W O R L D W I D E
wwPDB
P R O T E I N D A T A B A N K

wwPDB X-ray Structure Validation Summary Report (i)

Dec 16, 2024 – 12:17 PM EST

PDB ID	:	8VSR
Title	:	Co-crystal structure of Prx with Esub1
Authors	:	Muna, T.H.; Prehna, G.
Deposited on	:	2024-01-24
Resolution	:	2.69 Å(reported)

The following versions of software and data (see references (i)) were used in the production of this report:

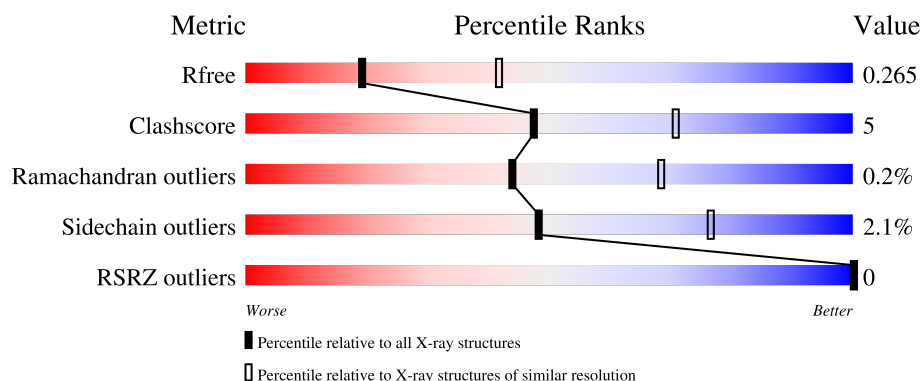|		
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.40

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 2.69 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	Percentile Ranks	Value
Rfree		0.265
Clashscore		5
Ramachandran outliers		0.2%
Sidechain outliers		2.1%
RSRZ outliers		0

Worse   *Better*

▮ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	3333 (2.70-2.70)
Clashscore	180529	3684 (2.70-2.70)
Ramachandran outliers	177936	3633 (2.70-2.70)
Sidechain outliers	177891	3633 (2.70-2.70)
RSRZ outliers	164620	3333 (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 <=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	68	71% 18% 12%
1	C	68	72% 16% 12%
2	B	198	80% 13% • 7%
2	D	198	78% 15% • 7%

2 Entry composition ⓘ

There are 5 unique types of molecules in this entry. The entry contains 8168 atoms, of which 4080 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 Prx.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	60	Total	C	H	N	O	S	0	0	0
			992	316	500	78	96	2			
1	C	60	Total	C	H	N	O	S	0	0	0
			992	316	500	78	96	2			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	61	LEU	-	expression tag	UNP A0A0H2UWN8
A	62	GLU	-	expression tag	UNP A0A0H2UWN8
A	63	HIS	-	expression tag	UNP A0A0H2UWN8
A	64	HIS	-	expression tag	UNP A0A0H2UWN8
A	65	HIS	-	expression tag	UNP A0A0H2UWN8
A	66	HIS	-	expression tag	UNP A0A0H2UWN8
A	67	HIS	-	expression tag	UNP A0A0H2UWN8
A	68	HIS	-	expression tag	UNP A0A0H2UWN8
C	61	LEU	-	expression tag	UNP A0A0H2UWN8
C	62	GLU	-	expression tag	UNP A0A0H2UWN8
C	63	HIS	-	expression tag	UNP A0A0H2UWN8
C	64	HIS	-	expression tag	UNP A0A0H2UWN8
C	65	HIS	-	expression tag	UNP A0A0H2UWN8
C	66	HIS	-	expression tag	UNP A0A0H2UWN8
C	67	HIS	-	expression tag	UNP A0A0H2UWN8
C	68	HIS	-	expression tag	UNP A0A0H2UWN8

- Molecule 2 is a protein called Esub1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	185	Total	C	H	N	O	S	0	0	0
			3027	951	1541	262	269	4			
2	D	184	Total	C	H	N	O	S	0	0	0
			3007	945	1529	261	268	4			

WORLDWIDE
wwPDB
PROTEIN DATA BANK

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	9	MET	-	initiating methionine	UNP A0A0H2UUU0
B	199	LEU	-	expression tag	UNP A0A0H2UUU0
B	200	GLU	-	expression tag	UNP A0A0H2UUU0
B	201	HIS	-	expression tag	UNP A0A0H2UUU0
B	202	HIS	-	expression tag	UNP A0A0H2UUU0
B	203	HIS	-	expression tag	UNP A0A0H2UUU0
B	204	HIS	-	expression tag	UNP A0A0H2UUU0
B	205	HIS	-	expression tag	UNP A0A0H2UUU0
B	206	HIS	-	expression tag	UNP A0A0H2UUU0
D	9	MET	-	initiating methionine	UNP A0A0H2UUU0
D	199	LEU	-	expression tag	UNP A0A0H2UUU0
D	200	GLU	-	expression tag	UNP A0A0H2UUU0
D	201	HIS	-	expression tag	UNP A0A0H2UUU0
D	202	HIS	-	expression tag	UNP A0A0H2UUU0
D	203	HIS	-	expression tag	UNP A0A0H2UUU0
D	204	HIS	-	expression tag	UNP A0A0H2UUU0
D	205	HIS	-	expression tag	UNP A0A0H2UUU0
D	206	HIS	-	expression tag	UNP A0A0H2UUU0

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).
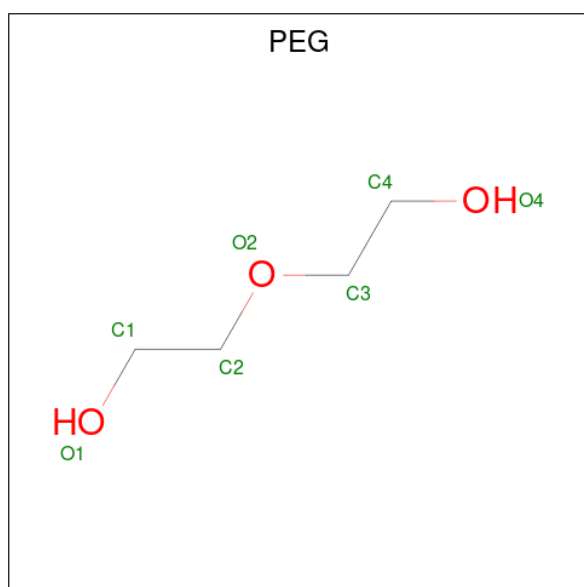
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
3	B	1	17	4	10	3	0	0

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).

PO4

O3

O⁻

O2 O⁻ — P = O O1

P

O⁻

O4

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total 5	O 4	P 1	0	0
4	C	1	Total 5	O 4	P 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	21	Total 21	O 21	0	0
5	B	43	Total 43	O 43	0	0
5	C	16	Total 16	O 16	0	0
5	D	43	Total 43	O 43	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.
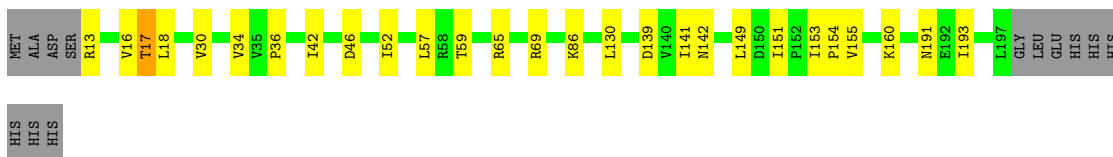
- Molecule 1: Prx

Chain A:

- Molecule 1: Prx

Chain C:

- Molecule 2: Esub1

Chain B:

- Molecule 2: Esub1

Chain D:

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α, β, γ	67.49Å 67.49Å 232.21Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.70 – 2.69 38.70 – 2.69	Depositor EDS
% Data completeness (in resolution range)	99.1 (38.70-2.69) 99.5 (38.70-2.69)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.30 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.228 , 0.265 0.228 , 0.265	Depositor DCC
R_{free} test set	826 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	52.1	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.40 , 53.4	EDS
L-test for twinning[2]	$< |L| > = 0.45, < L^2 > = 0.27$	Xtriage
Estimated twinning fraction	0.479 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	8168	wwPDB-VP
Average B, all atoms (Å^2)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.32% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: PO4, PEG

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	0.26	0/498	0.47	0/670
1	C	0.27	0/498	0.47	0/670
2	B	0.25	0/1511	0.48	0/2040
2	D	0.25	0/1503	0.50	0/2029
All	All	0.26	0/4010	0.49	0/5409

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 ⓘ

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	492	500	500	7	0
1	C	492	500	500	5	0
2	B	1486	1541	1540	16	1
2	D	1478	1529	1529	20	1
3	B	7	10	10	0	0
4	B	5	0	0	0	0
4	C	5	0	0	0	0
5	A	21	0	0	1	0
5	B	43	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	16	0	0	0	0
5	D	43	0	0	1	0
All	All	4088	4080	4079	44	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 44 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:B:142:ASN:OD1	5:B:401:HOH:O	2.06	0.74
1:A:19:ASP:OD2	1:A:50:ARG:NH1	2.26	0.69
1:A:44:GLU:OE1	2:B:65:ARG:NH2	2.27	0.68
1:A:44:GLU:OE2	5:A:101:HOH:O	2.13	0.66
2:B:16:VAL:HG13	2:B:155:VAL:HG21	1.78	0.64

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)
2:B:160:LYS:NZ	2:D:133:ALA:O[1_455]	2.00	0.20

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	58/68 (85%)	58 (100%)	0	0	100	100
1	C	58/68 (85%)	58 (100%)	0	0	100	100
2	B	183/198 (92%)	177 (97%)	6 (3%)	0	100	100
2	D	182/198 (92%)	176 (97%)	5 (3%)	1 (0%)	25	49

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	481/532 (90%)	469 (98%)	11 (2%)	1 (0%)	44	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	36	PRO

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	55/63 (87%)	53 (96%)	2 (4%)	30	59
1	C	55/63 (87%)	53 (96%)	2 (4%)	30	59
2	B	165/176 (94%)	162 (98%)	3 (2%)	54	80
2	D	164/176 (93%)	162 (99%)	2 (1%)	67	86
All	All	439/478 (92%)	430 (98%)	9 (2%)	48	76

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	17	THR
2	D	39	VAL
2	B	17	THR
2	B	191	ASN
1	C	29	LYS

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

5.3.3   RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

3 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	$\#|Z| > 2$
4	PO4	C	101	-	4,4,4	0.94	0	6,6,6	0.47	0
3	PEG	B	301	-	6,6,6	0.13	0	5,5,5	0.11	0
4	PO4	B	302	-	4,4,4	0.93	0	6,6,6	0.45	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
3	PEG	B	301	-	-	1/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	301	PEG	O2-C3-C4-O4

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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(Å²)	Q<0.9
1	A	60/68 (88%)	-1.57	0	100	100	45, 69, 89, 97	0
1	C	60/68 (88%)	-1.52	0	100	100	46, 73, 101, 116	0
2	B	185/198 (93%)	-1.54	0	100	100	41, 62, 106, 142	0
2	D	184/198 (92%)	-1.57	0	100	100	38, 63, 98, 145	0
All	All	489/532 (91%)	-1.55	0	100	100	38, 64, 100, 145	0

There are no RSRZ outliers to report.

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3   Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4   Ligands ⓘ

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(Å²)	Q<0.9
4	PO4	B	302	5/5	0.98	0.06	61,64,67,85	0
3	PEG	B	301	7/7	0.99	0.06	44,54,60,60	0
4	PO4	C	101	5/5	1.00	0.03	49,59,64,74	0

6.5 Other polymers ⓘ

There are no such residues in this entry.