



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 10:17 am BST

PDB ID : 2EC0
Title : RNA-dependent RNA polymerase of foot-and-mouth disease virus in complex with a template-primer RNA and ATP
Authors : Ferrer-Orta, C.; Arias, A.; Perez-Luque, R.; Escarmis, C.; Domingo, E.; Verdaguier, N.
Deposited on : 2007-02-09
Resolution : 2.75 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.11
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) : 2.11

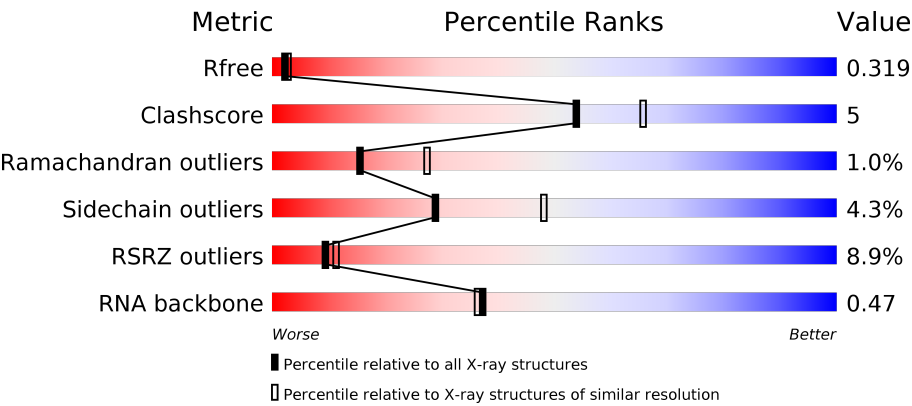
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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.



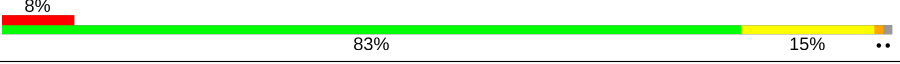

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	130704	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)
RNA backbone	3102	1060 (3.02-2.50)

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 on 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	B	8	<div><div>25%</div><div><div></div><div></div><div></div><div></div></div><div>25%13%63%</div></div>
1	E	8	<div><div>38%</div><div><div></div><div></div><div></div><div></div></div><div>25%38%25%13%</div></div>
2	C	7	<div><div>43%</div><div><div></div><div></div><div></div><div></div></div><div>57%29%14%</div></div>
2	F	7	<div><div>43%</div><div><div></div><div></div><div></div><div></div></div><div>57%29%14%</div></div>

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Mol	Chain	Length	Quality of chain
3	A	476	
3	D	476	

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
5	PPV	A	930	-	-	-	X
5	PPV	D	2930	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8200 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 RNA chain called 5'-R(P*AP*UP*GP*GP*GP*CP*CP*C)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	8	Total	C	N	O	P	0	0	0
			171	76	31	56	8			
1	E	8	Total	C	N	O	P	0	0	0
			171	76	31	56	8			

- Molecule 2 is a RNA chain called 5'-R(*GP*GP*GP*CP*CP*CP*A)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	7	Total	C	N	O	P	0	0	0
			148	67	29	46	6			
2	F	7	Total	C	N	O	P	0	0	0
			148	67	29	46	6			

- Molecule 3 is a protein called RNA-dependent RNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	470	Total	C	N	O	S	0	0	0
			3714	2361	643	689	21			
3	D	470	Total	C	N	O	S	0	0	0
			3714	2361	643	689	21			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	471	ALA	-	CLONING ARTIFACT	UNP Q0QEE1
A	472	ALA	-	CLONING ARTIFACT	UNP Q0QEE1
A	473	LEU	-	CLONING ARTIFACT	UNP Q0QEE1
A	474	GLU	-	CLONING ARTIFACT	UNP Q0QEE1
A	475	HIS	-	CLONING ARTIFACT	UNP Q0QEE1
A	476	HIS	-	CLONING ARTIFACT	UNP Q0QEE1
D	471	ALA	-	CLONING ARTIFACT	UNP Q0QEE1
D	472	ALA	-	CLONING ARTIFACT	UNP Q0QEE1

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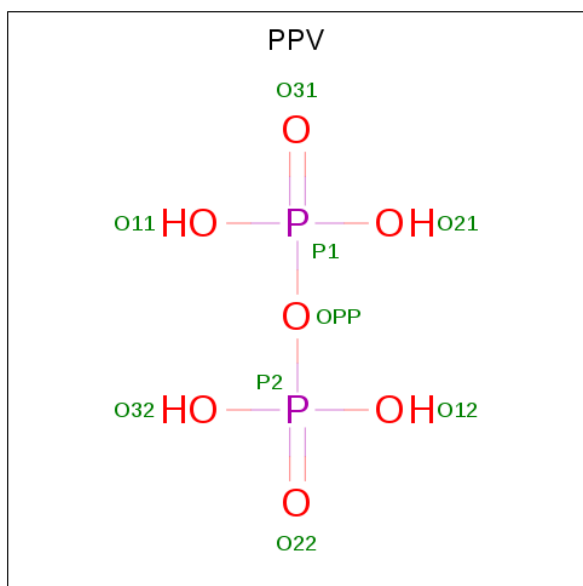
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Chain	Residue	Modelled	Actual	Comment	Reference
D	473	LEU	-	CLONING ARTIFACT	UNP Q0QEE1
D	474	GLU	-	CLONING ARTIFACT	UNP Q0QEE1
D	475	HIS	-	CLONING ARTIFACT	UNP Q0QEE1
D	476	HIS	-	CLONING ARTIFACT	UNP Q0QEE1

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Mg 1 1	0	0
4	D	1	Total Mg 1 1	0	0

- Molecule 5 is PYROPHOSPHATE (three-letter code: PPV) (formula: $H_4O_7P_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total O P 9 7 2	0	0
5	D	1	Total O P 9 7 2	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	3	Total O 3 3	0	0

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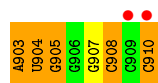
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	2	Total 2	O 2	0	0
6	E	4	Total 4	O 4	0	0
6	F	1	Total 1	O 1	0	0
6	A	49	Total 49	O 49	0	0
6	D	55	Total 55	O 55	0	0

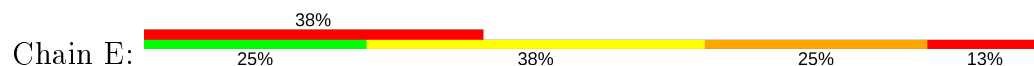
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

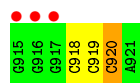
- Molecule 1: 5'-R(P*AP*UP*GP*GP*GP*CP*CP*C)-3'



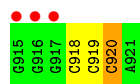
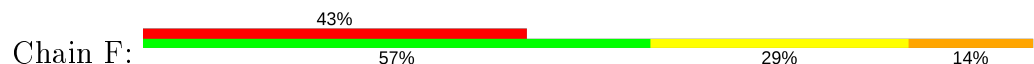
- Molecule 1: 5'-R(P*AP*UP*GP*GP*GP*CP*CP*C)-3'



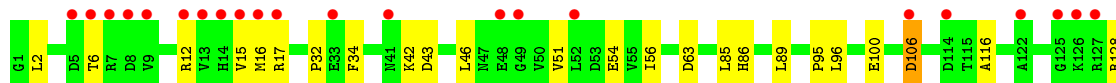
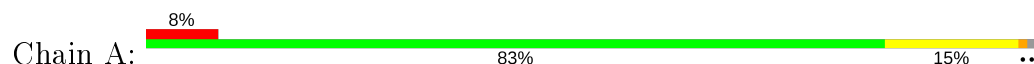
- Molecule 2: 5'-R(*GP*GP*GP*CP*CP*CP*A)-3'

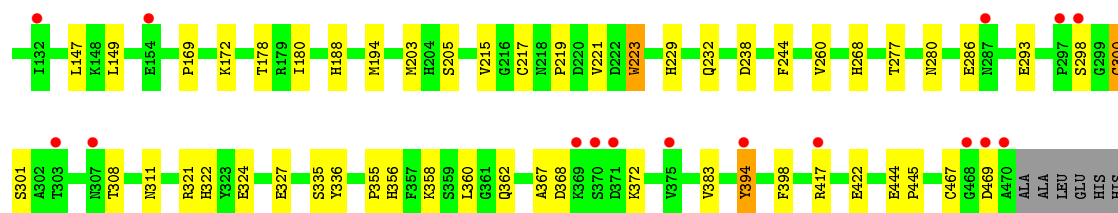


- Molecule 2: 5'-R(*GP*GP*GP*CP*CP*CP*A)-3'

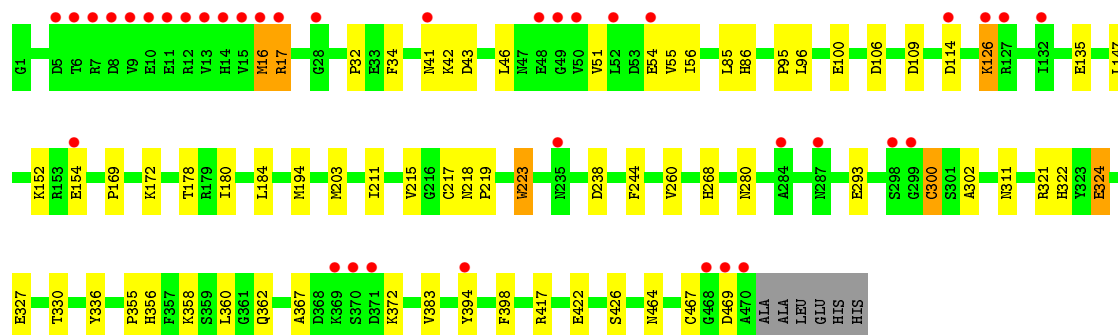
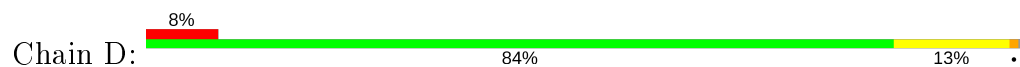


- Molecule 3: RNA-dependent RNA polymerase





● Molecule 3: RNA-dependent RNA polymerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	95.66 Å 95.66 Å 201.46 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.75 29.90 – 2.70	Depositor EDS
% Data completeness (in resolution range)	85.5 (20.00-2.75) 100.0 (29.90-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 2.68 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.245 , 0.295 0.277 , 0.319	Depositor DCC
R_{free} test set	1526 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	55.2	Xtriage
Anisotropy	0.118	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.035 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8200	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 91.29 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.5928e-08. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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

Bond lengths and bond angles in the following residue types are not validated in this section: MG, PPV

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	B	0.70	0/190	1.35	3/294 (1.0%)
1	E	0.69	0/190	1.39	5/294 (1.7%)
2	C	0.59	0/165	1.16	1/256 (0.4%)
2	F	0.59	0/165	1.25	1/256 (0.4%)
3	A	0.42	0/3803	0.55	0/5149
3	D	0.50	2/3803 (0.1%)	0.57	2/5149 (0.0%)
All	All	0.48	2/8316 (0.0%)	0.67	12/11398 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	126	LYS	CD-CE	13.60	1.85	1.51
3	D	126	LYS	CE-NZ	5.13	1.61	1.49

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	920	C	O4'-C1'-N1	7.23	113.98	108.20
2	C	920	C	O4'-C1'-N1	5.73	112.79	108.20
1	E	905	G	C5'-C4'-O4'	5.70	115.94	109.10
1	E	908	C	O4'-C1'-N1	5.54	112.63	108.20
3	D	126	LYS	CD-CE-NZ	-5.39	99.29	111.70
1	B	908	C	O4'-C1'-N1	5.31	112.45	108.20
1	E	905	G	O4'-C4'-C3'	-5.30	98.70	104.00
1	E	905	G	C1'-O4'-C4'	-5.22	105.72	109.90
1	B	903	A	C3'-C2'-C1'	-5.15	97.38	101.50
3	D	126	LYS	CG-CD-CE	-5.07	96.70	111.90
1	E	905	G	C4'-C3'-C2'	-5.05	97.55	102.60
1	B	910	C	O4'-C1'-N1	5.02	112.21	108.20

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	B	171	0	88	2	0
1	E	171	0	88	4	0
2	C	148	0	79	2	0
2	F	148	0	79	3	0
3	A	3714	0	3642	36	3
3	D	3714	0	3642	38	3
4	A	1	0	0	0	0
4	D	1	0	0	0	0
5	A	9	0	0	0	0
5	D	9	0	0	0	0
6	A	49	0	0	0	0
6	B	3	0	0	0	0
6	C	2	0	0	0	0
6	D	55	0	0	3	0
6	E	4	0	0	0	0
6	F	1	0	0	0	0
All	All	8200	0	7618	83	3

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.

All (83) 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)
3:D:126:LYS:CD	3:D:126:LYS:CE	1.85	1.50
3:A:85:LEU:HD11	3:A:203:MET:HE1	1.46	0.95
3:A:51:VAL:HB	3:A:54:GLU:HG2	1.55	0.87
3:A:85:LEU:HD11	3:A:203:MET:CE	2.13	0.78
3:D:126:LYS:CG	3:D:126:LYS:CE	2.62	0.78
3:D:85:LEU:HD11	3:D:203:MET:HE1	1.64	0.77
3:D:85:LEU:HD11	3:D:203:MET:CE	2.19	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:56:ILE:HG23	3:A:180:ILE:HG21	1.75	0.68
3:D:464:ASN:HB2	6:D:2969:HOH:O	1.96	0.65
3:D:238:ASP:HB2	3:D:383:VAL:HG12	1.80	0.63
2:F:918:C:H2'	2:F:919:C:C6	2.32	0.62
3:A:321:ARG:HH11	3:A:356:HIS:HD2	1.47	0.62
3:A:300:CYS:SG	3:A:301:SER:O	2.54	0.61
3:A:398:PHE:HE2	3:A:467:CYS:SG	2.24	0.61
3:D:126:LYS:NZ	3:D:126:LYS:CD	2.64	0.60
3:A:96:LEU:HG	3:A:100:GLU:HB3	1.84	0.60
2:C:918:C:H2'	2:C:919:C:C6	2.38	0.58
3:D:51:VAL:HB	3:D:54:GLU:HG2	1.86	0.56
3:D:302:ALA:HA	6:D:2014:HOH:O	2.05	0.55
3:A:238:ASP:HB2	3:A:383:VAL:HG12	1.89	0.55
3:A:43:ASP:HB3	3:A:46:LEU:HG	1.89	0.54
3:D:56:ILE:HG23	3:D:180:ILE:HG21	1.89	0.54
3:A:116:ALA:O	3:A:128:ARG:NH1	2.40	0.54
3:D:244:PHE:HA	3:D:362:GLN:HE22	1.72	0.53
2:F:919:C:H2'	2:F:920:C:C6	2.43	0.53
3:A:95:PRO:HG3	3:A:268:HIS:HB2	1.92	0.52
3:A:355:PRO:HA	3:A:358:LYS:HB2	1.91	0.52
3:A:238:ASP:HB3	3:A:367:ALA:HB3	1.91	0.51
3:D:96:LEU:HG	3:D:100:GLU:HB3	1.92	0.51
3:D:355:PRO:HA	3:D:358:LYS:HB2	1.91	0.51
3:A:280:ASN:HD22	3:A:293:GLU:HA	1.76	0.51
3:D:321:ARG:HH11	3:D:356:HIS:HD2	1.58	0.50
3:A:56:ILE:HD11	3:A:178:THR:HB	1.93	0.50
3:A:244:PHE:HA	3:A:362:GLN:HE22	1.76	0.49
3:D:169:PRO:HG2	3:D:172:LYS:HD2	1.95	0.49
3:D:238:ASP:HB3	3:D:367:ALA:HB3	1.94	0.49
3:A:229:HIS:O	3:A:232:GLN:HG2	2.13	0.49
3:D:280:ASN:HD22	3:D:293:GLU:HA	1.78	0.49
3:D:322:HIS:ND1	3:D:356:HIS:HE1	2.10	0.48
3:D:43:ASP:HB3	3:D:46:LEU:HG	1.94	0.48
3:D:95:PRO:HG3	3:D:268:HIS:HB2	1.95	0.48
1:E:904:U:H2'	1:E:905:G:O4'	2.14	0.47
3:A:308:THR:HG23	3:A:335:SER:OG	2.13	0.47
3:A:89:LEU:HD11	3:A:203:MET:CE	2.45	0.47
2:C:919:C:H2'	2:C:920:C:C6	2.50	0.47
3:A:169:PRO:HG2	3:A:172:LYS:HD2	1.96	0.47
3:A:12:ARG:HE	3:A:286:GLU:HB3	1.80	0.46
3:A:96:LEU:HB2	3:A:194:MET:HA	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:467:CYS:O	3:D:467:CYS:SG	2.73	0.46
3:A:322:HIS:ND1	3:A:356:HIS:HE1	2.13	0.46
1:B:904:U:H2'	1:B:905:G:O4'	2.16	0.45
1:E:906:G:OP2	3:D:109:ASP:HB2	2.16	0.45
3:D:398:PHE:HE2	3:D:467:CYS:SG	2.39	0.45
3:A:327:GLU:OE1	3:A:327:GLU:HA	2.17	0.45
3:D:96:LEU:HB2	3:D:194:MET:HA	1.99	0.45
3:D:219:PRO:O	3:D:223:TRP:HB2	2.16	0.45
3:A:215:VAL:HA	3:A:336:TYR:CE1	2.51	0.45
1:B:907:G:H2'	1:B:908:C:O4'	2.17	0.45
3:D:215:VAL:HA	3:D:336:TYR:CE1	2.51	0.45
3:D:86:HIS:HE1	3:D:260:VAL:O	1.99	0.45
3:D:184:LEU:HD22	3:D:300:CYS:HB2	1.98	0.44
1:E:907:G:H4'	3:D:218:ASN:HB2	1.99	0.44
3:A:444:GLU:N	3:A:445:PRO:CD	2.80	0.44
3:D:324:GLU:HG3	6:D:2009:HOH:O	2.18	0.44
3:D:56:ILE:HD11	3:D:178:THR:HB	1.99	0.44
3:A:86:HIS:HE1	3:A:260:VAL:O	2.00	0.43
3:A:467:CYS:SG	3:A:467:CYS:O	2.76	0.43
3:D:322:HIS:HB2	3:D:356:HIS:CE1	2.53	0.43
3:D:203:MET:HE1	3:D:211:ILE:HD12	2.01	0.43
3:D:51:VAL:O	3:D:55:VAL:HG23	2.18	0.43
3:D:32:PRO:C	3:D:34:PHE:H	2.22	0.43
3:A:32:PRO:C	3:A:34:PHE:H	2.22	0.42
3:A:298:SER:OG	3:A:298:SER:O	2.36	0.42
1:E:907:G:H2'	1:E:908:C:O4'	2.19	0.42
3:D:85:LEU:HD11	3:D:203:MET:HE3	2.01	0.42
3:A:2:LEU:HB2	3:A:63:ASP:N	2.35	0.41
3:A:219:PRO:O	3:A:223:TRP:HB2	2.20	0.41
3:D:152:LYS:NZ	3:D:154:GLU:HG3	2.36	0.40
3:A:89:LEU:HD11	3:A:203:MET:HE3	2.04	0.40
2:F:918:C:H2'	2:F:919:C:H6	1.81	0.40
3:A:188:HIS:HE1	3:A:277:THR:OG1	2.03	0.40
3:A:2:LEU:HB2	3:A:63:ASP:H	1.87	0.40
3:D:327:GLU:H	3:D:330:THR:HG1	1.68	0.40

All (3) 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:394:TYR:OH	3:D:16:MET:CB[3_664]	1.85	0.35

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:394:TYR:CE2	3:D:16:MET:SD[3_664]	1.89	0.31
3:A:394:TYR:OH	3:D:17:ARG:N[3_664]	2.10	0.10

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	
3	A	468/476 (98%)	444 (95%)	19 (4%)	5 (1%)	14	25
3	D	468/476 (98%)	444 (95%)	20 (4%)	4 (1%)	17	31
All	All	936/952 (98%)	888 (95%)	39 (4%)	9 (1%)	15	27

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	16	MET
3	D	16	MET
3	A	372	LYS
3	D	372	LYS
3	A	106	ASP
3	A	17	ARG
3	D	17	ARG
3	A	394	TYR
3	D	106	ASP

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	
3	A	395/399 (99%)	377 (95%)	18 (5%)	27	46
3	D	395/399 (99%)	379 (96%)	16 (4%)	30	50
All	All	790/798 (99%)	756 (96%)	34 (4%)	29	48

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	6	THR
3	A	15	VAL
3	A	42	LYS
3	A	106	ASP
3	A	147	LEU
3	A	149	LEU
3	A	205	SER
3	A	217	CYS
3	A	221	VAL
3	A	223	TRP
3	A	300	CYS
3	A	311	ASN
3	A	324	GLU
3	A	360	LEU
3	A	368	ASP
3	A	417	ARG
3	A	422	GLU
3	A	469	ASP
3	D	41	ASN
3	D	42	LYS
3	D	114	ASP
3	D	135	GLU
3	D	147	LEU
3	D	217	CYS
3	D	223	TRP
3	D	300	CYS
3	D	311	ASN
3	D	324	GLU
3	D	360	LEU
3	D	394	TYR
3	D	417	ARG
3	D	422	GLU
3	D	426	SER
3	D	469	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	60	HIS
3	A	86	HIS
3	A	188	HIS
3	A	254	ASN
3	A	280	ASN
3	A	311	ASN
3	A	356	HIS
3	A	362	GLN
3	D	60	HIS
3	D	86	HIS
3	D	124	GLN
3	D	188	HIS
3	D	210	GLN
3	D	254	ASN
3	D	280	ASN
3	D	311	ASN
3	D	356	HIS
3	D	362	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	B	8/8 (100%)	3 (37%)	1 (12%)
1	E	8/8 (100%)	2 (25%)	2 (25%)
2	C	6/7 (85%)	0	0
2	F	6/7 (85%)	0	0
All	All	28/30 (93%)	5 (17%)	3 (10%)

All (5) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	B	904	U
1	B	905	G
1	B	910	C
1	E	904	U
1	E	905	G

All (3) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	B	903	A
1	E	903	A
1	E	905	G

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
5	PPV	A	930	-	6,8,8	1.48	2 (33%)	13,13,13	1.51	1 (7%)
5	PPV	D	2930	-	6,8,8	0.70	0	13,13,13	1.15	1 (7%)

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
5	PPV	A	930	-	-	0/6/6/6	-
5	PPV	D	2930	-	-	0/6/6/6	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	930	PPV	P1-O31	2.30	1.58	1.50
5	A	930	PPV	P2-O22	2.29	1.57	1.50

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	930	PPV	P2-OPP-P1	-4.39	117.77	132.83
5	D	2930	PPV	P2-OPP-P1	-2.39	124.63	132.83

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 ⓘ

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	B	8/8 (100%)	1.48	2 (25%) 0 0	55, 64, 81, 87	8 (100%)
1	E	8/8 (100%)	1.65	3 (37%) 0 0	55, 64, 81, 87	8 (100%)
2	C	7/7 (100%)	1.81	3 (42%) 0 0	42, 48, 76, 82	7 (100%)
2	F	7/7 (100%)	1.81	3 (42%) 0 0	42, 48, 76, 82	7 (100%)
3	A	470/476 (98%)	0.48	38 (8%) 12 14	34, 49, 70, 80	0
3	D	470/476 (98%)	0.51	37 (7%) 12 15	33, 49, 70, 80	0
All	All	970/982 (98%)	0.53	86 (8%) 9 11	33, 49, 72, 87	30 (3%)

All (86) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	16	MET	9.8
3	D	469	ASP	8.4
3	A	469	ASP	8.0
3	A	369	LYS	6.4
3	D	369	LYS	5.6
3	A	468	GLY	5.3
3	A	16	MET	5.1
3	A	394	TYR	4.8
3	D	468	GLY	4.8
3	D	13	VAL	4.5
3	D	12	ARG	4.5
2	F	916	G	4.5
3	A	470	ALA	4.3
3	A	49	GLY	4.2
3	D	49	GLY	4.2
2	C	917	G	4.0
3	A	287	ASN	3.9
2	C	916	G	3.9
3	D	5	ASP	3.9

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Mol	Chain	Res	Type	RSRZ
2	F	917	G	3.6
3	D	370	SER	3.6
3	A	14	HIS	3.6
3	D	14	HIS	3.5
3	A	12	ARG	3.5
3	D	17	ARG	3.5
3	D	287	ASN	3.4
2	C	915	G	3.4
3	A	5	ASP	3.3
3	A	41	ASN	3.3
2	F	915	G	3.2
3	D	15	VAL	3.2
1	B	910	C	3.1
3	D	154	GLU	3.1
3	A	8	ASP	3.1
3	A	17	ARG	3.1
3	D	7	ARG	3.0
3	D	9	VAL	3.0
3	A	7	ARG	3.0
3	D	48	GLU	3.0
3	D	470	ALA	2.9
1	E	903	A	2.9
3	A	13	VAL	2.8
3	A	371	ASP	2.8
3	A	48	GLU	2.8
3	A	52	LEU	2.8
3	D	10	GLU	2.8
1	E	909	C	2.8
3	A	303	THR	2.7
3	D	41	ASN	2.7
3	D	371	ASP	2.7
3	D	394	TYR	2.6
3	D	114	ASP	2.6
3	D	52	LEU	2.6
3	A	9	VAL	2.6
3	A	6	THR	2.5
3	D	127	ARG	2.5
3	D	28	GLY	2.5
3	A	370	SER	2.5
3	D	6	THR	2.4
3	A	375	VAL	2.4
3	A	417	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
3	A	122	ALA	2.3
1	E	910	C	2.3
3	A	114	ASP	2.3
3	D	132	ILE	2.3
3	D	8	ASP	2.3
3	A	127	ARG	2.3
3	A	297	PRO	2.2
3	A	15	VAL	2.2
3	D	11	GLU	2.2
3	A	33	GLU	2.2
3	D	235	ASN	2.2
3	D	126	LYS	2.2
1	B	909	C	2.2
3	D	54	GLU	2.2
3	A	154	GLU	2.1
3	D	298	SER	2.1
3	D	299	GLY	2.1
3	A	126	LYS	2.1
3	A	298	SER	2.1
3	D	284	ALA	2.1
3	A	106	ASP	2.1
3	A	307	ASN	2.1
3	D	50	VAL	2.0
3	A	132	ILE	2.0
3	A	125	GLY	2.0

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 carbohydrates 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(\AA^2)	Q<0.9
5	PPV	D	2930	9/9	0.53	0.51	101,102,102,102	9
5	PPV	A	930	9/9	0.58	0.48	129,129,129,129	9
4	MG	D	3990	1/1	0.90	0.27	47,47,47,47	0
4	MG	A	1990	1/1	0.92	0.17	44,44,44,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.