Full wwPDB X-ray Structure Validation Report (i)

Mar 5, 2024 – 03:34 PM EST

PDB ID	:	3AQJ
Title	:	Crystal Structure of a C-terminal domain of the bacteriophage P2 tail spike protein, gpV
Authors	:	Takeda, S.; Yamashita, E.; Nakagawa, A.
Deposited on	:	2010-11-06
Resolution	:	1.27 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:

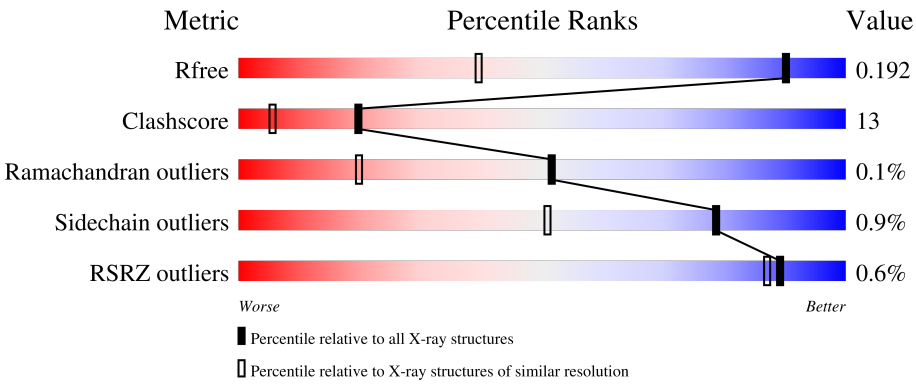
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation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

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.27 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}	130704	1850 (1.30-1.26)
Clashscore	141614	1926 (1.30-1.26)
Ramachandran outliers	138981	1860 (1.30-1.26)
Sidechain outliers	138945	1859 (1.30-1.26)
RSRZ outliers	127900	1807 (1.30-1.26)

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	134	% 74% 10% • 13%
1	B	134	% 66% 14% • 15%
1	C	134	69% 16% •• 13%
1	P	134	% 71% 13% • 15%
1	Q	134	% 73% 12% • 13%

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Mol	Chain	Length	Quality of chain
1	R	134	69% 16% • 13%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 6102 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 Baseplate assembly protein V.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	117	902	548	158	191	5	0	19	0
1	B	114	852	523	149	177	3	0	8	0
1	C	117	871	535	151	182	3	0	11	0
1	P	114	859	525	151	180	3	0	10	0
1	Q	116	859	528	152	176	3	0	7	0
1	R	117	857	525	151	178	3	0	6	0

There are 54 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP P31340
A	2	HIS	-	expression tag	UNP P31340
A	3	HIS	-	expression tag	UNP P31340
A	4	HIS	-	expression tag	UNP P31340
A	5	HIS	-	expression tag	UNP P31340
A	6	HIS	-	expression tag	UNP P31340
A	7	HIS	-	expression tag	UNP P31340
A	8	HIS	-	expression tag	UNP P31340
A	9	HIS	-	expression tag	UNP P31340
B	1	MET	-	expression tag	UNP P31340
B	2	HIS	-	expression tag	UNP P31340
B	3	HIS	-	expression tag	UNP P31340
B	4	HIS	-	expression tag	UNP P31340
B	5	HIS	-	expression tag	UNP P31340
B	6	HIS	-	expression tag	UNP P31340
B	7	HIS	-	expression tag	UNP P31340
B	8	HIS	-	expression tag	UNP P31340

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Chain	Residue	Modelled	Actual	Comment	Reference
B	9	HIS	-	expression tag	UNP P31340
C	1	MET	-	expression tag	UNP P31340
C	2	HIS	-	expression tag	UNP P31340
C	3	HIS	-	expression tag	UNP P31340
C	4	HIS	-	expression tag	UNP P31340
C	5	HIS	-	expression tag	UNP P31340
C	6	HIS	-	expression tag	UNP P31340
C	7	HIS	-	expression tag	UNP P31340
C	8	HIS	-	expression tag	UNP P31340
C	9	HIS	-	expression tag	UNP P31340
P	1	MET	-	expression tag	UNP P31340
P	2	HIS	-	expression tag	UNP P31340
P	3	HIS	-	expression tag	UNP P31340
P	4	HIS	-	expression tag	UNP P31340
P	5	HIS	-	expression tag	UNP P31340
P	6	HIS	-	expression tag	UNP P31340
P	7	HIS	-	expression tag	UNP P31340
P	8	HIS	-	expression tag	UNP P31340
P	9	HIS	-	expression tag	UNP P31340
Q	1	MET	-	expression tag	UNP P31340
Q	2	HIS	-	expression tag	UNP P31340
Q	3	HIS	-	expression tag	UNP P31340
Q	4	HIS	-	expression tag	UNP P31340
Q	5	HIS	-	expression tag	UNP P31340
Q	6	HIS	-	expression tag	UNP P31340
Q	7	HIS	-	expression tag	UNP P31340
Q	8	HIS	-	expression tag	UNP P31340
Q	9	HIS	-	expression tag	UNP P31340
R	1	MET	-	expression tag	UNP P31340
R	2	HIS	-	expression tag	UNP P31340
R	3	HIS	-	expression tag	UNP P31340
R	4	HIS	-	expression tag	UNP P31340
R	5	HIS	-	expression tag	UNP P31340
R	6	HIS	-	expression tag	UNP P31340
R	7	HIS	-	expression tag	UNP P31340
R	8	HIS	-	expression tag	UNP P31340
R	9	HIS	-	expression tag	UNP P31340

- Molecule 2 is FE (II) ION (three-letter code: FE2) (formula: Fe).

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Fe 1 1	0	0
2	P	1	Total Fe 1 1	0	0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Cl 1 1	0	0
3	P	1	Total Cl 1 1	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	1	Total Ca 1 1	0	0
4	P	1	Total Ca 1 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	138	Total O 138 138	0	0
5	B	142	Total O 142 142	0	0
5	C	155	Total O 155 155	0	0
5	P	159	Total O 159 159	0	0
5	Q	145	Total O 145 145	0	0
5	R	157	Total O 157 157	0	0

3 Residue-property plots (i)

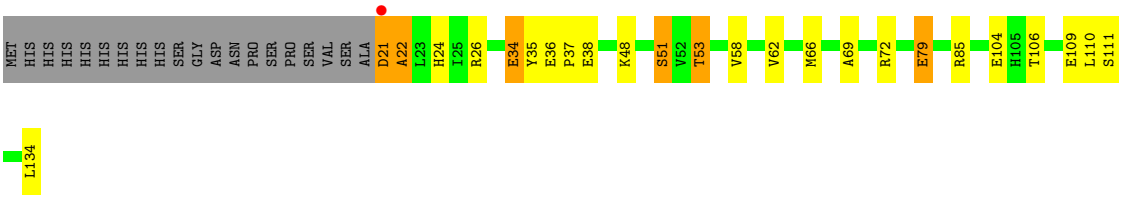
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Baseplate assembly protein V

Chain A:

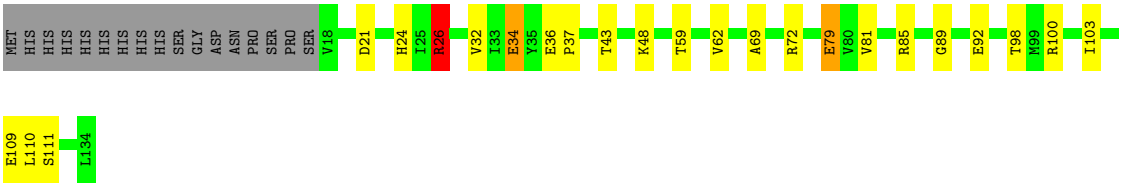
• Molecule 1: Baseplate assembly protein V

Chain B:



• Molecule 1: Baseplate assembly protein V

Chain C:

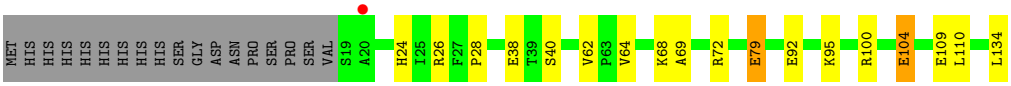


• Molecule 1: Baseplate assembly protein V

Chain P:

• Molecule 1: Baseplate assembly protein V

Chain Q:



• Molecule 1: Baseplate assembly protein V

Chain R:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	66.91Å 64.46Å 67.19Å 90.00° 91.32° 90.00°	Depositor
Resolution (Å)	46.42 – 1.27 46.42 – 1.27	Depositor EDS
% Data completeness (in resolution range)	99.5 (46.42-1.27) 99.5 (46.42-1.27)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	3.21 (at 1.27Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.160 , 0.190 0.162 , 0.192	Depositor DCC
R_{free} test set	7523 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å^2)	9.0	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 48.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.000 for l,k,-h 0.014 for -h,-l,-k 0.000 for -h,l,k 0.002 for k,h,-l 0.000 for -k,-h,-l 0.000 for l,h,k 0.000 for k,l,h 0.000 for -l,-h,k 0.000 for -k,-l,h 0.015 for h,-k,-l 0.011 for l,-k,h	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	6102	wwPDB-VP
Average B, all atoms (Å^2)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.43% 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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: CL, CA, FE2

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	1.53	12/1008 (1.2%)	1.25	3/1369 (0.2%)
1	B	1.46	7/900 (0.8%)	1.37	9/1222 (0.7%)
1	C	1.49	7/938 (0.7%)	1.20	3/1275 (0.2%)
1	P	1.49	4/909 (0.4%)	1.24	7/1236 (0.6%)
1	Q	1.42	5/902 (0.6%)	1.11	2/1226 (0.2%)
1	R	1.55	9/889 (1.0%)	1.27	4/1208 (0.3%)
All	All	1.49	44/5546 (0.8%)	1.24	28/7536 (0.4%)

All (44) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	P	111[A]	SER	CB-OG	-10.82	1.28	1.42
1	P	111[B]	SER	CB-OG	-10.82	1.28	1.42
1	B	51	SER	CB-OG	-9.41	1.30	1.42
1	R	111[A]	SER	CB-OG	-8.44	1.31	1.42
1	R	111[B]	SER	CB-OG	-8.44	1.31	1.42
1	C	111[A]	SER	CB-OG	-7.63	1.32	1.42
1	C	111[B]	SER	CB-OG	-7.63	1.32	1.42
1	A	34	GLU	CB-CG	-7.36	1.38	1.52
1	B	51	SER	CA-CB	-7.21	1.42	1.52
1	R	51	SER	CB-OG	-7.09	1.33	1.42
1	C	34	GLU	CB-CG	-7.02	1.38	1.52
1	A	45[A]	SER	CB-OG	6.79	1.51	1.42
1	A	45[B]	SER	CB-OG	6.79	1.51	1.42
1	A	45[C]	SER	CB-OG	6.79	1.51	1.42
1	C	32	VAL	CB-CG1	6.72	1.67	1.52
1	R	79	GLU	CB-CG	-6.56	1.39	1.52
1	B	34	GLU	CD-OE1	6.53	1.32	1.25
1	P	51	SER	CA-CB	-6.30	1.43	1.52
1	B	111[A]	SER	CB-OG	-6.16	1.34	1.42
1	B	111[B]	SER	CB-OG	-6.16	1.34	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	89	GLY	N-CA	6.08	1.55	1.46
1	R	40	SER	CB-OG	5.96	1.50	1.42
1	Q	79	GLU	CD-OE1	-5.91	1.19	1.25
1	Q	38	GLU	CD-OE2	5.89	1.32	1.25
1	B	58	VAL	CB-CG1	5.69	1.64	1.52
1	R	55	SER	CB-OG	-5.61	1.34	1.42
1	B	35	TYR	CG-CD2	5.39	1.46	1.39
1	R	34	GLU	CB-CG	-5.33	1.42	1.52
1	A	109[A]	GLU	CG-CD	5.33	1.59	1.51
1	A	109[B]	GLU	CG-CD	5.33	1.59	1.51
1	A	109[A]	GLU	CB-CG	5.25	1.62	1.52
1	A	109[B]	GLU	CB-CG	5.25	1.62	1.52
1	A	111[A]	SER	CB-OG	-5.25	1.35	1.42
1	A	111[B]	SER	CB-OG	-5.25	1.35	1.42
1	C	79[A]	GLU	CD-OE1	5.25	1.31	1.25
1	C	79[B]	GLU	CD-OE1	5.25	1.31	1.25
1	Q	64	VAL	CB-CG1	-5.24	1.41	1.52
1	R	24	HIS	C-O	5.19	1.33	1.23
1	P	34	GLU	CB-CG	-5.18	1.42	1.52
1	Q	40	SER	CB-OG	5.11	1.48	1.42
1	R	51	SER	CA-CB	-5.08	1.45	1.52
1	Q	104	GLU	CD-OE1	-5.06	1.20	1.25
1	A	43[A]	THR	CB-OG1	-5.06	1.33	1.43
1	A	43[B]	THR	CB-OG1	-5.06	1.33	1.43

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	53[A]	THR	CA-CB-CG2	-12.10	95.46	112.40
1	B	53[B]	THR	CA-CB-CG2	-12.10	95.46	112.40
1	P	66	MET	CG-SD-CE	-11.85	81.24	100.20
1	R	85	ARG	NE-CZ-NH1	-9.47	115.57	120.30
1	P	29	ASP	CB-CG-OD2	-7.78	111.30	118.30
1	A	85[A]	ARG	NE-CZ-NH2	-7.35	116.63	120.30
1	A	85[B]	ARG	NE-CZ-NH2	-7.35	116.63	120.30
1	B	22	ALA	CB-CA-C	-6.92	99.73	110.10
1	B	72	ARG	NE-CZ-NH2	-6.12	117.24	120.30
1	A	26	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	P	52[A]	VAL	CA-CB-CG2	-5.81	102.18	110.90
1	P	52[B]	VAL	CA-CB-CG2	-5.81	102.18	110.90
1	Q	92	GLU	OE1-CD-OE2	5.77	130.22	123.30
1	R	72	ARG	NE-CZ-NH2	-5.74	117.43	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	R	19	SER	N-CA-CB	5.71	119.07	110.50
1	R	27	PHE	CB-CG-CD1	-5.71	116.81	120.80
1	B	134	LEU	CA-CB-CG	-5.41	102.85	115.30
1	C	26	ARG	NE-CZ-NH1	5.33	122.97	120.30
1	B	79[A]	GLU	CA-CB-CG	5.26	124.97	113.40
1	B	79[B]	GLU	CA-CB-CG	5.26	124.97	113.40
1	P	100	ARG	NE-CZ-NH1	-5.17	117.71	120.30
1	P	100	ARG	NE-CZ-NH2	5.17	122.89	120.30
1	P	100	ARG	CG-CD-NE	5.11	122.53	111.80
1	Q	95	LYS	CD-CE-NZ	5.09	123.42	111.70
1	B	53[A]	THR	OG1-CB-CG2	5.04	121.59	110.00
1	B	53[B]	THR	OG1-CB-CG2	5.04	121.59	110.00
1	C	79[A]	GLU	CG-CD-OE1	5.01	128.33	118.30
1	C	79[B]	GLU	CG-CD-OE1	5.01	128.33	118.30

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	902	0	918	30	0
1	B	852	0	866	31	1
1	C	871	0	892	32	0
1	P	859	0	867	23	0
1	Q	859	0	878	39	0
1	R	857	0	864	29	1
2	A	1	0	0	0	0
2	P	1	0	0	0	0
3	A	1	0	0	0	0
3	P	1	0	0	0	0
4	C	1	0	0	0	0
4	P	1	0	0	0	0
5	A	138	0	0	10	0
5	B	142	0	0	12	3
5	C	155	0	0	11	5

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	P	159	0	0	9	2
5	Q	145	0	0	7	0
5	R	157	0	0	9	5
All	All	6102	0	5285	135	10

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 13.

All (135) 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:R:121[B]:LYS:CE	1:R:121[B]:LYS:NZ	1.68	1.56
1:A:79[B]:GLU:OE2	1:C:72:ARG:NE	1.72	1.22
1:P:72[B]:ARG:NH1	5:P:850:HOH:O	1.72	1.20
1:A:79[B]:GLU:OE2	1:C:72:ARG:CZ	1.91	1.16
1:Q:72[B]:ARG:NH1	5:Q:382:HOH:O	1.79	1.16
1:Q:72[A]:ARG:NE	1:R:79:GLU:OE2	1.80	1.13
1:Q:68[B]:LYS:HE2	5:R:307:HOH:O	1.50	1.11
1:Q:72[A]:ARG:CZ	1:R:79:GLU:OE2	1.99	1.10
1:A:92[B]:GLU:OE2	5:A:835:HOH:O	1.67	1.09
1:A:21[B]:ASP:OD1	5:A:367:HOH:O	1.72	1.07
1:A:79[B]:GLU:OE2	1:C:72:ARG:NH2	1.90	1.05
1:Q:72[A]:ARG:NH2	1:R:79:GLU:OE2	1.91	1.04
1:A:21[A]:ASP:HB2	5:A:795:HOH:O	1.56	1.01
1:A:21[B]:ASP:HB3	5:A:795:HOH:O	1.60	1.01
1:C:48[A]:LYS:NZ	5:C:138:HOH:O	1.97	0.97
1:R:36:GLU:HG2	5:R:830:HOH:O	1.61	0.96
1:A:38:GLU:OE2	1:Q:100:ARG:HD2	1.66	0.96
1:A:110[B]:LEU:HD23	1:C:103:ILE:HB	1.48	0.94
1:P:26:ARG:HH11	1:P:26:ARG:HG2	1.33	0.93
1:A:45[A]:SER:OG	1:B:53[A]:THR:OG1	1.88	0.90
1:A:38:GLU:CD	1:Q:100:ARG:HD2	1.91	0.89
1:B:24:HIS:NE2	1:B:26:ARG:NH2	2.23	0.86
1:C:92:GLU:HB2	5:C:438:HOH:O	1.74	0.86
1:P:104[B]:GLU:OE1	5:P:150:HOH:O	1.93	0.86
1:C:26:ARG:CZ	5:C:556:HOH:O	2.25	0.83
5:A:558:HOH:O	1:C:98:THR:HG21	1.81	0.80
1:B:21:ASP:N	5:B:585:HOH:O	2.14	0.79
1:R:121[B]:LYS:NZ	1:R:121[B]:LYS:CD	2.46	0.78
1:A:45[A]:SER:HG	1:B:53[A]:THR:HG1	1.30	0.78
1:B:85[A]:ARG:HD3	5:C:438:HOH:O	1.87	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:72[A]:ARG:HE	1:R:79:GLU:CD	1.90	0.74
1:B:104[B]:GLU:CG	5:B:829:HOH:O	2.36	0.74
1:B:24:HIS:NE2	1:B:26:ARG:CZ	2.52	0.73
1:P:26:ARG:HH11	1:P:26:ARG:CG	2.01	0.73
1:P:72[B]:ARG:HH12	1:R:66:MET:HE2	1.56	0.71
1:C:36[A]:GLU:HG3	5:C:145:HOH:O	1.91	0.70
1:B:104[B]:GLU:HG3	5:B:829:HOH:O	1.91	0.70
5:P:593:HOH:O	1:Q:68[A]:LYS:HD2	1.90	0.70
1:Q:109[A]:GLU:OE2	5:Q:391:HOH:O	2.11	0.69
1:B:104[B]:GLU:OE2	5:B:829:HOH:O	2.08	0.69
1:P:79:GLU:OE2	5:P:581:HOH:O	2.12	0.68
1:P:21:ASP:HB3	1:Q:28:PRO:HG3	1.76	0.68
1:Q:72[A]:ARG:HG3	1:R:79:GLU:HG3	1.75	0.67
1:R:104:GLU:HG3	5:R:899:HOH:O	1.94	0.66
1:A:72[B]:ARG:NE	5:A:179:HOH:O	2.29	0.65
1:A:110[B]:LEU:HD23	1:C:103:ILE:CB	2.27	0.63
1:B:24:HIS:ND1	5:B:751:HOH:O	2.17	0.62
1:A:38:GLU:OE1	1:Q:100:ARG:CD	2.48	0.62
1:C:34:GLU:CG	1:C:43[B]:THR:OG1	2.48	0.62
1:A:87:ILE:HD12	1:C:79[B]:GLU:OE2	2.00	0.61
1:C:26:ARG:CG	1:C:26:ARG:HH11	2.14	0.60
1:Q:72[B]:ARG:HG3	5:Q:382:HOH:O	2.00	0.60
1:Q:109[B]:GLU:CG	5:Q:391:HOH:O	2.49	0.60
1:Q:24:HIS:NE2	1:Q:26:ARG:NE	2.50	0.60
1:C:43[B]:THR:HG22	5:C:799:HOH:O	2.01	0.59
1:R:104:GLU:CG	5:R:899:HOH:O	2.48	0.59
1:P:21:ASP:CB	1:Q:28:PRO:HG3	2.32	0.59
5:A:140:HOH:O	1:B:53[A]:THR:HG23	2.01	0.59
1:R:84:ASN:OD1	5:R:601:HOH:O	2.17	0.59
1:Q:100:ARG:NH2	5:Q:541:HOH:O	2.33	0.58
1:A:79[B]:GLU:CD	1:C:72:ARG:HE	2.05	0.58
1:C:109[A]:GLU:HG3	5:C:603:HOH:O	2.04	0.58
1:A:38:GLU:OE1	1:Q:100:ARG:HD2	2.04	0.57
1:P:90:THR:HG23	5:R:601:HOH:O	2.04	0.57
1:R:24:HIS:HE1	5:R:456:HOH:O	1.86	0.57
1:Q:134:LEU:HD12	1:R:121[B]:LYS:HD2	1.85	0.57
1:R:95:LYS:HE2	5:R:821:HOH:O	2.04	0.57
1:B:85[A]:ARG:NH1	5:C:438:HOH:O	2.30	0.57
1:B:38:GLU:OE2	5:B:513:HOH:O	2.17	0.56
1:B:109[A]:GLU:OE2	5:B:854:HOH:O	2.18	0.56
1:Q:134:LEU:HD12	1:R:121[B]:LYS:HG3	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:21:ASP:N	5:B:844:HOH:O	2.39	0.55
1:Q:134:LEU:HD12	1:R:121[B]:LYS:CG	2.37	0.54
1:A:72[A]:ARG:NE	5:A:495:HOH:O	2.40	0.54
1:C:43[B]:THR:CG2	5:C:799:HOH:O	2.55	0.54
1:B:48:LYS:CE	1:Q:79:GLU:OE1	2.56	0.54
1:P:69:ALA:HA	1:R:62:VAL:O	2.08	0.54
1:P:72[B]:ARG:HH12	1:R:66:MET:CE	2.20	0.54
1:P:62:VAL:O	1:Q:69:ALA:HA	2.08	0.53
1:A:38:GLU:CD	1:Q:100:ARG:CD	2.72	0.53
1:Q:134:LEU:HD12	1:R:121[B]:LYS:CD	2.38	0.53
1:B:36:GLU:OE2	1:B:38:GLU:HG3	2.09	0.52
1:P:26:ARG:HB2	5:P:545:HOH:O	2.08	0.52
1:Q:62:VAL:O	1:R:69:ALA:HA	2.10	0.52
1:B:53[A]:THR:HG21	5:B:777:HOH:O	2.08	0.52
1:A:69:ALA:HA	1:C:62:VAL:O	2.10	0.52
1:A:21[A]:ASP:OD1	5:A:535:HOH:O	2.19	0.51
1:P:21:ASP:N	5:P:772:HOH:O	2.43	0.51
1:C:34:GLU:HG2	1:C:43[B]:THR:OG1	2.11	0.50
1:A:79[B]:GLU:HG3	1:C:72:ARG:HG3	1.94	0.49
1:B:34:GLU:OE1	5:B:818:HOH:O	2.20	0.49
1:A:38:GLU:OE1	1:Q:100:ARG:HD3	2.12	0.48
1:C:26:ARG:HH11	1:C:26:ARG:HG3	1.78	0.48
1:A:62:VAL:O	1:B:69:ALA:HA	2.13	0.48
1:Q:134:LEU:CD1	1:R:121[B]:LYS:HD2	2.44	0.48
1:C:43[B]:THR:HG23	5:C:155:HOH:O	2.13	0.48
1:P:26:ARG:CG	1:P:26:ARG:NH1	2.71	0.48
1:P:21:ASP:HB2	1:Q:28:PRO:CG	2.43	0.48
1:C:79[B]:GLU:OE2	1:C:81:VAL:CG2	2.62	0.47
1:Q:100:ARG:NE	5:Q:541:HOH:O	2.41	0.47
1:C:100:ARG:NE	1:R:38:GLU:HG3	2.30	0.46
1:P:24:HIS:HB2	5:P:683:HOH:O	2.14	0.46
1:P:43[A]:THR:HG23	5:P:146:HOH:O	2.14	0.46
1:B:79[A]:GLU:OE2	1:C:85:ARG:NH1	2.49	0.45
1:P:34:GLU:CG	1:P:43[A]:THR:OG1	2.64	0.45
1:B:21:ASP:O	1:B:37:PRO:HG2	2.16	0.45
1:Q:104:GLU:OE1	1:R:111[A]:SER:OG	2.30	0.45
1:A:34:GLU:CG	1:A:43[A]:THR:OG1	2.65	0.45
1:C:26:ARG:NE	5:C:556:HOH:O	2.47	0.45
1:B:21:ASP:CA	5:B:585:HOH:O	2.63	0.44
1:C:21:ASP:O	1:C:37:PRO:HG2	2.19	0.43
1:B:62:VAL:O	1:C:69:ALA:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:HIS:CD2	1:B:26:ARG:CZ	3.02	0.42
1:C:24:HIS:HB3	1:C:26:ARG:NH1	2.33	0.42
1:P:72[B]:ARG:HG3	5:P:850:HOH:O	2.18	0.42
1:R:92[A]:GLU:OE2	5:R:578:HOH:O	2.22	0.42
1:P:102:ASN:HB3	1:P:104[B]:GLU:OE2	2.20	0.42
1:A:72[B]:ARG:HG3	5:A:179:HOH:O	2.19	0.42
1:P:110:LEU:HD21	1:Q:110[B]:LEU:HD11	2.02	0.42
1:Q:134:LEU:CD1	1:R:121[B]:LYS:CD	2.98	0.41
1:B:48:LYS:HE3	1:Q:79:GLU:OE1	2.20	0.41
1:P:21:ASP:CB	1:Q:28:PRO:CG	2.98	0.41
1:B:51:SER:HA	1:C:59:THR:O	2.19	0.41
1:Q:72[A]:ARG:NE	1:R:79:GLU:CD	2.57	0.41
1:R:23:LEU:HD11	1:R:25:ILE:HD11	2.02	0.41
1:A:59:THR:HA	1:B:66:MET:O	2.21	0.41
1:B:38:GLU:HG2	5:B:513:HOH:O	2.20	0.41
1:Q:134:LEU:CD1	1:R:121[B]:LYS:HG3	2.49	0.41
1:Q:109[B]:GLU:CD	5:Q:391:HOH:O	2.59	0.41

All (10) 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)
1:B:21:ASP:N	1:R:18:VAL:N[1_455]	1.90	0.30
5:B:654:HOH:O	5:R:480:HOH:O[1_556]	1.90	0.30
5:R:633:HOH:O	5:R:831:HOH:O[2_746]	1.96	0.24
5:B:884:HOH:O	5:P:772:HOH:O[1_455]	1.97	0.23
5:C:875:HOH:O	5:R:601:HOH:O[2_757]	2.03	0.17
5:P:669:HOH:O	5:P:745:HOH:O[2_746]	2.03	0.17
5:C:806:HOH:O	5:R:241:HOH:O[1_455]	2.11	0.09
5:C:238:HOH:O	5:C:614:HOH:O[2_747]	2.13	0.07
5:B:660:HOH:O	5:C:614:HOH:O[2_747]	2.18	0.02
5:C:875:HOH:O	5:R:439:HOH:O[2_757]	2.19	0.01

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	135/134 (101%)	129 (96%)	6 (4%)	0	100	100
1	B	120/134 (90%)	114 (95%)	5 (4%)	1 (1%)	19	2
1	C	126/134 (94%)	123 (98%)	3 (2%)	0	100	100
1	P	122/134 (91%)	119 (98%)	3 (2%)	0	100	100
1	Q	121/134 (90%)	118 (98%)	3 (2%)	0	100	100
1	R	121/134 (90%)	116 (96%)	5 (4%)	0	100	100
All	All	745/804 (93%)	719 (96%)	25 (3%)	1 (0%)	51	20

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	22	ALA

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	115/111 (104%)	112 (97%)	3 (3%)	46	9
1	B	101/111 (91%)	100 (99%)	1 (1%)	76	46
1	C	106/111 (96%)	105 (99%)	1 (1%)	78	51
1	P	102/111 (92%)	101 (99%)	1 (1%)	76	46
1	Q	101/111 (91%)	101 (100%)	0	100	100
1	R	99/111 (89%)	99 (100%)	0	100	100
All	All	624/666 (94%)	618 (99%)	6 (1%)	78	46

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	VAL

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Mol	Chain	Res	Type
1	A	21[A]	ASP
1	A	21[B]	ASP
1	B	21	ASP
1	C	26	ARG
1	P	26	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	ASN
1	B	84	ASN
1	R	24	HIS

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)

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

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 (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	117/134 (87%)	-0.45	1 (0%)	84	82	6, 9, 16, 24	0
1	B	114/134 (85%)	-0.44	1 (0%)	84	82	5, 10, 19, 31	0
1	C	117/134 (87%)	-0.45	0	100	100	6, 9, 17, 22	0
1	P	114/134 (85%)	-0.52	1 (0%)	84	82	6, 9, 20, 38	0
1	Q	116/134 (86%)	-0.49	1 (0%)	84	82	5, 10, 21, 36	0
1	R	117/134 (87%)	-0.46	0	100	100	5, 9, 18, 22	0
All	All	695/804 (86%)	-0.47	4 (0%)	89	87	5, 9, 18, 38	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	P	21	ASP	3.0
1	Q	20	ALA	2.6
1	B	21	ASP	2.6
1	A	22	ALA	2.4

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, 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
3	CL	A	136	1/1	0.99	0.05	8,8,8,8	0
4	CA	C	135	1/1	0.99	0.03	9,9,9,9	0
2	FE2	A	135	1/1	1.00	0.04	5,5,5,5	0
3	CL	P	137	1/1	1.00	0.03	8,8,8,8	0
2	FE2	P	135	1/1	1.00	0.04	4,4,4,4	0
4	CA	P	136	1/1	1.00	0.04	8,8,8,8	0

6.5 Other polymers (i)

There are no such residues in this entry.