



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 5E4V / pdb_00005e4v
Title : Crystal structure of measles N0-P complex
Authors : Guryanov, S.G.; Liljeroos, L.; Kasaragod, P.; Kajander, T.; Butcher, S.J.
Deposited on : 2015-10-07
Resolution : 2.71 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

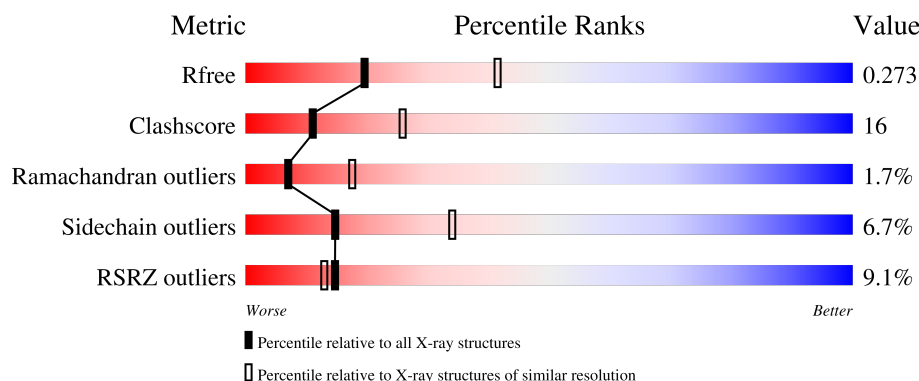
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.71 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}	180053	4348 (2.74-2.70)
Clashscore	190562	4665 (2.74-2.70)
Ramachandran outliers	187476	4584 (2.74-2.70)
Sidechain outliers	187428	4585 (2.74-2.70)
RSRZ outliers	180081	4348 (2.74-2.70)

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 $\leq 5\%$. 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	437	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5817 atoms, of which 2833 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 Nucleoprotein,Phosphoprotein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	407	Total	C	H	N	O	S	0	0	0
			5808	1908	2833	490	563	14			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	SER	-	expression tag	UNP Q89933

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	9	Total	O	0	0
			9	9		

- Molecule 1: Nucleoprotein, Phosphoprotein



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	91.14Å 91.14Å 94.21Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	78.93 – 2.71 78.93 – 2.71	Depositor EDS
% Data completeness (in resolution range)	99.8 (78.93-2.71) 99.8 (78.93-2.71)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 2.73Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.211 , 0.266 0.218 , 0.273	Depositor DCC
R_{free} test set	700 reflections (5.52%)	wwPDB-VP
Wilson B-factor (Å ²)	62.5	Xtriage
Anisotropy	0.369	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 89.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.039 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5817	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

5 Model quality

5.1 Standard geometry

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	0.39	0/3025	0.81	4/4117 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	171	ASP	N-CA-C	-6.90	102.92	111.75
1	A	291	ALA	CA-C-N	-5.51	112.95	119.84
1	A	291	ALA	C-N-CA	-5.51	112.95	119.84
1	A	413	ILE	N-CA-C	5.08	115.22	108.27

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	2975	2833	2830	95	0
2	A	9	0	0	0	0
All	All	2984	2833	2830	95	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 16.

All (95) 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:327:SER:HA	1:A:328:MET:CB	2.00	0.89
1:A:69:ILE:O	1:A:73:THR:HG23	1.79	0.83
1:A:327:SER:HA	1:A:328:MET:HB3	1.59	0.83
1:A:99:SER:C	1:A:102:THR:HA	2.12	0.74
1:A:323:GLU:HG3	1:A:347:GLN:HA	1.69	0.74
1:A:165:ALA:HB3	1:A:166:PRO:HD2	1.70	0.72
1:A:226:GLY:CA	1:A:227:ASN:CB	2.69	0.71
1:A:165:ALA:CB	1:A:166:PRO:CD	2.69	0.70
1:A:122:GLN:HA	1:A:125:PHE:CD2	2.27	0.69
1:A:130:ASN:HB3	1:A:133:ILE:CD1	2.21	0.69
1:A:150:THR:HG21	1:A:190:PHE:HB2	1.80	0.64
1:A:263:PRO:HG2	1:A:349:MET:HE1	1.80	0.62
1:A:106:ASP:O	1:A:107:GLU:C	2.41	0.62
1:A:323:GLU:HG3	1:A:347:GLN:CA	2.30	0.61
1:A:226:GLY:HA2	1:A:227:ASN:CB	2.30	0.60
1:A:146:MET:SD	1:A:187:VAL:HG21	2.42	0.59
1:A:169:ALA:O	1:A:173:GLU:HG3	2.05	0.57
1:A:130:ASN:OD1	1:A:132:GLU:HB2	2.05	0.57
1:A:122:GLN:HA	1:A:125:PHE:HD2	1.68	0.56
1:A:318:MET:O	1:A:322:VAL:HG13	2.04	0.56
1:A:310:TYR:N	1:A:311:PRO:HD3	2.20	0.56
1:A:123:SER:C	1:A:125:PHE:H	2.15	0.55
1:A:405:ILE:HG22	1:A:406:ARG:N	2.21	0.55
1:A:71:ARG:NH2	1:A:220:ASP:OD1	2.40	0.54
1:A:327:SER:HA	1:A:328:MET:HB2	1.87	0.54
1:A:165:ALA:HB2	1:A:302:GLN:NE2	2.23	0.54
1:A:170:ALA:C	1:A:172:SER:N	2.61	0.54
1:A:122:GLN:HA	1:A:125:PHE:CE2	2.44	0.53
1:A:127:TRP:CG	1:A:128:PHE:H	2.27	0.53
1:A:234:MET:HG3	1:A:235:ILE:N	2.24	0.52
1:A:323:GLU:CG	1:A:347:GLN:HA	2.39	0.51
1:A:67:GLN:O	1:A:71:ARG:HG2	2.09	0.51
1:A:55:ILE:HG23	1:A:213:PHE:CE2	2.46	0.51
1:A:273:GLU:OE1	1:A:313:LEU:HB2	2.10	0.51
1:A:291:ALA:O	1:A:294:MET:HG2	2.10	0.51
1:A:327:SER:CA	1:A:328:MET:CB	2.78	0.51
1:A:121:ASP:O	1:A:122:GLN:C	2.54	0.50
1:A:165:ALA:HB1	1:A:166:PRO:CD	2.40	0.50
1:A:86:VAL:HG22	1:A:87:GLN:N	2.26	0.49
1:A:165:ALA:CB	1:A:166:PRO:HD2	2.34	0.49
1:A:258:ILE:HD12	1:A:258:ILE:N	2.29	0.48
1:A:160:ALA:HA	1:A:242:ILE:HG12	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:421:ALA:O	1:A:424:ALA:HB3	2.14	0.47
1:A:426:SER:OG	1:A:427:GLU:N	2.48	0.47
1:A:83:LEU:HD23	1:A:97:PHE:CE1	2.50	0.47
1:A:409:LYS:O	1:A:410:ALA:O	2.32	0.47
1:A:42:ASN:HB3	1:A:45:VAL:HG23	1.97	0.47
1:A:264:ALA:H	1:A:349:MET:HE2	1.78	0.47
1:A:68:LEU:O	1:A:72:ILE:HG12	2.14	0.46
1:A:236:CYS:O	1:A:240:THR:HG23	2.14	0.46
1:A:121:ASP:O	1:A:123:SER:N	2.49	0.46
1:A:319:GLY:HA2	1:A:343:PHE:CE1	2.51	0.46
1:A:29:ARG:HG2	1:A:144:PHE:CZ	2.50	0.46
1:A:165:ALA:CB	1:A:302:GLN:HE22	2.29	0.45
1:A:324:LEU:HD13	1:A:350:VAL:HG13	1.98	0.45
1:A:97:PHE:CE2	1:A:108:ALA:HA	2.52	0.44
1:A:327:SER:CA	1:A:328:MET:HB2	2.46	0.44
1:A:316:TYR:HA	1:A:342:TYR:O	2.16	0.44
1:A:323:GLU:HG2	1:A:347:GLN:CG	2.48	0.44
1:A:51:THR:HG23	1:A:210:LEU:HB2	1.98	0.44
1:A:16:ILE:HG22	1:A:18:VAL:HG23	2.00	0.43
1:A:98:ALA:HA	1:A:99:SER:HA	1.79	0.43
1:A:260:THR:HB	1:A:262:TYR:CZ	2.53	0.43
1:A:177:TRP:CD1	1:A:181:THR:HG1	2.36	0.43
1:A:346:GLY:O	1:A:350:VAL:HG23	2.19	0.43
1:A:18:VAL:HG22	1:A:52:GLY:HA2	2.01	0.43
1:A:23:ASP:OD1	1:A:25:SER:HB3	2.18	0.43
1:A:291:ALA:O	1:A:294:MET:HE2	2.19	0.43
1:A:30:SER:HA	1:A:144:PHE:CE1	2.54	0.42
1:A:45:VAL:HG13	1:A:49:LYS:HD3	2.00	0.42
1:A:233:GLU:O	1:A:236:CYS:HB3	2.19	0.42
1:A:236:CYS:O	1:A:239:ASP:HB2	2.19	0.42
1:A:177:TRP:CD1	1:A:177:TRP:C	2.97	0.42
1:A:390:MET:HE2	1:A:390:MET:HB2	1.95	0.42
1:A:165:ALA:HB1	1:A:166:PRO:HD3	2.00	0.42
1:A:31:ARG:O	1:A:35:ARG:HG2	2.20	0.42
1:A:319:GLY:N	1:A:343:PHE:CD1	2.87	0.42
1:A:277:LEU:O	1:A:281:MET:HG3	2.20	0.41
1:A:30:SER:HA	1:A:144:PHE:HE1	1.85	0.41
1:A:86:VAL:HG22	1:A:87:GLN:H	1.85	0.41
1:A:123:SER:C	1:A:125:PHE:N	2.76	0.41
1:A:187:VAL:HG12	1:A:189:GLU:H	1.85	0.41
1:A:176:ARG:O	1:A:179:LYS:HB3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:LEU:O	1:A:161:LYS:HB2	2.21	0.41
1:A:153:ALA:HB2	1:A:234:MET:HE1	2.02	0.41
1:A:183:GLN:HA	1:A:184:ARG:HA	1.82	0.41
1:A:130:ASN:HB3	1:A:133:ILE:HD13	1.99	0.41
1:A:340:PRO:O	1:A:343:PHE:HB3	2.21	0.41
1:A:107:GLU:O	1:A:111:TYR:HB2	2.21	0.41
1:A:140:ASP:HA	1:A:141:PRO:HD3	1.94	0.41
1:A:215:VAL:HA	1:A:218:ILE:HD12	2.03	0.40
1:A:348:GLU:O	1:A:352:ARG:HG3	2.20	0.40
1:A:260:THR:HB	1:A:262:TYR:CE2	2.57	0.40
1:A:367:ILE:HG22	1:A:368:THR:N	2.36	0.40
1:A:122:GLN:CD	1:A:125:PHE:CE2	3.00	0.40

There are no symmetry-related clashes.

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	401/437 (92%)	363 (90%)	31 (8%)	7 (2%)	7	17

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	122	GLN
1	A	165	ALA
1	A	410	ALA
1	A	329	GLY
1	A	412	PRO
1	A	227	ASN
1	A	327	SER

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	284/368 (77%)	265 (93%)	19 (7%)	15	34

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	SER
1	A	44	ASP
1	A	46	SER
1	A	71	ARG
1	A	77	ASP
1	A	80	ILE
1	A	122	GLN
1	A	123	SER
1	A	138	VAL
1	A	244	GLU
1	A	269	GLU
1	A	275	SER
1	A	285	GLN
1	A	322	VAL
1	A	336	SER
1	A	359	SER
1	A	368	THR
1	A	405	ILE
1	A	413	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	87	GLN
1	A	302	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](#)

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

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, 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	407/437 (93%)	0.56	37 (9%) 15 13	43, 69, 119, 200	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	242	ILE	5.7
1	A	171	ASP	5.4
1	A	239	ASP	4.8
1	A	385	ASP	4.5
1	A	328	MET	4.1
1	A	241	TYR	4.1
1	A	387	ILE	3.5
1	A	102	THR	3.3
1	A	243	VAL	3.3
1	A	371	ASP	3.2
1	A	104	MET	3.2
1	A	105	GLU	3.1
1	A	240	THR	3.1
1	A	296	ILE	3.1
1	A	390	MET	3.1
1	A	244	GLU	3.0
1	A	413	ILE	3.0
1	A	12	ILE	2.9
1	A	184	ARG	2.8
1	A	386	LYS	2.7
1	A	186	VAL	2.7
1	A	169	ALA	2.6
1	A	297	LEU	2.6
1	A	425	TRP	2.5
1	A	303	ASN	2.5
1	A	227	ASN	2.4
1	A	99	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	103	ASN	2.3
1	A	427	GLU	2.3
1	A	98	ALA	2.3
1	A	183	GLN	2.2
1	A	157	VAL	2.2
1	A	238	ILE	2.2
1	A	180	TYR	2.1
1	A	407	ALA	2.1
1	A	146	MET	2.0
1	A	293	TYR	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 oligosaccharides 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.