Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 06:50 AM GMT

PDB ID	:	2YI8
Title	:	Structure of the RNA polymerase VP1 from Infectious Pancreatic Necrosis Virus
Authors	:	Graham, S.C.; Sarin, L.P.; Bahar, M.W.; Myers, R.A.; Stuart, D.I.; Bamford, D.H.; Grimes, J.M.
Deposited on	:	2011-05-11
Resolution	:	2.30 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

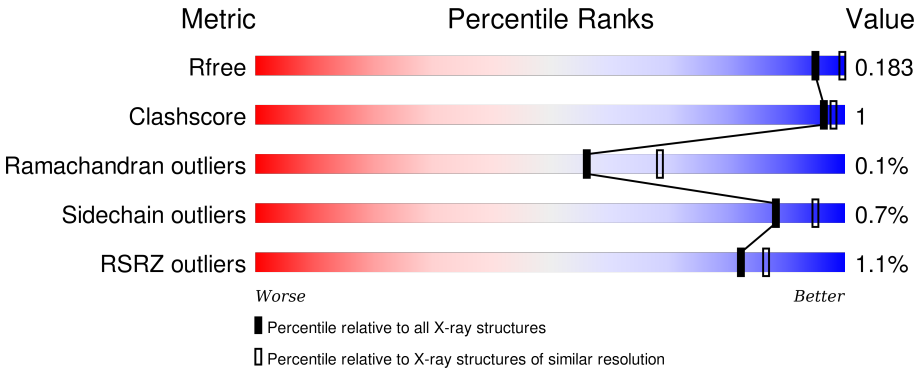
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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. 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	799	3% 92% 5% ·
1	B	799	% 92% 5% ·
1	C	799	% 92% · ·
1	D	799	% 92% 5% ·
1	E	799	92% 5% ·

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 33471 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 RNA-DIRECTED RNA POLYMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	771	Total	C	N	O	S	0	6	0
			6054	3845	1021	1161	27			
1	B	771	Total	C	N	O	S	0	7	0
			6054	3843	1021	1163	27			
1	C	771	Total	C	N	O	S	0	7	0
			6061	3849	1022	1163	27			
1	D	771	Total	C	N	O	S	0	8	0
			6063	3850	1022	1164	27			
1	E	771	Total	C	N	O	S	0	9	0
			6085	3865	1024	1169	27			

There are 45 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	791	LYS	-	EXPRESSION TAG	UNP P22173
A	792	THR	-	EXPRESSION TAG	UNP P22173
A	793	GLY	-	EXPRESSION TAG	UNP P22173
A	794	HIS	-	EXPRESSION TAG	UNP P22173
A	795	HIS	-	EXPRESSION TAG	UNP P22173
A	796	HIS	-	EXPRESSION TAG	UNP P22173
A	797	HIS	-	EXPRESSION TAG	UNP P22173
A	798	HIS	-	EXPRESSION TAG	UNP P22173
A	799	HIS	-	EXPRESSION TAG	UNP P22173
B	791	LYS	-	EXPRESSION TAG	UNP P22173
B	792	THR	-	EXPRESSION TAG	UNP P22173
B	793	GLY	-	EXPRESSION TAG	UNP P22173
B	794	HIS	-	EXPRESSION TAG	UNP P22173
B	795	HIS	-	EXPRESSION TAG	UNP P22173
B	796	HIS	-	EXPRESSION TAG	UNP P22173
B	797	HIS	-	EXPRESSION TAG	UNP P22173
B	798	HIS	-	EXPRESSION TAG	UNP P22173
B	799	HIS	-	EXPRESSION TAG	UNP P22173
C	791	LYS	-	EXPRESSION TAG	UNP P22173

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Chain	Residue	Modelled	Actual	Comment	Reference
C	792	THR	-	EXPRESSION TAG	UNP P22173
C	793	GLY	-	EXPRESSION TAG	UNP P22173
C	794	HIS	-	EXPRESSION TAG	UNP P22173
C	795	HIS	-	EXPRESSION TAG	UNP P22173
C	796	HIS	-	EXPRESSION TAG	UNP P22173
C	797	HIS	-	EXPRESSION TAG	UNP P22173
C	798	HIS	-	EXPRESSION TAG	UNP P22173
C	799	HIS	-	EXPRESSION TAG	UNP P22173
D	791	LYS	-	EXPRESSION TAG	UNP P22173
D	792	THR	-	EXPRESSION TAG	UNP P22173
D	793	GLY	-	EXPRESSION TAG	UNP P22173
D	794	HIS	-	EXPRESSION TAG	UNP P22173
D	795	HIS	-	EXPRESSION TAG	UNP P22173
D	796	HIS	-	EXPRESSION TAG	UNP P22173
D	797	HIS	-	EXPRESSION TAG	UNP P22173
D	798	HIS	-	EXPRESSION TAG	UNP P22173
D	799	HIS	-	EXPRESSION TAG	UNP P22173
E	791	LYS	-	EXPRESSION TAG	UNP P22173
E	792	THR	-	EXPRESSION TAG	UNP P22173
E	793	GLY	-	EXPRESSION TAG	UNP P22173
E	794	HIS	-	EXPRESSION TAG	UNP P22173
E	795	HIS	-	EXPRESSION TAG	UNP P22173
E	796	HIS	-	EXPRESSION TAG	UNP P22173
E	797	HIS	-	EXPRESSION TAG	UNP P22173
E	798	HIS	-	EXPRESSION TAG	UNP P22173
E	799	HIS	-	EXPRESSION TAG	UNP P22173

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total K 1 1	0	0
2	A	1	Total K 1 1	0	0
2	D	1	Total K 1 1	0	0
2	C	1	Total K 1 1	0	0
2	E	1	Total K 1 1	0	0

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Cl 1 1	0	0
3	A	1	Total Cl 1 1	0	0
3	D	1	Total Cl 1 1	0	0
3	C	1	Total Cl 1 1	0	0
3	E	2	Total Cl 2 2	0	0

- Molecule 4 is water.

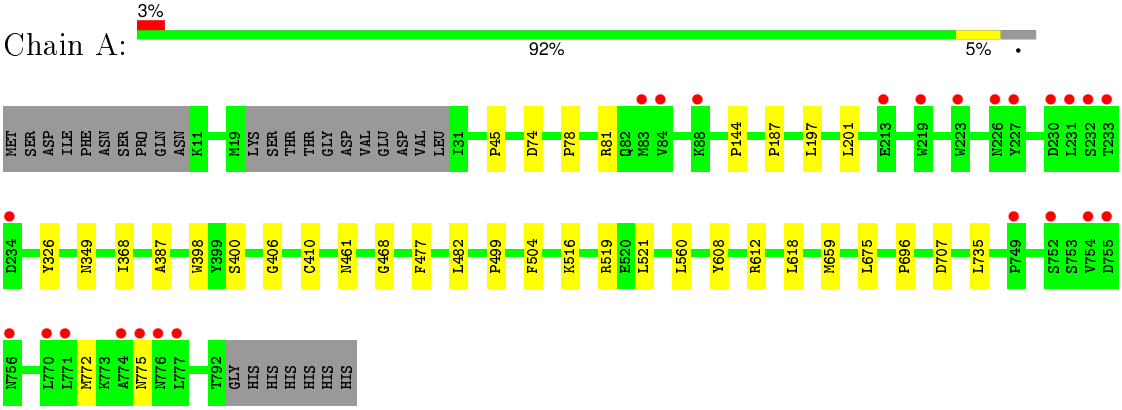
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	563	Total O 563 563	0	0
4	B	622	Total O 622 622	0	0
4	C	611	Total O 611 611	0	0
4	D	637	Total O 637 637	0	0
4	E	710	Total O 710 710	0	0

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 errors 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.

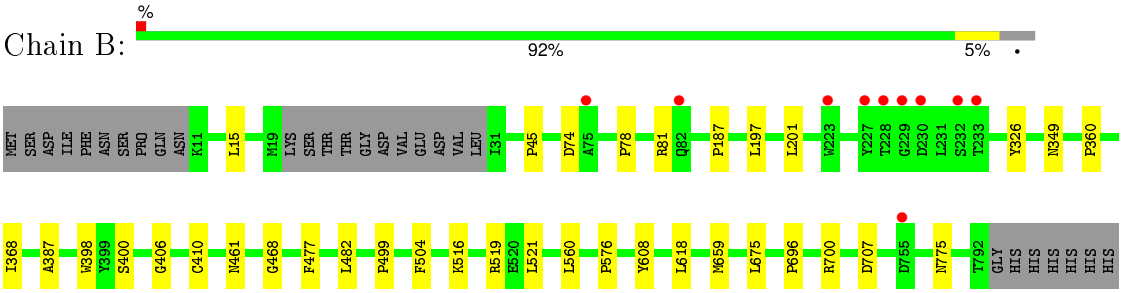
- Molecule 1: RNA-DIRECTED RNA POLYMERASE

Chain A:



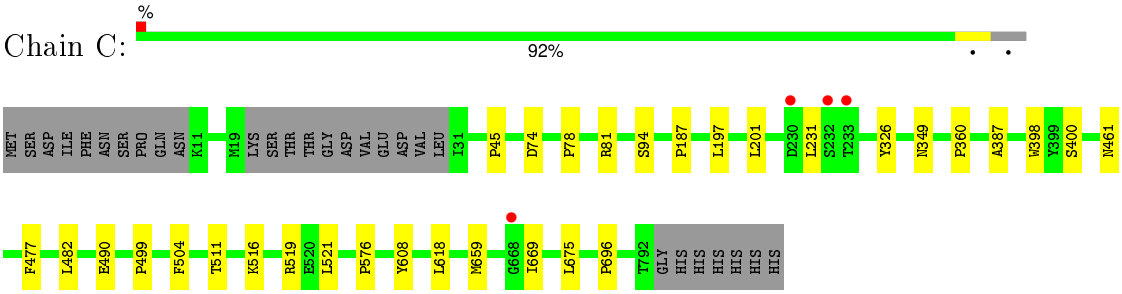
- Molecule 1: RNA-DIRECTED RNA POLYMERASE

Chain B:



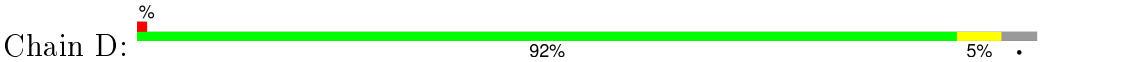
- Molecule 1: RNA-DIRECTED RNA POLYMERASE

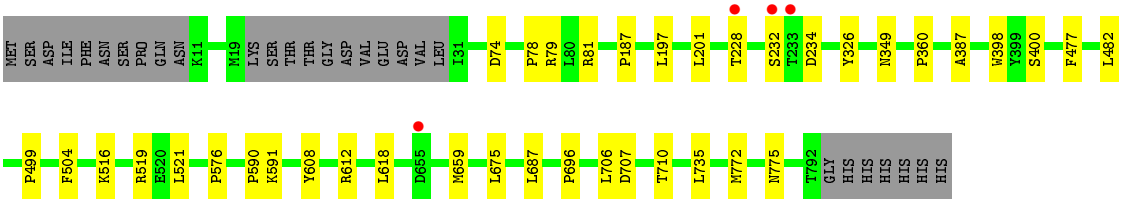
Chain C:



- Molecule 1: RNA-DIRECTED RNA POLYMERASE

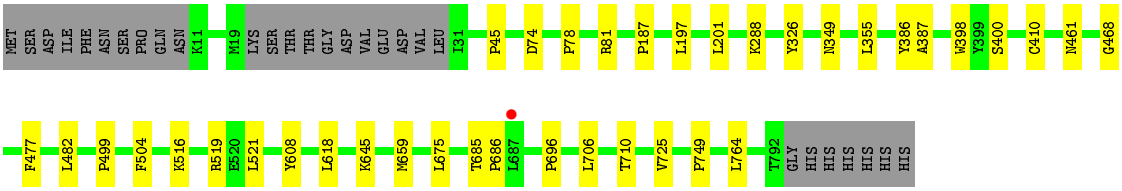
Chain D:





- Molecule 1: RNA-DIRECTED RNA POLYMERASE

Chain E:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	134.05Å 183.87Å 244.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	90.33 – 2.30 81.08 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (90.33-2.30) 100.0 (81.08-2.30)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.86 (at 2.29Å)	Xtriage
Refinement program	BUSTER 2.9.2	Depositor
R, R_{free}	0.166 , 0.187 0.163 , 0.183	Depositor DCC
R_{free} test set	2655 reflections (1.01%)	DCC
Wilson B-factor (Å^2)	12.0	Xtriage
Anisotropy	0.473	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 46.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 266799 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	33471	wwPDB-VP
Average B, all atoms (Å^2)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.18% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: K, CL

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	0.52	0/6207	0.61	0/8448
1	B	0.54	0/6210	0.61	0/8453
1	C	0.54	0/6217	0.62	0/8461
1	D	0.54	0/6222	0.62	0/8468
1	E	0.57	0/6247	0.63	0/8498
All	All	0.54	0/31103	0.62	0/42328

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	6054	0	6023	18	0
1	B	6054	0	6014	19	0
1	C	6061	0	6034	15	0
1	D	6063	0	6034	17	0
1	E	6085	0	6078	20	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	1	0	0	0	0
2	E	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	2	0	0	0	0
4	A	563	0	0	2	0
4	B	622	0	0	2	0
4	C	611	0	0	0	0
4	D	637	0	0	0	0
4	E	710	0	0	2	0
All	All	33471	0	30183	87	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 1.

All (87) 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:E:288:LYS:HE2	4:E:2350:HOH:O	1.69	0.91
1:B:187:PRO:HB3	1:B:197:LEU:HD11	1.78	0.66
1:C:187:PRO:HB3	1:C:197:LEU:HD11	1.78	0.64
1:E:706:LEU:O	1:E:710[B]:THR:HG23	1.97	0.64
1:D:187:PRO:HB3	1:D:197:LEU:HD11	1.78	0.64
1:E:187:PRO:HB3	1:E:197:LEU:HD11	1.79	0.64
1:A:187:PRO:HB3	1:A:197:LEU:HD11	1.80	0.62
1:A:735:LEU:HD21	1:A:772:MET:HG2	1.85	0.58
1:D:74:ASP:HB2	1:D:81:ARG:HG2	1.89	0.55
1:C:74:ASP:HB2	1:C:81:ARG:HG2	1.88	0.55
1:C:201:LEU:HD11	1:C:482[A]:LEU:HD13	1.89	0.54
1:A:201:LEU:HD11	1:A:482[A]:LEU:HD13	1.90	0.54
1:D:608:TYR:HB2	1:D:659:MET:HE2	1.88	0.54
1:E:201:LEU:HD11	1:E:482[A]:LEU:HD13	1.90	0.54
1:B:74:ASP:HB2	1:B:81:ARG:HG2	1.89	0.54
1:D:201:LEU:HD11	1:D:482[A]:LEU:HD13	1.90	0.54
1:D:590:PRO:HB2	1:D:612:ARG:HD2	1.88	0.54
1:E:618:LEU:HG	1:E:675:LEU:HD12	1.91	0.53
1:E:74:ASP:HB2	1:E:81:ARG:HG2	1.89	0.53
1:D:618:LEU:HG	1:D:675:LEU:HD12	1.91	0.53
1:B:700:ARG:NH2	4:B:2412:HOH:O	2.37	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:ASP:HB2	1:A:81:ARG:HG2	1.90	0.53
1:B:618:LEU:HG	1:B:675:LEU:HD12	1.90	0.53
1:B:201:LEU:HD11	1:B:482[A]:LEU:HD13	1.90	0.52
1:C:618:LEU:HG	1:C:675:LEU:HD12	1.92	0.52
1:A:618:LEU:HG	1:A:675:LEU:HD12	1.91	0.51
1:A:608:TYR:HB2	1:A:659:MET:HE3	1.92	0.51
1:E:608:TYR:HB2	1:E:659:MET:HE3	1.92	0.51
1:C:608:TYR:HB2	1:C:659:MET:HE2	1.92	0.50
1:D:706:LEU:O	1:D:710[B]:THR:HG23	2.12	0.49
1:D:349:ASN:O	1:D:696:PRO:HD2	2.13	0.48
1:D:400:SER:HB3	1:D:519:ARG:HB3	1.96	0.48
1:C:398:TRP:HB3	1:C:521:LEU:HB3	1.95	0.48
1:E:349:ASN:O	1:E:696:PRO:HD2	2.13	0.48
1:A:400:SER:HB3	1:A:519:ARG:HB3	1.96	0.48
1:D:78:PRO:O	1:D:81:ARG:HG3	2.14	0.47
1:D:398:TRP:HB3	1:D:521:LEU:HB3	1.97	0.47
1:E:398:TRP:HB3	1:E:521:LEU:HB3	1.96	0.47
1:B:398:TRP:HB3	1:B:521:LEU:HB3	1.96	0.47
1:A:349:ASN:O	1:A:696:PRO:HD2	2.15	0.47
1:A:387:ALA:HB2	1:A:477:PHE:CD2	2.50	0.47
1:A:612:ARG:HG2	4:A:2473:HOH:O	2.15	0.47
1:C:387:ALA:HB2	1:C:477:PHE:CD2	2.50	0.47
1:E:749:PRO:HG2	1:E:764:LEU:CD1	2.45	0.47
1:A:398:TRP:HB3	1:A:521:LEU:HB3	1.96	0.47
1:D:591:LYS:O	1:D:608:TYR:OH	2.25	0.46
1:A:78:PRO:O	1:A:81:ARG:HG3	2.16	0.46
1:B:400:SER:HB3	1:B:519:ARG:HB3	1.97	0.46
1:B:387:ALA:HB2	1:B:477:PHE:CD2	2.51	0.46
1:E:400:SER:HB3	1:E:519:ARG:HB3	1.96	0.46
1:B:775:ASN:ND2	4:B:2611:HOH:O	2.49	0.46
1:B:499:PRO:HA	1:B:504:PHE:CG	2.51	0.46
1:E:387:ALA:HB2	1:E:477:PHE:CD2	2.50	0.46
1:E:78:PRO:O	1:E:81:ARG:HG3	2.16	0.45
1:B:349:ASN:O	1:B:696:PRO:HD2	2.16	0.45
1:C:499:PRO:HA	1:C:504:PHE:CG	2.51	0.45
1:D:387:ALA:HB2	1:D:477:PHE:CD2	2.51	0.45
1:A:612:ARG:CG	4:A:2473:HOH:O	2.64	0.45
1:C:78:PRO:O	1:C:81:ARG:HG3	2.17	0.45
1:B:78:PRO:O	1:B:81:ARG:HG3	2.16	0.45
1:C:349:ASN:O	1:C:696:PRO:HD2	2.15	0.45
1:B:608:TYR:HB2	1:B:659:MET:HE3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:499:PRO:HA	1:E:504:PHE:CG	2.52	0.44
1:A:499:PRO:HA	1:A:504:PHE:CG	2.51	0.44
1:D:499:PRO:HA	1:D:504:PHE:CG	2.53	0.44
1:D:735:LEU:HD21	1:D:772:MET:HG2	1.99	0.43
1:E:725:VAL:HG23	4:E:2654:HOH:O	2.17	0.43
1:A:144:PRO:HG3	1:E:645:LYS:HA	2.00	0.42
1:C:400:SER:HB3	1:C:519:ARG:HB3	1.99	0.42
1:B:187:PRO:CB	1:B:197:LEU:HD11	2.46	0.42
1:D:360:PRO:HB2	1:D:576:PRO:HB3	2.01	0.42
1:C:490:GLU:HG3	1:C:511:THR:HG22	2.01	0.41
1:A:45:PRO:HB3	1:A:461:ASN:OD1	2.21	0.41
1:C:45:PRO:HB3	1:C:461:ASN:OD1	2.21	0.41
1:B:15:LEU:HD21	1:D:234:ASP:HB3	2.03	0.41
1:E:685:THR:HA	1:E:686:PRO:HD3	2.00	0.41
1:B:360:PRO:HB2	1:B:576:PRO:HB3	2.03	0.41
1:B:368:ILE:HG21	1:B:560:LEU:HD22	2.03	0.41
1:C:187:PRO:CB	1:C:197:LEU:HD11	2.47	0.40
1:A:410:CYS:O	1:A:468:GLY:HA2	2.21	0.40
1:C:360:PRO:HB2	1:C:576:PRO:HB3	2.03	0.40
1:E:45:PRO:HB3	1:E:461:ASN:OD1	2.21	0.40
1:E:355:LEU:HB2	1:E:386:TYR:HB2	2.04	0.40
1:B:410:CYS:O	1:B:468:GLY:HA2	2.22	0.40
1:A:368:ILE:HG21	1:A:560:LEU:HD22	2.04	0.40
1:E:410:CYS:O	1:E:468:GLY:HA2	2.22	0.40
1:B:45:PRO:HB3	1:B:461:ASN:OD1	2.22	0.40

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	773/799 (97%)	756 (98%)	16 (2%)	1 (0%)	56 68

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	774/799 (97%)	758 (98%)	15 (2%)	1 (0%)	56 68
1	C	774/799 (97%)	758 (98%)	16 (2%)	0	100 100
1	D	775/799 (97%)	757 (98%)	18 (2%)	0	100 100
1	E	776/799 (97%)	761 (98%)	15 (2%)	0	100 100
All	All	3872/3995 (97%)	3790 (98%)	80 (2%)	2 (0%)	56 68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	406	GLY
1	B	406	GLY

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	660/692 (95%)	656 (99%)	4 (1%)	90 96
1	B	660/692 (95%)	657 (100%)	3 (0%)	92 97
1	C	662/692 (96%)	657 (99%)	5 (1%)	86 94
1	D	662/692 (96%)	654 (99%)	8 (1%)	78 89
1	E	668/692 (96%)	666 (100%)	2 (0%)	94 98
All	All	3312/3460 (96%)	3290 (99%)	22 (1%)	88 95

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	326	TYR
1	A	516	LYS
1	A	707	ASP
1	A	775	ASN
1	B	326	TYR
1	B	516	LYS

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Mol	Chain	Res	Type
1	B	707	ASP
1	C	94	SER
1	C	231	LEU
1	C	326	TYR
1	C	516	LYS
1	C	669	ILE
1	D	79	ARG
1	D	228	THR
1	D	232	SER
1	D	326	TYR
1	D	516	LYS
1	D	687	LEU
1	D	707	ASP
1	D	775	ASN
1	E	326	TYR
1	E	516	LYS

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

Mol	Chain	Res	Type
1	B	775	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry

Of 11 ligands modelled in this entry, 11 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

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, 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	771/799 (96%)	-0.10	24 (3%) 52 62	5, 17, 46, 72	0
1	B	771/799 (96%)	-0.21	10 (1%) 79 84	3, 14, 41, 63	0
1	C	771/799 (96%)	-0.27	4 (0%) 91 94	4, 14, 39, 58	0
1	D	771/799 (96%)	-0.28	4 (0%) 91 94	3, 13, 39, 70	0
1	E	771/799 (96%)	-0.30	1 (0%) 95 97	3, 10, 33, 56	0
All	All	3855/3995 (96%)	-0.23	43 (1%) 82 86	3, 14, 40, 72	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	228	THR	5.4
1	A	774	ALA	4.5
1	A	776	ASN	4.3
1	A	775	ASN	4.3
1	B	232	SER	4.3
1	A	88	LYS	4.0
1	A	223	TRP	3.8
1	E	687	LEU	3.8
1	B	228	THR	3.7
1	C	232	SER	3.7
1	A	232	SER	3.7
1	D	232	SER	3.6
1	C	233	THR	3.3
1	A	233	THR	3.2
1	A	756	ASN	3.1
1	A	755	ASP	3.0
1	A	777	LEU	2.9
1	B	233	THR	2.9
1	D	233	THR	2.9
1	B	223	TRP	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	219	TRP	2.8
1	C	230	ASP	2.8
1	B	229	GLY	2.7
1	A	771	LEU	2.7
1	A	84	VAL	2.7
1	C	668	GLY	2.6
1	A	231	LEU	2.5
1	A	230	ASP	2.4
1	B	82	GLN	2.4
1	B	755	ASP	2.4
1	A	754	VAL	2.3
1	A	749	PRO	2.3
1	A	752	SER	2.2
1	A	226	ASN	2.2
1	B	230	ASP	2.2
1	B	75	ALA	2.2
1	A	227	TYR	2.1
1	A	770	LEU	2.1
1	A	213	GLU	2.1
1	A	234	ASP	2.1
1	B	227	TYR	2.1
1	D	655	ASP	2.0
1	A	83	MET	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å^2)	Q<0.9
2	K	B	1002	1/1	1.00	0.09	-2.02	12,12,12,12	0
2	K	E	1002	1/1	1.00	0.09	-2.70	10,10,10,10	0
2	K	D	1002	1/1	1.00	0.07	-3.27	10,10,10,10	0
3	CL	E	1005	1/1	0.99	0.08	-3.32	14,14,14,14	0
2	K	A	1002	1/1	0.99	0.07	-3.59	14,14,14,14	0
2	K	C	1002	1/1	0.99	0.09	-4.59	12,12,12,12	0
3	CL	A	1003	1/1	1.00	0.08	-	21,21,21,21	0
3	CL	E	1003	1/1	0.99	0.12	-	14,14,14,14	0
3	CL	B	1003	1/1	1.00	0.09	-	14,14,14,14	0
3	CL	D	1003	1/1	0.99	0.09	-	16,16,16,16	0
3	CL	C	1003	1/1	0.99	0.11	-	17,17,17,17	0

6.5 Other polymers (i)

There are no such residues in this entry.