Full wwPDB X-ray Structure Validation Report (i)

Mar 31, 2025 – 12:19 PM EDT

PDB ID	:	8VP4 / pdb_00008vp4
Title	:	Crystal Structure of JF1cpCasp2 with Peptide Inhibitor AcVDVAD-CHO
Authors	:	Fuller, J.L.; Shi, K.; Walters, M.A.; Finzel, B.C.
Deposited on	:	2024-01-16
Resolution	:	1.51 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 (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:

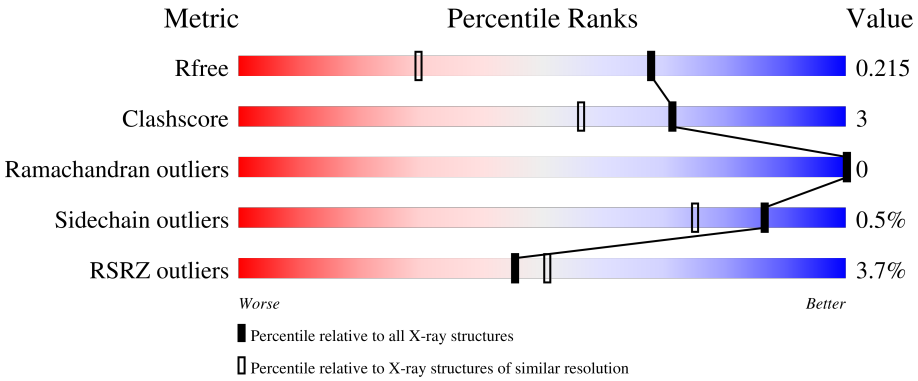
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.006 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.42

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.51 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}	164625	5293 (1.54-1.50)
Clashscore	180529	5759 (1.54-1.50)
Ramachandran outliers	177936	5653 (1.54-1.50)
Sidechain outliers	177891	5650 (1.54-1.50)
RSRZ outliers	164620	5293 (1.54-1.50)

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	282	3% 78% 6% 16%
1	B	282	4% 79% • 16%
2	C	6	100%
2	D	6	83% 17%

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	ASA	D	6	-	-	X	-

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4058 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 JF1cpCasp2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	S	0	3	0
			1847	1160	328	340	19			
1	B	237	Total	C	N	O	S	0	4	0
			1850	1162	328	340	20			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	327	MET	-	initiating methionine	UNP P42575
A	328	HIS	-	expression tag	UNP P42575
A	329	HIS	-	expression tag	UNP P42575
A	330	HIS	-	expression tag	UNP P42575
A	331	HIS	-	expression tag	UNP P42575
A	332	HIS	-	expression tag	UNP P42575
A	333	HIS	-	expression tag	UNP P42575
A	347	ALA	ASP	conflict	UNP P42575
A	1174	GLY	-	linker	UNP P42575
A	1175	GLY	-	linker	UNP P42575
B	327	MET	-	initiating methionine	UNP P42575
B	328	HIS	-	expression tag	UNP P42575
B	329	HIS	-	expression tag	UNP P42575
B	330	HIS	-	expression tag	UNP P42575
B	331	HIS	-	expression tag	UNP P42575
B	332	HIS	-	expression tag	UNP P42575
B	333	HIS	-	expression tag	UNP P42575
B	347	ALA	ASP	conflict	UNP P42575
B	1174	GLY	-	linker	UNP P42575
B	1175	GLY	-	linker	UNP P42575

- Molecule 2 is a protein called AcVDVAD-CHO.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
2	D	6	Total C N O 38 23 5 10	0	0	0
2	C	6	Total C N O 38 23 5 10	0	0	0

- Molecule 3 is water.

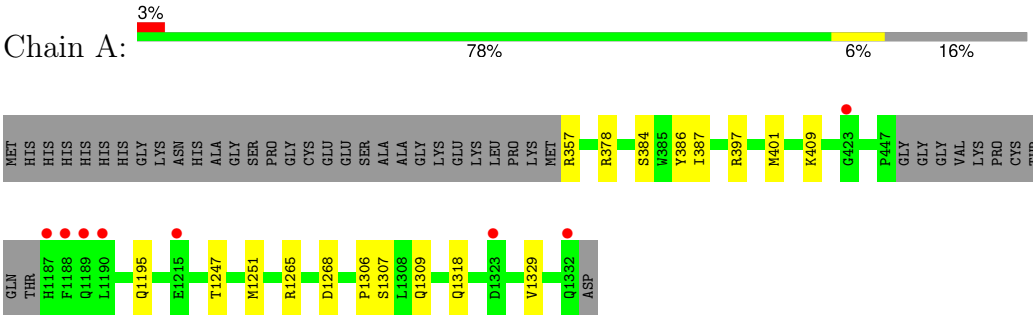
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	134	Total O 134 134	0	0
3	B	144	Total O 144 144	0	0
3	D	4	Total O 4 4	0	0
3	C	3	Total O 3 3	0	0

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.

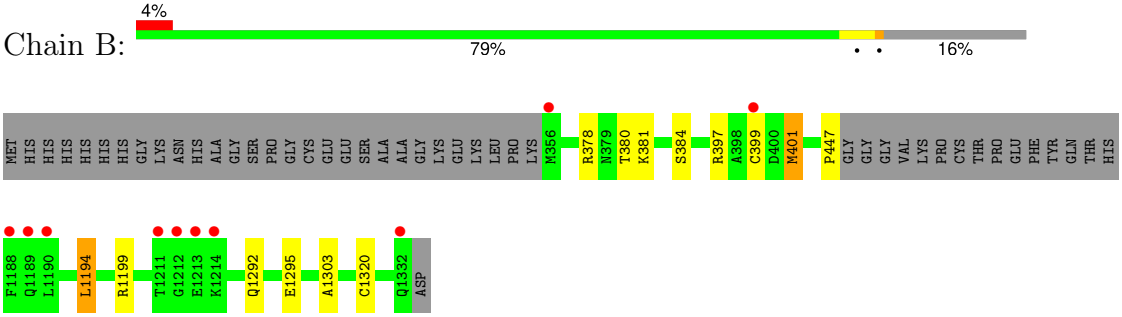
• Molecule 1: JF1cpCasp2

Chain A:



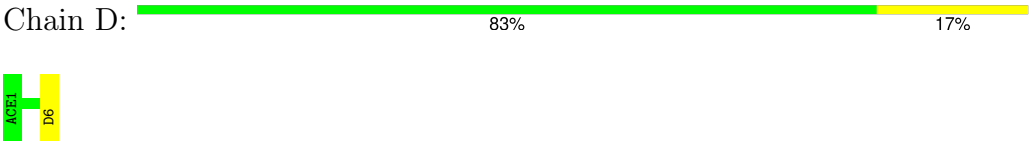
• Molecule 1: JF1cpCasp2

Chain B:



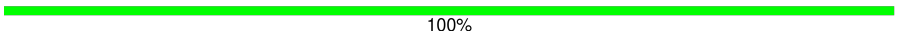
• Molecule 2: AcVDVAD-CHO

Chain D:



• Molecule 2: AcVDVAD-CHO

Chain C:



There are no outlier residues recorded for this chain.

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	55.87Å 101.55Å 112.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.70 – 1.51 37.70 – 1.51	Depositor EDS
% Data completeness (in resolution range)	68.6 (37.70-1.51) 64.6 (37.70-1.51)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	0.92 (at 1.50Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.188 , 0.214 0.188 , 0.215	Depositor DCC
R_{free} test set	5048 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	14.3	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 37.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	4058	wwPDB-VP
Average B, all atoms (Å^2)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.64% of the height of the origin peak. No significant pseudotranslation is detected.*

[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: ASA, ACE

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	0.40	0/1882	0.64	0/2542
1	B	0.37	0/1884	0.62	0/2545
2	C	0.33	0/27	0.62	0/37
2	D	0.31	0/27	0.56	0/37
All	All	0.38	0/3820	0.63	0/5161

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	1847	0	1781	13	0
1	B	1850	0	1785	12	0
2	C	38	0	33	0	0
2	D	38	0	33	4	0
3	A	134	0	0	1	0
3	B	144	0	0	0	0
3	C	3	0	0	0	0
3	D	4	0	0	0	0
All	All	4058	0	3632	24	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 (24) 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:1320:CYS:SG	2:D:6:ASA:C	2.03	1.45
1:A:386:TYR:HB3	1:A:1318:GLN:HE22	1.50	0.76
1:B:1320:CYS:SG	2:D:6:ASA:CA	2.81	0.69
1:A:1247:THR:O	1:A:1251:MET:HG3	1.92	0.69
1:B:397:ARG:HB3	1:B:401:MET:HE3	1.77	0.64
1:B:1320:CYS:SG	2:D:6:ASA:O	2.53	0.62
1:B:1194:LEU:HD22	1:B:1199:ARG:HD2	1.83	0.61
1:A:397:ARG:HB3	1:A:401:MET:HE3	1.90	0.54
1:B:378:ARG:HA	1:B:384:SER:HA	1.95	0.49
1:A:401:MET:HE1	1:A:409:LYS:HD2	1.96	0.48
1:B:380:THR:C	1:B:381:LYS:HD3	2.34	0.48
1:A:1265:ARG:HD3	1:A:1307[A]:SER:CB	2.43	0.48
1:B:1292:GLN:HB2	1:B:1295:GLU:HG3	1.93	0.48
1:A:386:TYR:HB3	1:A:1318:GLN:NE2	2.24	0.48
1:B:399[B]:CYS:SG	1:B:447:PRO:HG2	2.54	0.47
1:B:381:LYS:HD3	1:B:381:LYS:N	2.31	0.46
1:A:357:ARG:HD3	3:A:1404:HOH:O	2.15	0.46
1:A:387:ILE:HG12	1:A:1318:GLN:HE21	1.81	0.44
1:B:1320:CYS:SG	2:D:6:ASA:N	2.91	0.43
1:A:401:MET:HE2	1:A:401:MET:HB2	1.88	0.43
1:A:1306:PRO:HA	1:A:1309:GLN:HG3	1.99	0.43
1:A:378:ARG:HA	1:A:384:SER:HA	2.01	0.42
1:A:1195:GLN:NE2	1:A:1268:ASP:OD2	2.53	0.41
1:A:1329:VAL:HG11	1:B:1303:ALA:HB2	2.03	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	236/282 (84%)	232 (98%)	4 (2%)	0	100	100
1	B	237/282 (84%)	232 (98%)	5 (2%)	0	100	100
2	C	4/6 (67%)	4 (100%)	0	0	100	100
2	D	4/6 (67%)	4 (100%)	0	0	100	100
All	All	481/576 (84%)	472 (98%)	9 (2%)	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	195/238 (82%)	195 (100%)	0	100	100
1	B	195/238 (82%)	193 (99%)	2 (1%)	73	52
2	C	3/3 (100%)	3 (100%)	0	100	100
2	D	3/3 (100%)	3 (100%)	0	100	100
All	All	396/482 (82%)	394 (100%)	2 (0%)	86	75

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	401	MET
1	B	1194	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1318	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)

2 non-standard protein/DNA/RNA residues 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	ASA	C	6	2,1	6,7,7	1.12	0	4,8,8	1.11	0
2	ASA	D	6	2	6,7,7	1.08	0	4,8,8	0.88	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	ASA	C	6	2,1	-	3/5/6/6	-
2	ASA	D	6	2	-	3/5/6/6	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	6	ASA	O-C-CA-CB
2	D	6	ASA	C-CA-CB-CG
2	C	6	ASA	O-C-CA-CB
2	C	6	ASA	C-CA-CB-CG
2	D	6	ASA	N-CA-CB-CG
2	C	6	ASA	N-CA-CB-CG

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	6	ASA	4	0

5.5 Carbohydrates (i)

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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^{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	OWAB(Å^2)	Q<0.9
1	A	237/282 (84%)	0.12	8 (3%) 48 53	7, 18, 32, 47	3 (1%)
1	B	237/282 (84%)	0.14	10 (4%) 41 46	7, 18, 33, 55	4 (1%)
2	C	4/6 (66%)	0.47	0 100 100	19, 20, 20, 22	0
2	D	4/6 (66%)	0.41	0 100 100	17, 18, 19, 24	0
All	All	482/576 (83%)	0.14	18 (3%) 45 51	7, 18, 32, 55	7 (1%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1188	PHE	6.0
1	B	1190	LEU	5.8
1	B	1188	PHE	5.2
1	A	1190	LEU	4.6
1	B	1212	GLY	3.6
1	B	1214	LYS	3.6
1	A	1189	GLN	3.3
1	B	1189	GLN	3.3
1	B	1213	GLU	3.1
1	B	1211	THR	3.0
1	B	356	MET	2.8
1	A	1215	GLU	2.7
1	A	423	GLY	2.5
1	B	399[A]	CYS	2.5
1	A	1332	GLN	2.4
1	A	1187	HIS	2.4
1	B	1332	GLN	2.2
1	A	1323	ASP	2.1

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, 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	ASA	D	6	8/8	0.95	0.07	20,21,24,25	0
2	ASA	C	6	8/8	0.95	0.07	19,21,22,25	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.