28

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type					
1	А	127	GLU					
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Mol	Chain	Res	Type
1	С	140	PRO

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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	А	190/201~(94%)	190 (100%)	0	100 100
1	В	190/201~(94%)	187~(98%)	3~(2%)	62 41
1	С	192/201~(96%)	190~(99%)	2(1%)	76 62
1	D	192/201~(96%)	191 (100%)	1 (0%)	88 81
All	All	764/804~(95%)	758~(99%)	6 (1%)	81 70

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	С	145	ARG
1	С	244	SER
1	D	18	ARG
1	В	213	ARG
1	В	89	GLN

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)

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	B	ond leng	gths	В	ond ang	gles
IVI01	rybe		nam nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2
2	SO4	А	301	-	$4,\!4,\!4$	0.18	0	$6,\!6,\!6$	0.18	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.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	А	301	SO4	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 (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, 95^{th} 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	$\mathbf{OWAB}(\mathbf{A}^2)$	Q < 0.9
1	А	242/252~(96%)	0.01	6 (2%) 57 58	15, 23, 34, 49	0
1	В	245/252~(97%)	0.02	7 (2%) 51 52	15, 24, 38, 52	0
1	С	244/252~(96%)	0.57	30 (12%) 4 3	17, 30, 57, 67	0
1	D	243/252~(96%)	-0.08	0 100 100	17, 25, 35, 40	0
All	All	974/1008~(96%)	0.13	43 (4%) 34 34	15, 25, 45, 67	0

The worst 5 of 43 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	С	140	PRO	9.1
1	С	141	GLY	7.7
1	А	123	GLY	5.7
1	С	124	SER	5.6
1	С	213	ARG	5.6

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, 95^{th} 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	SO4	А	301	5/5	0.98	0.36	27,33,40,41	0

6.5 Other polymers (i)

There are no such residues in this entry.

