



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:39 AM GMT

PDB ID : 3MGV
Title : Cre recombinase-DNA transition state
Authors : Gibb, B.P.; Gupta, K.; Ghosh, K.; Sharp, R.; Chen, J.; Van Duyne, G.D.
Deposited on : 2010-04-07
Resolution : 2.29 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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

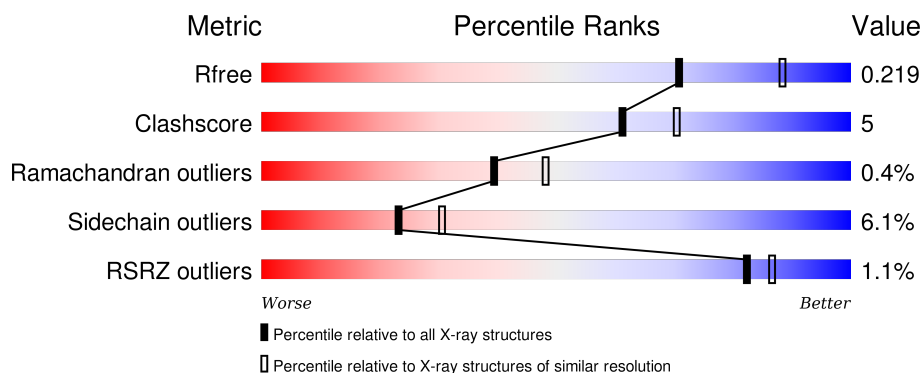
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.29 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 $\leq 5\%$. 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	343	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 7%, orange 0%, yellow 80%, green 81%, grey 81%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 7% 80% 11% 6% </div> </div>
1	B	343	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 7%, orange 0%, yellow 80%, green 81%, grey 81%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 7% 81% 11% 6% </div> </div>
1	C	343	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 7%, orange 0%, yellow 78%, green 78%, grey 78%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 7% 78% 14% 6% </div> </div>
1	D	343	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, orange 0%, yellow 83%, green 83%, grey 83%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 83% 9% 6% </div> </div>
2	E	15	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 7%, orange 0%, yellow 40%, green 40%, grey 40%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 7% 40% 60% </div> </div>

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Mol	Chain	Length	Quality of chain
2	G	15	
2	I	15	
2	L	15	
3	F	20	
3	H	20	
3	K	20	
3	M	20	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 13393 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 Recombinase cre.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	0	0	0
			2550	1584	486	465	15			
1	B	322	Total	C	N	O	S	0	0	0
			2550	1584	486	465	15			
1	C	322	Total	C	N	O	S	0	0	0
			2550	1584	486	465	15			
1	D	322	Total	C	N	O	S	0	0	0
			2550	1584	486	465	15			

- Molecule 2 is a DNA chain called DNA (5'-D(*TP*AP*TP*AP*AP*CP*TP*TP*CP*GP*TP*AP*TP*AP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	15	Total	C	N	O	P	0	0	0
			304	148	53	89	14			
2	G	15	Total	C	N	O	P	0	0	0
			304	148	53	89	14			
2	I	15	Total	C	N	O	P	0	0	0
			304	148	53	89	14			
2	L	15	Total	C	N	O	P	0	0	0
			304	148	53	89	14			

- Molecule 3 is a DNA chain called DNA (5'-D(*CP*AP*TP*AP*TP*GP*CP*TP*AP*TP*AP*CP*GP*AP*AP*GP*TP*TP*AP*T)-3').

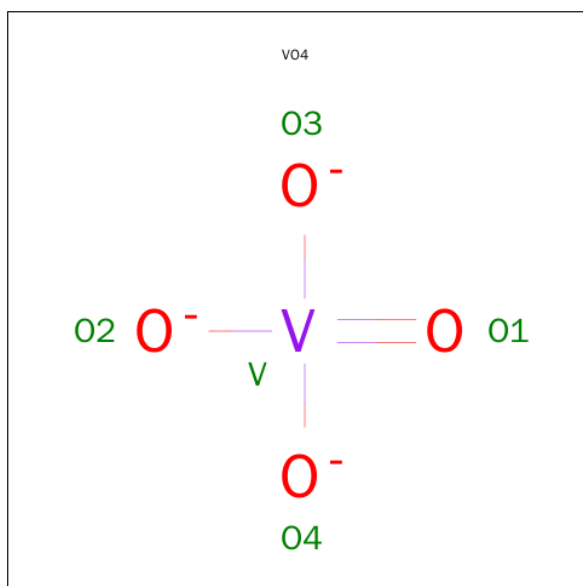
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	20	Total	C	N	O	P	0	0	0
			407	197	73	118	19			
3	H	20	Total	C	N	O	P	0	0	0
			407	197	73	118	19			
3	K	20	Total	C	N	O	P	0	0	0
			407	197	73	118	19			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	M	20	Total	C	N	O	P	0	0	0
			407	197	73	118	19			

- Molecule 4 is VANADATE ION (three-letter code: VO4) (formula: O₄V).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	V	0	0
			3	2	1		
4	D	1	Total	O	V	0	0
			3	2	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	75	Total	O	0	0
			75	75		
5	B	56	Total	O	0	0
			56	56		
5	C	58	Total	O	0	0
			58	58		
5	D	83	Total	O	0	0
			83	83		
5	E	8	Total	O	0	0
			8	8		
5	F	9	Total	O	0	0
			9	9		

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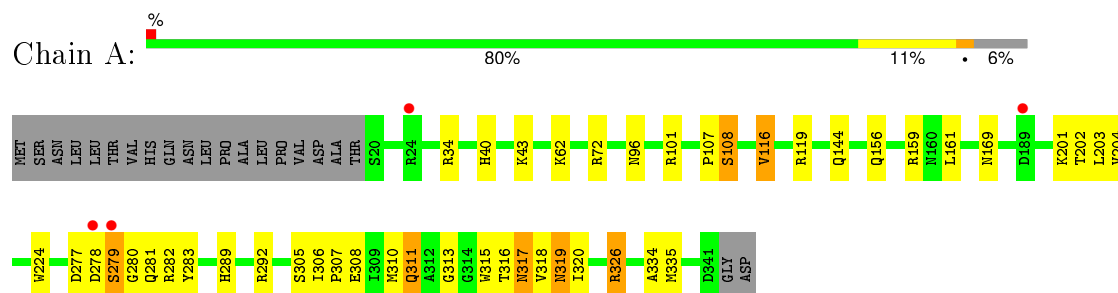
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	G	6	Total 6	O 6	0	0
5	H	7	Total 7	O 7	0	0
5	I	12	Total 12	O 12	0	0
5	K	7	Total 7	O 7	0	0
5	L	9	Total 9	O 9	0	0
5	M	13	Total 13	O 13	0	0

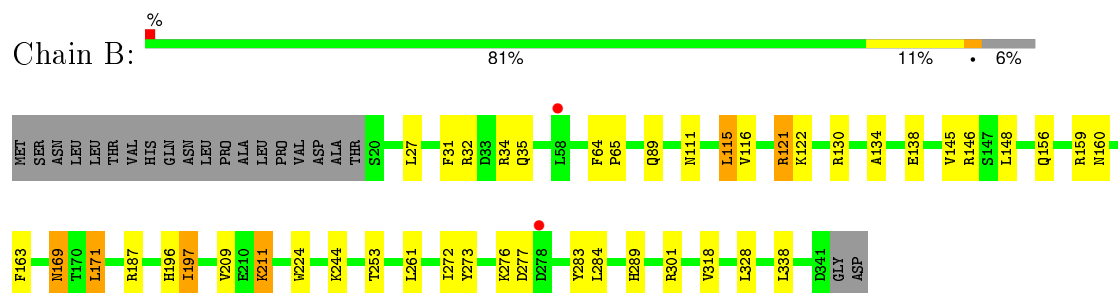
3 Residue-property plots [i](#)

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.

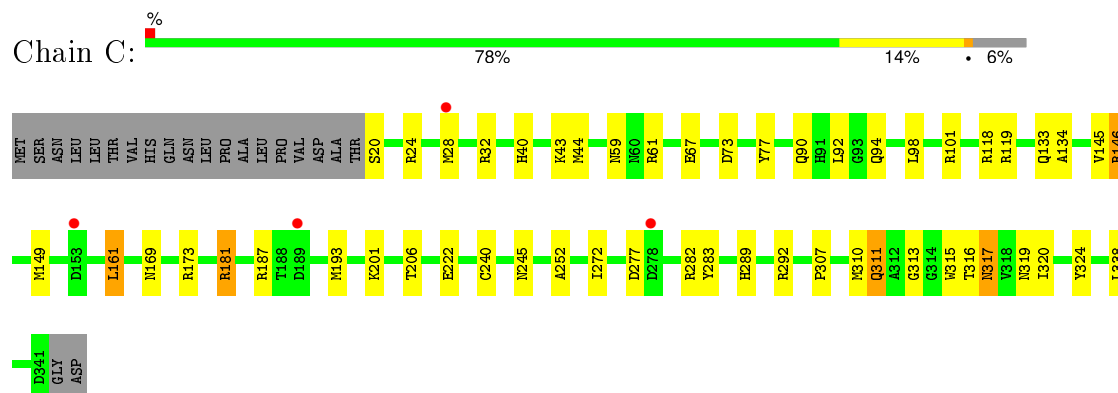
- Molecule 1: Recombinase cre



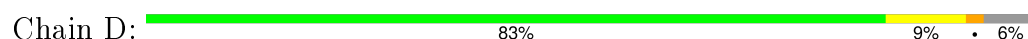
- Molecule 1: Recombinase cre

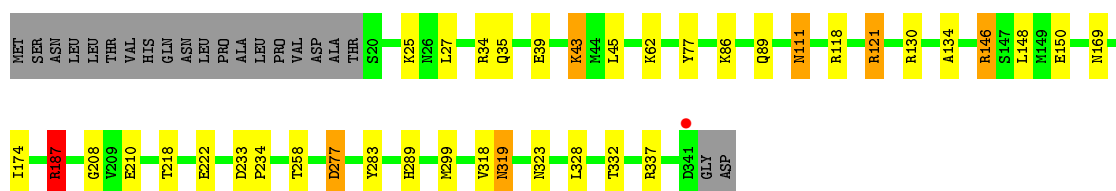


- Molecule 1: Recombinase cre

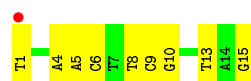


- Molecule 1: Recombinase cre

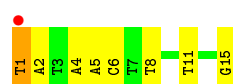




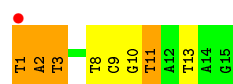
- Molecule 2: DNA (5'-D(*TP*AP*TP*AP*AP*CP*TP*TP*CP*GP*TP*AP*TP*AP*G)-3')



- Molecule 2: DNA (5'-D(*TP*AP*TP*AP*AP*CP*TP*TP*CP*GP*TP*AP*TP*AP*G)-3')



- Molecule 2: DNA (5'-D(*TP*AP*TP*AP*AP*CP*TP*TP*CP*GP*TP*AP*TP*AP*G)-3')



- Molecule 2: DNA (5'-D(*TP*AP*TP*AP*AP*CP*TP*TP*CP*GP*TP*AP*TP*AP*G)-3')

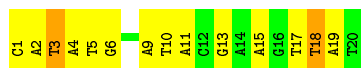


- Molecule 3: DNA (5'-D(*CP*AP*TP*AP*TP*GP*CP*TP*AP*TP*AP*CP*GP*AP*AP*GP*TP*TP*AP*T)-3')



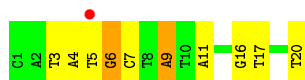
- Molecule 3: DNA (5'-D(*CP*AP*TP*AP*TP*GP*CP*TP*AP*TP*AP*CP*GP*AP*AP*GP*TP*TP*AP*T)-3')

Chain H:  30% 60% 10%



- Molecule 3: DNA (5'-D(*CP*AP*TP*AP*TP*GP*CP*TP*AP*TP*AP*CP*GP*AP*AP*GP*TP*TP*AP*T)-3')

Chain K:  5% 50% 40% 10%



- Molecule 3: DNA (5'-D(*CP*AP*TP*AP*TP*GP*CP*TP*AP*TP*AP*CP*GP*AP*AP*GP*TP*TP*AP*T)-3')

Chain M:  30% 40% 30%



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	136.05Å 136.05Å 218.43Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.57 – 2.29 49.54 – 2.29	Depositor EDS
% Data completeness (in resolution range)	97.1 (49.57-2.29) 97.1 (49.54-2.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.93 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.193 , 0.222 0.190 , 0.219	Depositor DCC
R_{free} test set	5105 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	30.7	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.8	EDS
Estimated twinning fraction	0.016 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 102010 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13393	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.91% 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, 0.375 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: VO4

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	0.63	0/2591	0.69	0/3493
1	B	0.61	0/2591	0.69	1/3493 (0.0%)
1	C	0.61	0/2591	0.70	2/3493 (0.1%)
1	D	0.67	0/2591	0.71	1/3493 (0.0%)
2	E	1.15	1/340 (0.3%)	1.87	10/523 (1.9%)
2	G	1.07	0/340	1.85	8/523 (1.5%)
2	I	1.20	1/340 (0.3%)	1.92	12/523 (2.3%)
2	L	1.06	0/340	2.46	14/523 (2.7%)
3	F	1.13	0/456	1.75	8/702 (1.1%)
3	H	1.07	1/456 (0.2%)	1.74	11/702 (1.6%)
3	K	1.09	0/456	1.90	14/702 (2.0%)
3	M	1.21	1/456 (0.2%)	2.02	15/702 (2.1%)
All	All	0.78	4/13548 (0.0%)	1.16	96/18872 (0.5%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	M	8	DT	N1-C6	5.18	1.41	1.38
3	H	9	DA	C3'-O3'	-5.17	1.37	1.44
2	I	11	DT	C3'-O3'	-5.12	1.37	1.44
2	E	1	DT	C1'-N1	5.04	1.55	1.49

All (96) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	2	DA	O5'-P-OP2	-25.25	80.40	110.70
2	L	2	DA	O5'-P-OP1	-20.87	85.66	110.70
2	L	2	DA	OP1-P-OP2	14.46	141.30	119.60
2	L	1	DT	OP2-P-O3'	-12.72	77.21	105.20
2	L	1	DT	O3'-P-O5'	11.92	126.65	104.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	K	5	DT	O4'-C1'-N1	11.88	116.31	108.00
3	M	11	DA	O4'-C1'-N9	-11.11	100.22	108.00
3	H	11	DA	O4'-C1'-N9	-10.90	100.37	108.00
2	L	1	DT	OP1-P-O3'	-10.18	82.81	105.20
3	M	15	DA	O4'-C1'-N9	9.40	114.58	108.00
3	F	5	DT	O4'-C1'-N1	9.24	114.47	108.00
2	E	5	DA	O4'-C1'-N9	8.82	114.17	108.00
2	G	8	DT	O4'-C1'-N1	-8.35	102.16	108.00
2	G	6	DC	O4'-C4'-C3'	-8.34	100.99	106.00
3	M	9	DA	O4'-C4'-C3'	-8.26	101.05	106.00
3	M	1	DC	O4'-C4'-C3'	-8.18	101.09	106.00
3	M	9	DA	O5'-P-OP2	-7.99	98.51	105.70
2	E	15	DG	C5-C6-O6	-7.93	123.84	128.60
2	I	1	DT	O4'-C1'-N1	7.92	113.54	108.00
3	H	6	DG	O4'-C1'-N9	7.90	113.53	108.00
2	L	5	DA	O4'-C1'-N9	7.83	113.48	108.00
2	I	8	DT	O4'-C1'-N1	-7.77	102.56	108.00
2	E	15	DG	C5-C6-N1	7.75	115.38	111.50
2	E	13	DT	O4'-C1'-N1	7.73	113.41	108.00
2	E	8	DT	O4'-C1'-N1	-7.71	102.60	108.00
1	B	171	LEU	CA-CB-CG	7.68	132.97	115.30
1	C	181	ARG	NE-CZ-NH2	-7.08	116.76	120.30
2	I	9	DC	P-O3'-C3'	7.05	128.16	119.70
3	F	3	DT	O4'-C1'-N1	7.03	112.92	108.00
3	K	9	DA	O5'-P-OP2	-6.94	99.46	105.70
2	G	1	DT	P-O3'-C3'	6.86	127.93	119.70
3	M	16	DG	O4'-C1'-N9	6.75	112.72	108.00
3	F	6	DG	O4'-C1'-N9	6.70	112.69	108.00
2	E	6	DC	P-O5'-C5'	6.65	131.54	120.90
3	F	4	DA	O4'-C1'-N9	6.61	112.62	108.00
1	D	187	ARG	NE-CZ-NH2	-6.59	117.01	120.30
2	G	4	DA	O4'-C1'-N9	-6.54	103.42	108.00
3	H	15	DA	P-O3'-C3'	6.54	127.54	119.70
2	L	15	DG	O4'-C1'-N9	-6.51	103.44	108.00
2	E	4	DA	O4'-C1'-N9	-6.47	103.47	108.00
3	F	18	DT	O4'-C1'-N1	-6.47	103.47	108.00
2	I	13	DT	P-O3'-C3'	6.41	127.39	119.70
3	H	17	DT	C5-C4-O4	-6.39	120.43	124.90
3	M	3	DT	C4-C5-C7	6.23	122.74	119.00
2	L	1	DT	O4'-C1'-N1	6.21	112.35	108.00
2	E	15	DG	C6-N1-C2	-6.19	121.39	125.10
3	K	3	DT	P-O3'-C3'	6.14	127.07	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	K	20	DT	O4'-C1'-N1	6.14	112.30	108.00
3	K	9	DA	O5'-P-OP1	6.09	118.01	110.70
3	K	7	DC	P-O3'-C3'	6.09	127.00	119.70
3	F	15	DA	P-O3'-C3'	6.04	126.94	119.70
2	E	6	DC	C5'-C4'-O4'	-6.03	97.85	109.30
3	K	5	DT	C1'-O4'-C4'	-6.02	104.08	110.10
2	I	10	DG	O4'-C1'-N9	5.97	112.18	108.00
3	H	3	DT	O4'-C1'-N1	5.92	112.15	108.00
2	G	5	DA	O4'-C1'-N9	5.88	112.12	108.00
2	L	13	DT	P-O3'-C3'	5.86	126.74	119.70
2	I	3	DT	N3-C2-O2	-5.84	118.80	122.30
3	M	20	DT	C4-C5-C7	5.72	122.43	119.00
3	H	10	DT	N3-C4-O4	5.65	123.29	119.90
2	I	1	DT	P-O3'-C3'	5.61	126.43	119.70
3	H	10	DT	C5-C4-O4	-5.49	121.06	124.90
2	I	11	DT	C4-C5-C7	5.48	122.28	119.00
3	M	19	DA	N1-C6-N6	5.47	121.88	118.60
3	F	9	DA	C5'-C4'-C3'	5.43	123.87	114.10
3	H	18	DT	C4-C5-C7	5.40	122.24	119.00
2	G	11	DT	P-O3'-C3'	5.39	126.17	119.70
2	E	5	DA	C1'-O4'-C4'	-5.36	104.74	110.10
3	F	19	DA	O4'-C1'-N9	5.35	111.74	108.00
3	M	5	DT	N3-C4-O4	5.33	123.10	119.90
3	K	7	DC	C5-C4-N4	5.31	123.92	120.20
2	L	8	DT	C6-C5-C7	-5.31	119.71	122.90
2	I	8	DT	C2-N3-C4	-5.29	124.03	127.20
3	K	4	DA	P-O3'-C3'	5.28	126.04	119.70
3	M	5	DT	C5-C4-O4	-5.27	121.21	124.90
2	G	15	DG	C8-N9-C4	-5.27	104.29	106.40
2	I	3	DT	P-O3'-C3'	5.26	126.02	119.70
2	I	1	DT	C4-C5-C7	5.26	122.16	119.00
3	K	17	DT	N3-C4-O4	5.23	123.04	119.90
3	K	6	DG	P-O3'-C3'	5.23	125.98	119.70
3	M	18	DT	N3-C2-O2	-5.22	119.17	122.30
2	L	5	DA	C1'-O4'-C4'	-5.22	104.88	110.10
3	M	10	DT	C5-C4-O4	-5.20	121.26	124.90
3	M	4	DA	N1-C6-N6	5.18	121.71	118.60
3	H	4	DA	N1-C6-N6	5.18	121.71	118.60
2	I	2	DA	P-O3'-C3'	-5.14	113.54	119.70
3	H	13	DG	C5-C6-N1	5.09	114.05	111.50
2	L	7	DT	N3-C2-O2	-5.06	119.26	122.30
2	G	5	DA	C1'-O4'-C4'	-5.05	105.05	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	181	ARG	NE-CZ-NH1	5.04	122.82	120.30
3	K	16	DG	P-O5'-C5'	-5.04	112.83	120.90
3	K	11	DA	O4'-C1'-N9	-5.03	104.48	108.00
3	H	5	DT	O4'-C1'-N1	-5.02	104.49	108.00
2	L	8	DT	O4'-C1'-N1	-5.02	104.49	108.00
3	M	10	DT	N3-C4-O4	5.01	122.91	119.90
3	K	6	DG	O4'-C1'-N9	5.00	111.50	108.00

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	2550	0	2571	33	0
1	B	2550	0	2570	26	0
1	C	2550	0	2571	39	0
1	D	2550	0	2570	29	0
2	E	304	0	172	1	0
2	G	304	0	173	1	0
2	I	304	0	172	5	0
2	L	304	0	173	1	0
3	F	407	0	228	5	0
3	H	407	0	229	6	0
3	K	407	0	228	3	0
3	M	407	0	229	12	0
4	B	3	0	0	0	0
4	D	3	0	0	0	0
5	A	75	0	0	3	0
5	B	56	0	0	2	0
5	C	58	0	0	4	0
5	D	83	0	0	3	0
5	E	8	0	0	0	0
5	F	9	0	0	1	0
5	G	6	0	0	0	0
5	H	7	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	I	12	0	0	1	0
5	K	7	0	0	0	0
5	L	9	0	0	0	0
5	M	13	0	0	1	0
All	All	13393	0	11886	135	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 5.

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:B:121:ARG:HH11	1:B:121:ARG:HG2	1.33	0.93
3:F:2:DA:H2"	3:F:3:DT:H5"	1.52	0.88
1:D:187:ARG:NH2	1:D:222:GLU:OE1	2.13	0.81
1:B:121:ARG:HG2	1:B:121:ARG:NH1	1.94	0.80
1:C:201:LYS:HE2	3:K:9:DA:H5'	1.67	0.76
1:A:62:LYS:HD3	5:A:409:HOH:O	1.86	0.73
1:D:89:GLN:HE22	1:D:121:ARG:NH2	1.86	0.72
1:C:307:PRO:HA	1:C:310:MET:HE2	1.72	0.71
1:A:311:GLN:HG2	1:B:328:LEU:HD12	1.74	0.69
1:C:307:PRO:HA	1:C:310:MET:CE	2.23	0.68
1:D:118:ARG:NH2	3:M:3:DT:H4'	2.09	0.68
1:A:334:ALA:HB3	1:D:299:MET:CE	2.24	0.67
3:K:6:DG:H21	3:M:2:DA:H1'	1.61	0.66
1:A:289:HIS:ND1	5:A:354:HOH:O	2.29	0.66
1:B:111:ASN:O	1:B:115:LEU:HB2	1.95	0.66
1:C:187:ARG:NH2	1:C:222:GLU:OE1	2.27	0.65
1:D:89:GLN:HE22	1:D:121:ARG:HH22	1.42	0.65
1:B:273:TYR:O	5:B:344:HOH:O	2.14	0.64
1:B:121:ARG:HG3	1:B:122:LYS:N	2.14	0.63
3:M:3:DT:H2"	3:M:4:DA:OP2	1.98	0.62
1:C:201:LYS:HD3	3:M:2:DA:OP1	1.98	0.62
3:M:3:DT:H6	3:M:3:DT:OP2	1.82	0.62
3:M:19:DA:C8	3:M:19:DA:H5"	2.34	0.61
1:D:332:THR:O	1:D:337:ARG:NH1	2.32	0.61
1:C:315:TRP:HA	3:M:1:DC:H5'	1.81	0.60
1:B:156:GLN:HE21	1:B:160:ASN:HD21	1.48	0.60
1:A:96:ASN:HD21	1:A:108:SER:H	1.49	0.60
1:D:89:GLN:NE2	1:D:121:ARG:NH1	2.51	0.58
1:D:146:ARG:O	1:D:150:GLU:HB2	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:LYS:HE2	3:F:9:DA:H5'	1.85	0.56
1:C:43:LYS:HG3	1:C:44:MET:N	2.21	0.56
1:A:334:ALA:HB3	1:D:299:MET:HE1	1.87	0.56
1:B:134:ALA:HA	1:B:283:TYR:CD2	2.41	0.56
1:A:279:SER:O	1:A:281:GLN:N	2.33	0.55
1:D:89:GLN:HE21	1:D:121:ARG:NH1	2.04	0.55
3:M:19:DA:H8	3:M:19:DA:H5''	1.71	0.55
2:G:1:DT:H4'	2:G:2:DA:OP1	2.07	0.55
1:A:335:MET:HG2	1:D:299:MET:HE3	1.88	0.55
1:C:317:ASN:HB3	1:C:320:ILE:HG12	1.88	0.55
1:D:89:GLN:NE2	1:D:121:ARG:HH12	2.03	0.55
1:A:40:HIS:HE1	5:F:28:HOH:O	1.90	0.55
2:I:1:DT:H1'	2:I:2:DA:OP1	2.09	0.53
1:A:144:GLN:NE2	5:A:351:HOH:O	2.40	0.53
1:D:277:ASP:C	1:D:277:ASP:OD1	2.47	0.53
1:A:96:ASN:HD22	1:A:107:PRO:HD2	1.73	0.53
1:B:122:LYS:HE2	3:F:8:DT:OP1	2.10	0.52
3:M:2:DA:H2''	3:M:3:DT:H5'	1.92	0.52
1:C:67:GLU:CG	5:C:402:HOH:O	2.58	0.51
1:D:89:GLN:HE21	1:D:121:ARG:HH12	1.58	0.51
1:B:121:ARG:NH1	3:H:3:DT:H3'	2.26	0.51
1:C:149:MET:CE	1:C:161:LEU:HG	2.40	0.51
1:A:282:ARG:O	1:A:283:TYR:HB2	2.11	0.51
1:C:193:MET:CE	1:C:222:GLU:HG3	2.42	0.50
1:C:90:GLN:NE2	1:C:94:GLN:HE21	2.09	0.50
1:C:206:THR:HG21	1:C:316:THR:CG2	2.42	0.50
1:C:24:ARG:O	1:C:28:MET:HG2	2.11	0.50
1:C:193:MET:HE3	1:C:222:GLU:HG3	1.94	0.49
1:C:317:ASN:C	1:C:317:ASN:HD22	2.16	0.49
1:A:159:ARG:HB2	1:A:224:TRP:CZ3	2.47	0.49
1:C:289:HIS:HE1	1:C:292:ARG:NH1	2.11	0.48
1:A:334:ALA:HB3	1:D:299:MET:HE2	1.95	0.48
1:D:319:ASN:HB2	5:D:344:HOH:O	2.14	0.48
1:A:202:THR:HG22	1:A:204:VAL:H	1.77	0.48
1:C:317:ASN:HD22	1:C:319:ASN:H	1.62	0.48
2:I:3:DT:P	5:I:135:HOH:O	2.72	0.48
1:C:61:ARG:NH2	1:C:73:ASP:OD1	2.39	0.48
1:A:289:HIS:CE1	1:A:292:ARG:HE	2.33	0.47
1:B:35:GLN:HB2	1:C:119:ARG:HG3	1.96	0.47
1:C:317:ASN:ND2	1:C:319:ASN:H	2.13	0.47
1:C:311:GLN:HG2	1:D:328:LEU:HD12	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:169:ASN:C	1:B:169:ASN:HD22	2.18	0.47
3:H:2:DA:H2'	3:H:3:DT:C6	2.50	0.46
1:B:197:ILE:HD13	1:B:211:LYS:HG3	1.96	0.46
1:A:305:SER:OG	1:A:308:GLU:HG3	2.16	0.46
1:D:218:THR:O	1:D:222:GLU:HG3	2.16	0.46
2:I:1:DT:H2''	2:I:2:DA:C8	2.51	0.46
1:A:317:ASN:HD22	1:A:319:ASN:H	1.64	0.46
3:H:18:DT:H2''	3:H:19:DA:N7	2.31	0.45
1:B:276:LYS:HD2	1:B:284:LEU:HB2	1.98	0.45
1:D:43:LYS:HE3	2:I:11:DT:O4	2.16	0.45
1:B:89:GLN:HE22	1:B:121:ARG:NH2	2.13	0.45
1:B:138:GLU:CD	1:B:301:ARG:HH22	2.19	0.45
5:D:365:HOH:O	3:M:9:DA:H5''	2.15	0.45
1:C:40:HIS:O	1:C:43:LYS:HG2	2.15	0.45
1:A:310:MET:SD	1:A:318:VAL:HG12	2.57	0.45
1:C:324:TYR:OH	2:L:15:DG:OP1	2.34	0.45
1:B:163:PHE:CE1	1:B:261:LEU:HD22	2.52	0.45
1:B:159:ARG:HB2	1:B:224:TRP:CZ3	2.51	0.45
1:A:305:SER:HB2	1:A:307:PRO:HD2	1.99	0.45
1:D:318:VAL:HG12	5:D:423:HOH:O	2.16	0.45
1:A:326:ARG:NH2	1:D:208:GLY:O	2.50	0.45
2:I:2:DA:H5''	2:I:2:DA:H8	1.82	0.44
1:C:59:ASN:O	1:C:61:ARG:HD3	2.17	0.44
1:C:240:CYS:HB3	5:C:377:HOH:O	2.17	0.44
1:A:326:ARG:HG3	1:D:210:GLU:HG3	1.98	0.44
1:A:101:ARG:NH1	1:B:115:LEU:HD13	2.33	0.44
1:B:31:PHE:O	1:B:34:ARG:HB3	2.16	0.44
1:B:145:VAL:HG22	1:B:272:ILE:HD11	2.01	0.43
1:D:134:ALA:HA	1:D:283:TYR:CD2	2.53	0.43
3:F:9:DA:H2'	3:F:10:DT:C6	2.54	0.43
1:C:282:ARG:O	1:C:283:TYR:HB2	2.18	0.43
1:B:64:PHE:HA	1:B:65:PRO:C	2.39	0.43
1:A:306:ILE:HB	1:A:307:PRO:HD3	2.00	0.43
1:A:201:LYS:HG2	3:H:2:DA:OP1	2.19	0.42
1:D:89:GLN:HE22	1:D:121:ARG:CZ	2.32	0.42
1:A:119:ARG:HG3	1:D:35:GLN:HB2	2.01	0.42
1:C:67:GLU:HG2	5:C:402:HOH:O	2.17	0.42
1:C:173:ARG:NH1	1:C:201:LYS:HD2	2.35	0.42
1:C:181:ARG:NH2	1:C:252:ALA:O	2.53	0.42
1:B:121:ARG:CG	1:B:121:ARG:NH1	2.67	0.42
1:A:279:SER:OG	1:A:279:SER:O	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:3:DT:OP2	3:H:3:DT:H6	2.03	0.42
1:C:134:ALA:HA	1:C:283:TYR:CD2	2.54	0.42
1:A:72:ARG:HG3	1:A:116:VAL:HG11	2.02	0.42
1:C:173:ARG:CZ	1:C:201:LYS:HD2	2.50	0.42
3:K:6:DG:N2	3:M:2:DA:H1'	2.33	0.41
1:D:89:GLN:NE2	1:D:121:ARG:CZ	2.82	0.41
1:B:187:ARG:HH11	1:B:187:ARG:HG3	1.85	0.41
1:C:307:PRO:CA	1:C:310:MET:HE2	2.46	0.41
1:C:145:VAL:HG22	1:C:272:ILE:HD11	2.02	0.41
1:C:149:MET:HE2	1:C:161:LEU:HG	2.03	0.41
1:D:174:ILE:HD12	1:D:258:THR:HB	2.02	0.41
1:A:313:GLY:HA3	1:A:315:TRP:CZ3	2.55	0.41
2:E:9:DC:H2''	2:E:10:DG:C8	2.56	0.41
1:C:146:ARG:NH2	5:C:398:HOH:O	2.46	0.41
1:B:196:HIS:HE1	5:B:397:HOH:O	2.02	0.41
1:A:311:GLN:CG	1:B:328:LEU:HD12	2.49	0.41
1:C:101:ARG:HH11	1:D:111:ASN:ND2	2.19	0.41
1:C:313:GLY:HA3	1:C:315:TRP:CZ3	2.56	0.41
1:A:320:ILE:HD11	3:H:1:DC:C6	2.56	0.41
1:D:233:ASP:HA	1:D:234:PRO:HD3	1.96	0.41
3:M:20:DT:H4'	5:M:331:HOH:O	2.21	0.40
1:A:317:ASN:ND2	1:A:319:ASN:H	2.17	0.40
1:C:292:ARG:NH2	1:C:313:GLY:O	2.52	0.40
3:F:1:DC:H2'	3:F:2:DA:C8	2.56	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	320/343 (93%)	314 (98%)	4 (1%)	2 (1%)	30 36

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	320/343 (93%)	315 (98%)	4 (1%)	1 (0%)	46	57
1	C	320/343 (93%)	314 (98%)	6 (2%)	0	100	100
1	D	320/343 (93%)	316 (99%)	2 (1%)	2 (1%)	30	36
All	All	1280/1372 (93%)	1259 (98%)	16 (1%)	5 (0%)	39	48

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	279	SER
1	A	280	GLY
1	D	277	ASP
1	D	34	ARG
1	B	244	LYS

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	269/287 (94%)	254 (94%)	15 (6%)	26	35
1	B	269/287 (94%)	251 (93%)	18 (7%)	20	26
1	C	269/287 (94%)	254 (94%)	15 (6%)	26	35
1	D	269/287 (94%)	251 (93%)	18 (7%)	20	26
All	All	1076/1148 (94%)	1010 (94%)	66 (6%)	23	30

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	ARG
1	A	43	LYS
1	A	108	SER
1	A	116	VAL
1	A	156	GLN
1	A	161	LEU

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Mol	Chain	Res	Type
1	A	169	ASN
1	A	203	LEU
1	A	277	ASP
1	A	278	ASP
1	A	311	GLN
1	A	316	THR
1	A	317	ASN
1	A	319	ASN
1	A	326	ARG
1	B	27	LEU
1	B	32	ARG
1	B	115	LEU
1	B	116	VAL
1	B	121	ARG
1	B	130	ARG
1	B	146	ARG
1	B	148	LEU
1	B	169	ASN
1	B	171	LEU
1	B	197	ILE
1	B	209	VAL
1	B	211	LYS
1	B	253	THR
1	B	277	ASP
1	B	289	HIS
1	B	318	VAL
1	B	338	LEU
1	C	20	SER
1	C	32	ARG
1	C	77	TYR
1	C	92	LEU
1	C	98	LEU
1	C	118	ARG
1	C	133	GLN
1	C	146	ARG
1	C	161	LEU
1	C	169	ASN
1	C	245	ASN
1	C	277	ASP
1	C	311	GLN
1	C	317	ASN
1	C	338	LEU

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Mol	Chain	Res	Type
1	D	25	LYS
1	D	27	LEU
1	D	39	GLU
1	D	43	LYS
1	D	45	LEU
1	D	62	LYS
1	D	77	TYR
1	D	86	LYS
1	D	111	ASN
1	D	121	ARG
1	D	130	ARG
1	D	146	ARG
1	D	148	LEU
1	D	169	ASN
1	D	187	ARG
1	D	289	HIS
1	D	319	ASN
1	D	323	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	HIS
1	A	60	ASN
1	A	96	ASN
1	A	144	GLN
1	A	235	ASN
1	A	317	ASN
1	A	319	ASN
1	A	323	ASN
1	B	89	GLN
1	B	156	GLN
1	B	196	HIS
1	B	319	ASN
1	B	323	ASN
1	C	59	ASN
1	C	89	GLN
1	C	90	GLN
1	C	133	GLN
1	C	144	GLN
1	C	245	ASN
1	C	317	ASN

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Mol	Chain	Res	Type
1	C	323	ASN
1	D	89	GLN
1	D	111	ASN
1	D	144	GLN
1	D	281	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

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	$\# Z > 2$
4	VO4	B	690	1,3,2	0,2,4	0.00	-	0,1,6	0.00	-
4	VO4	D	690	1,3,2	0,2,4	0.00	-	0,1,6	0.00	-

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
4	VO4	B	690	1,3,2	-	0/0/0/0	0/0/0/0
4	VO4	D	690	1,3,2	-	0/0/0/0	0/0/0/0

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

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	322/343 (93%)	-0.39	4 (1%) 81 85	16, 27, 44, 59	0
1	B	322/343 (93%)	-0.42	2 (0%) 90 93	16, 28, 42, 54	0
1	C	322/343 (93%)	-0.36	4 (1%) 81 85	16, 28, 45, 58	0
1	D	322/343 (93%)	-0.46	1 (0%) 94 96	13, 24, 38, 52	0
2	E	15/15 (100%)	-0.30	1 (6%) 21 29	17, 28, 70, 98	0
2	G	15/15 (100%)	-0.13	1 (6%) 21 29	19, 26, 75, 105	0
2	I	15/15 (100%)	-0.39	1 (6%) 21 29	15, 23, 64, 92	0
2	L	15/15 (100%)	-0.38	1 (6%) 21 29	19, 29, 77, 102	0
3	F	20/20 (100%)	-0.31	0 100 100	15, 41, 63, 76	0
3	H	20/20 (100%)	-0.49	0 100 100	16, 35, 60, 65	0
3	K	20/20 (100%)	-0.21	1 (5%) 32 41	18, 38, 71, 81	0
3	M	20/20 (100%)	-0.40	0 100 100	14, 33, 63, 65	0
All	All	1428/1512 (94%)	-0.40	16 (1%) 82 86	13, 27, 46, 105	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	G	1	DT	6.5
2	E	1	DT	5.8
2	L	1	DT	5.6
1	A	189	ASP	3.9
1	A	278	ASP	3.6
2	I	1	DT	3.3
1	C	278	ASP	3.3
1	C	153	ASP	3.1
1	D	341	ASP	2.8
1	B	278	ASP	2.7
1	C	28	MET	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	24	ARG	2.7
1	B	58	LEU	2.6
3	K	5	DT	2.5
1	C	189	ASP	2.3
1	A	279	SER	2.3

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, 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
4	VO4	B	690	3/5	1.00	0.13	0.91	15,15,17,18	0
4	VO4	D	690	3/5	1.00	0.13	0.65	15,15,15,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.