W O R L D W I D E
wwPDB
PROTEIN DATA BANK

Full wwPDB X-ray Structure Validation Report (i)

Feb 19, 2016 – 09:00 PM GMT

PDB ID	:	4Z3U
Title	:	PRV nuclear egress complex
Authors	:	Bigalke, J.M.; Heldwein, E.E.
Deposited on	:	2015-03-31
Resolution	:	2.71 Å(reported)

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

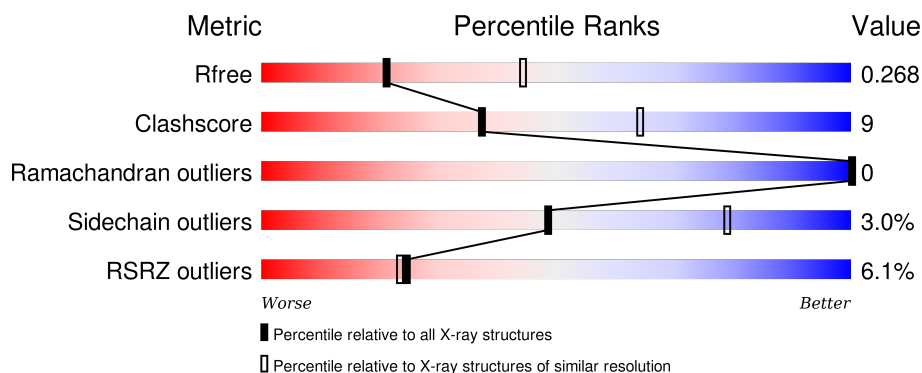|		
MolProbity	:	4.02b-467
Mogul	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026982
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20026982

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	Percentile Ranks	Value
Rfree		0.268
Clashscore		9
Ramachandran outliers		0
Sidechain outliers		3.0%
RSRZ outliers		6.1%

Worse ⬛ *Better*

⬛ Percentile relative to all X-ray structures
▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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 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	181	3% — 73% / 19% / • 7%
1	C	181	6% — 66% / 26% / • 6%
2	B	258	6% — 80% / 16% / • •
2	D	258	7% — 85% / 10% / • •

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
3	CL	A	201	-	-	-	X
3	CL	C	202	-	-	X	-
4	ZN	B	301	-	-	-	X
5	NA	C	201	-	-	-	X

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 6611 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 UL34 protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	168	Total	C	N	O	S	Se	0	0	0
			1310	817	244	238	5	6			
1	C	170	Total	C	N	O	S	Se	0	0	0
			1324	824	249	240	5	6			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP G3G8R3
A	-3	PRO	-	expression tag	UNP G3G8R3
A	-2	LEU	-	expression tag	UNP G3G8R3
A	-1	GLY	-	expression tag	UNP G3G8R3
A	0	SER	-	expression tag	UNP G3G8R3
C	-4	GLY	-	expression tag	UNP G3G8R3
C	-3	PRO	-	expression tag	UNP G3G8R3
C	-2	LEU	-	expression tag	UNP G3G8R3
C	-1	GLY	-	expression tag	UNP G3G8R3
C	0	SER	-	expression tag	UNP G3G8R3

- Molecule 2 is a protein called UL31.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	252	Total	C	N	O	S	Se	0	1	0
			1981	1266	347	357	7	4			
2	D	249	Total	C	N	O	S	Se	0	0	0
			1950	1250	341	348	7	4			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	14	GLY	-	expression tag	UNP G3G955
B	15	PRO	-	expression tag	UNP G3G955

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Chain	Residue	Modelled	Actual	Comment	Reference
B	16	GLY	-	expression tag	UNP G3G955
B	17	SER	-	expression tag	UNP G3G955
D	14	GLY	-	expression tag	UNP G3G955
D	15	PRO	-	expression tag	UNP G3G955
D	16	GLY	-	expression tag	UNP G3G955
D	17	SER	-	expression tag	UNP G3G955

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	2	Total Cl 2 2	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

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Zn 1 1	0	0
4	D	1	Total Zn 1 1	0	0

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	1	Total Na 1 1	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	9	Total O 9 9	0	0
6	B	9	Total O 9 9	0	0
6	C	11	Total O 11 11	0	0

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| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	9	Total 9	O 9	0	0

3 Residue-property plots ⓘ

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.
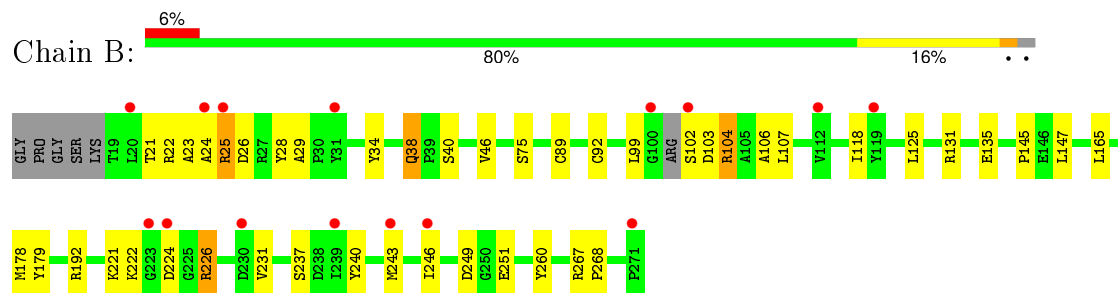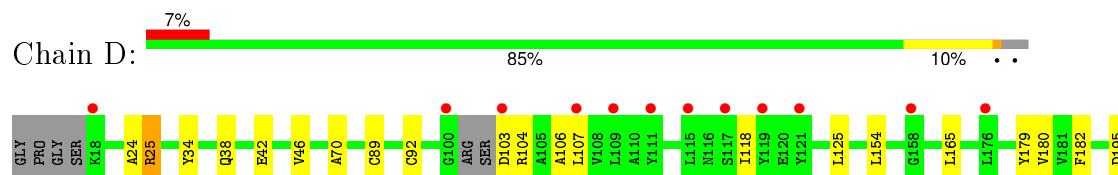
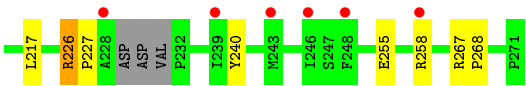- Molecule 1: UL34 protein
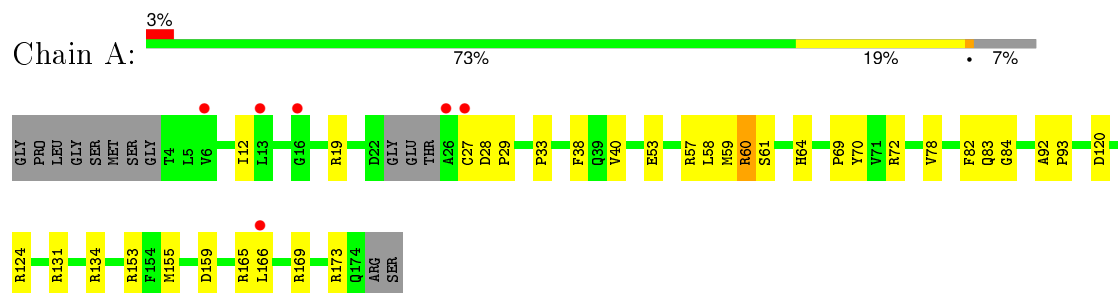
- Molecule 1: UL34 protein
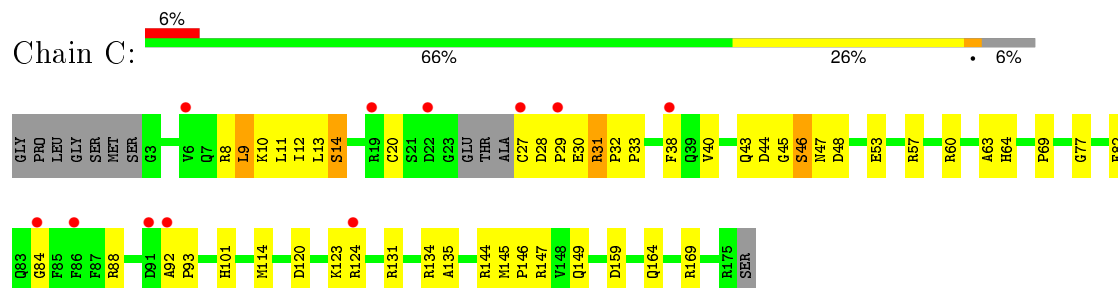
- Molecule 2: UL31

- Molecule 2: UL31

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α, β, γ	125.46Å 125.46Å 109.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.91 – 2.71 49.91 – 2.71	Depositor EDS
% Data completeness (in resolution range)	99.3 (49.91-2.71) 99.6 (49.91-2.71)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$<I/\sigma(I)>$ [1]	1.29 (at 2.69Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.212 , 0.273 0.213 , 0.268	Depositor DCC
R_{free} test set	1995 reflections (8.24%)	DCC
Wilson B-factor (Å^2)	65.2	Xtriage
Anisotropy	0.392	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 44.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$<|L|> = 0.47, <L^2> = 0.30$	Xtriage
Outliers	6 of 24293 reflections (0.025%)	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	6611	wwPDB-VP
Average B, all atoms (Å^2)	81.0	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 61.21 % 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 1.3565e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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[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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: NA, ZN, 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	0.44	0/1330	0.56	0/1791
1	C	0.44	0/1344	0.58	1/1808 (0.1%)
2	B	0.42	0/2025	0.53	0/2744
2	D	0.32	0/1993	0.46	0/2697
All	All	0.40	0/6692	0.53	1/9040 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	47	ASN	N-CA-C	-5.43	96.33	111.00

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	1310	0	1301	28	1
1	C	1324	0	1317	41	1
2	B	1981	0	1955	43	1
2	D	1950	0	1934	19	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	1	0	0	2	0
3	D	1	0	0	0	0
4	B	1	0	0	0	0
4	D	1	0	0	0	0
5	C	1	0	0	0	0
6	A	9	0	0	2	0
6	B	9	0	0	1	0
6	C	11	0	0	0	0
6	D	9	0	0	0	0
All	All	6611	0	6507	122	2

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 9.

All (122) 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)
2:D:103:ASP:HB3	2:D:106:ALA:HB3	1.29	1.09
1:C:144:ARG:NH1	2:D:195:ASP:OD2	1.88	1.06
1:C:31:ARG:NH1	1:C:32:PRO:HD2	1.81	0.96
1:C:11:LEU:O	1:C:14:SER:OG	1.84	0.95
2:B:104:ARG:HH11	2:B:104:ARG:HG3	1.37	0.89
2:B:22:ARG:HG2	2:B:26:ASP:CB	2.06	0.86
2:B:22:ARG:HG2	2:B:26:ASP:HB3	1.57	0.85
2:D:103:ASP:HB3	2:D:106:ALA:CB	2.06	0.85
2:B:25:ARG:HA	2:B:28:TYR:HB2	1.63	0.81
1:C:53:GLU:OE2	1:C:57:ARG:NH1	2.14	0.81
2:B:25:ARG:O	2:B:29:ALA:HB2	1.81	0.79
1:A:166:LEU:HD23	1:A:166:LEU:O	1.84	0.77
2:B:135:GLU:OE2	2:B:237:SER:OG	2.04	0.75
2:B:22:ARG:HG2	2:B:26:ASP:CG	2.07	0.75
1:C:63:ALA:O	2:D:25:ARG:NH2	2.19	0.75
1:C:88:ARG:NH1	1:C:92:ALA:O	2.19	0.74
2:B:25:ARG:O	2:B:29:ALA:N	2.19	0.74
1:C:43:GLN:O	1:C:46:SER:HB3	1.88	0.74
2:B:22:ARG:CG	2:B:26:ASP:CG	2.57	0.72
2:B:99:LEU:HB2	2:B:192:ARG:HH11	1.56	0.70
1:C:60:ARG:HH22	2:D:24:ALA:HB3	1.56	0.70
1:C:120:ASP:HB3	1:C:124:ARG:HH12	1.56	0.70
1:C:27:CYS:HB2	1:C:33:PRO:CD	2.22	0.70
2:B:21:THR:O	2:B:22:ARG:C	2.28	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:120:ASP:HB3	1:C:124:ARG:NH1	2.08	0.68
2:B:145:PRO:HG2	2:B:231:VAL:HG11	1.74	0.68
1:A:166:LEU:C	1:A:166:LEU:HD23	2.16	0.66
1:C:77:GLY:O	1:C:145:MSE:HE2	1.97	0.65
2:B:89:CYS:HB3	2:B:92:CYS:HB2	1.78	0.64
1:C:10:LYS:HG2	1:C:20:CYS:HB3	1.80	0.64
2:B:25:ARG:O	2:B:29:ALA:CB	2.45	0.63
1:C:101:HIS:CE1	3:C:202:CL:CL	2.89	0.63
1:C:12:ILE:HG23	2:D:46:VAL:HG11	1.81	0.62
2:B:22:ARG:HG3	2:B:26:ASP:OD2	2.00	0.61
2:B:267:ARG:NH1	2:B:268:PRO:O	2.34	0.61
1:C:145:MSE:HE3	1:C:146:PRO:HA	1.82	0.60
1:A:12:ILE:HG23	2:B:46:VAL:HG11	1.84	0.60
1:A:153:ARG:HB2	1:A:155:MSE:HE2	1.84	0.59
1:C:40:VAL:HB	1:C:82:PHE:HB2	1.85	0.59
1:C:29:PRO:HB3	1:C:64:HIS:ND1	2.18	0.58
2:B:104:ARG:HH11	2:B:104:ARG:CG	2.10	0.58
1:A:57:ARG:HH11	1:A:60:ARG:HH12	1.50	0.58
1:A:59:MSE:HE3	1:A:69:PRO:HB3	1.86	0.58
1:C:27:CYS:HB2	1:C:33:PRO:HD2	1.85	0.57
2:B:25:ARG:HD2	2:B:29:ALA:HB2	1.86	0.57
1:C:134:ARG:NH2	1:C:164:GLN:OE1	2.38	0.56
1:A:165:ARG:NE	6:A:301:HOH:O	2.23	0.56
1:A:120:ASP:HB3	1:A:124:ARG:NH1	2.21	0.56
1:A:120:ASP:HB3	1:A:124:ARG:HH12	1.72	0.55
2:D:118:ILE:HD12	2:D:125:LEU:HD11	1.88	0.55
2:B:103:ASP:O	2:B:106:ALA:HB3	2.07	0.55
1:C:8:ARG:NH2	2:D:42:GLU:OE2	2.29	0.54
1:A:40:VAL:HB	1:A:82:PHE:HB2	1.89	0.54
2:B:40:SER:OG	6:B:402:HOH:O	2.19	0.54
2:B:249:ASP:OD1	2:B:251:GLU:HB2	2.09	0.52
1:C:29:PRO:CB	1:C:64:HIS:ND1	2.73	0.52
2:B:22:ARG:CG	2:B:26:ASP:OD2	2.58	0.52
2:B:24:ALA:O	2:B:25:ARG:HB3	2.09	0.52
2:B:104:ARG:HG3	2:B:107:LEU:HD12	1.92	0.52
1:A:131:ARG:HB2	1:A:134:ARG:HG3	1.92	0.52
1:C:131:ARG:HH21	1:C:134:ARG:CZ	2.23	0.52
2:B:25:ARG:O	2:B:25:ARG:HD2	2.10	0.51
1:C:123:LYS:NZ	1:C:135:ALA:O	2.36	0.50
1:C:101:HIS:ND1	3:C:202:CL:CL	2.81	0.50
1:C:9:LEU:O	1:C:13:LEU:HG	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:TYR:CZ	1:A:92:ALA:HB1	2.47	0.49
1:A:57:ARG:NH1	1:A:60:ARG:HH12	2.09	0.49
2:D:255:GLU:HA	2:D:258:ARG:HH11	1.77	0.49
1:A:27:CYS:SG	1:A:33:PRO:HG3	2.53	0.49
2:D:154:LEU:HD13	2:D:180:VAL:HG11	1.94	0.49
2:B:103:ASP:HB2	2:B:106:ALA:HB3	1.94	0.48
2:B:147:LEU:HB3	2:B:178:MSE:SE	2.64	0.48
1:C:92:ALA:HA	1:C:93:PRO:HD3	1.77	0.47
2:B:103:ASP:CB	2:B:106:ALA:HB3	2.44	0.47
2:B:21:THR:O	2:B:23:ALA:N	2.47	0.47
1:A:53:GLU:O	1:A:57:ARG:HG3	2.13	0.47
1:C:38:PHE:CE2	1:C:84:GLY:HA3	2.49	0.47
1:C:60:ARG:NH2	2:D:24:ALA:HB3	2.27	0.46
2:B:118:ILE:HD12	2:B:125:LEU:HD11	1.97	0.46
2:D:165:LEU:HB2	2:D:179:TYR:HB2	1.98	0.46
2:B:165:LEU:HB2	2:B:179:TYR:HB2	1.96	0.46
1:C:159:ASP:OD1	1:C:159:ASP:N	2.48	0.46
2:B:104:ARG:HH11	2:B:107:LEU:HD12	1.80	0.46
2:D:103:ASP:O	2:D:106:ALA:N	2.49	0.46
1:A:159:ASP:N	1:A:159:ASP:OD1	2.49	0.46
1:C:31:ARG:CZ	1:C:32:PRO:HD2	2.45	0.45
1:A:173:ARG:NH2	2:B:38:GLN:OE1	2.45	0.45
1:C:60:ARG:HB2	1:C:69:PRO:HG3	1.99	0.45
2:B:104:ARG:NH1	2:B:104:ARG:CG	2.74	0.45
1:C:44:ASP:OD1	1:C:45:GLY:N	2.50	0.45
1:A:29:PRO:CB	1:A:64:HIS:ND1	2.80	0.44
1:A:92:ALA:HA	1:A:93:PRO:HD3	1.71	0.44
1:A:166:LEU:C	1:A:166:LEU:CD2	2.85	0.44
1:A:60:ARG:HB2	1:A:69:PRO:HG2	1.98	0.44
1:A:58:LEU:O	1:A:61:SER:OG	2.22	0.44
2:B:221:LYS:NZ	2:B:226:ARG:NH1	2.66	0.44
1:C:27:CYS:HB3	1:C:33:PRO:HG2	1.99	0.44
1:A:19:ARG:NH2	6:A:303:HOH:O	2.47	0.43
1:C:114:MSE:HE3	1:C:114:MSE:HA	1.99	0.43
2:B:243:MET:HA	2:B:246:ILE:HD12	2.00	0.43
2:B:104:ARG:O	2:B:107:LEU:HB2	2.19	0.42
1:C:145:MSE:HB3	1:C:145:MSE:HE3	1.83	0.42
1:A:29:PRO:HB3	1:A:64:HIS:ND1	2.35	0.42
1:A:38:PHE:CE2	1:A:84:GLY:HA3	2.54	0.42
1:A:72:ARG:NH2	1:A:83:GLN:OE1	2.53	0.42
2:D:34:TYR:CZ	2:D:38:GLN:HG3	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:149:GLN:OE1	2:D:70:ALA:N	2.53	0.41
2:D:226:ARG:N	2:D:227:PRO:HD3	2.35	0.41
1:C:30:GLU:OE2	1:C:30:GLU:HA	2.19	0.41
1:C:11:LEU:C	1:C:14:SER:HG	2.21	0.41
2:D:267:ARG:HA	2:D:268:PRO:HD2	1.92	0.41
2:B:131:ARG:HD2	2:B:260:TYR:CE1	2.55	0.41
2:B:267:ARG:HA	2:B:268:PRO:HD2	1.92	0.41
1:C:144:ARG:O	1:C:147:ARG:N	2.53	0.41
1:C:27:CYS:SG	1:C:27:CYS:O	2.79	0.41
1:A:60:ARG:HB3	1:A:60:ARG:HE	1.68	0.41
1:A:165:ARG:CG	1:A:165:ARG:HH11	2.34	0.40
2:D:182:PHE:HE2	2:D:217:LEU:HG	1.85	0.40
2:D:89:CYS:HB3	2:D:92:CYS:HB2	2.03	0.40
2:B:24:ALA:O	2:B:25:ARG:CB	2.70	0.40
2:B:34:TYR:CZ	2:B:38:GLN:HG3	2.55	0.40
2:B:25:ARG:HA	2:B:28:TYR:CB	2.42	0.40

All (2) 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)
2:B:222:LYS:NZ	2:B:224:ASP:OD2[8_665]	1.80	0.40
1:A:169:ARG:NH2	1:C:169:ARG:NH2[4_455]	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	164/181 (91%)	162 (99%)	2 (1%)	0	100	100
1	C	166/181 (92%)	164 (99%)	2 (1%)	0	100	100
2	B	249/258 (96%)	243 (98%)	6 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	243/258 (94%)	238 (98%)	5 (2%)	0	100	100
All	All	822/878 (94%)	807 (98%)	15 (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	145/148 (98%)	142 (98%)	3 (2%)	61	87
1	C	146/148 (99%)	140 (96%)	6 (4%)	37	69
2	B	205/204 (100%)	198 (97%)	7 (3%)	44	75
2	D	201/204 (98%)	196 (98%)	5 (2%)	55	84
All	All	697/704 (99%)	676 (97%)	21 (3%)	48	79

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	ASP
1	A	60	ARG
1	A	78	VAL
2	B	25	ARG
2	B	38	GLN
2	B	75	SER
2	B	102	SER
2	B	104	ARG
2	B	226	ARG
2	B	240	TYR
1	C	9	LEU
1	C	14	SER
1	C	28	ASP
1	C	31	ARG
1	C	46	SER
1	C	48	ASP

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Mol	Chain	Res	Type
2	D	25	ARG
2	D	104	ARG
2	D	107	LEU
2	D	226	ARG
2	D	240	TYR

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

Mol	Chain	Res	Type
2	D	188	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 carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 8 ligands modelled in this entry, 8 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	162/181 (89%)	0.13	6 (3%)	45	45	40, 77, 121, 161	0
1	C	164/181 (90%)	0.34	11 (6%)	21	19	51, 79, 121, 191	0
2	B	248/258 (96%)	0.28	15 (6%)	25	24	39, 72, 141, 186	0
2	D	245/258 (94%)	0.23	18 (7%)	18	16	42, 70, 143, 171	0
All	All	819/878 (93%)	0.25	50 (6%)	25	23	39, 75, 140, 191	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	92	ALA	5.9
2	B	25	ARG	5.6
1	C	38	PHE	5.2
2	D	228	ALA	5.0
2	B	24	ALA	5.0
2	B	246	ILE	4.4
2	B	119	TYR	4.3
2	B	100	GLY	4.2
2	D	107	LEU	3.8
1	C	91	ASP	3.7
1	A	26	ALA	3.4
2	D	243	MET	3.3
2	D	119	TYR	3.3
1	C	27	CYS	3.2
2	B	239	ILE	3.2
2	B	112	VAL	3.1
2	B	243	MET	3.1
1	A	6	VAL	3.1
2	D	239	ILE	3.1
2	D	176	LEU	3.0
2	B	224	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	27	CYS	3.0
2	D	18	LYS	2.9
2	B	20	LEU	2.9
1	A	16	GLY	2.9
2	D	117	SER	2.8
1	C	19	ARG	2.7
2	B	102	SER	2.7
2	B	271	PRO	2.6
1	A	166	LEU	2.6
2	D	158	GLY	2.5
2	D	248	PHE	2.5
2	D	115	LEU	2.4
2	B	230	ASP	2.4
2	D	103	ASP	2.4
2	D	246	ILE	2.3
1	C	86	PHE	2.2
2	D	100	GLY	2.2
1	C	6	VAL	2.2
2	B	223	GLY	2.2
1	C	84	GLY	2.1
2	B	31	TYR	2.1
1	A	13	LEU	2.1
2	D	121	TYR	2.1
1	C	29	PRO	2.1
2	D	111	TYR	2.1
2	D	109	LEU	2.1
2	D	258	ARG	2.1
1	C	124	ARG	2.0
1	C	22	ASP	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.

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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, 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	LLDF	B-factors(Å^2)	Q<0.9
5	NA	C	201	1/1	0.90	0.32	3.84	82,82,82,82	0
3	CL	A	201	1/1	0.90	0.30	3.75	106,106,106,106	0
4	ZN	B	301	1/1	0.98	0.27	2.70	67,67,67,67	0
4	ZN	D	301	1/1	0.99	0.18	0.85	52,52,52,52	0
3	CL	B	302	1/1	0.92	0.21	0.37	77,77,77,77	0
3	CL	C	202	1/1	0.85	0.17	0.15	108,108,108,108	0
3	CL	B	303	1/1	0.92	0.20	0.05	88,88,88,88	0
3	CL	D	302	1/1	0.94	0.16	-0.29	79,79,79,79	0

6.5 Other polymers (i)

There are no such residues in this entry.