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PDB ID	:	6DBQ
EMDB ID	:	EMD-7847
Title	:	Cryo-EM structure of RAG in complex with 12-RSS and 23-RSS substrate
		DNAs
Authors	:	Wu, H.; Liao, M.; Ru, H.; Mi, W.
Deposited on	:	2018-05-03
Resolution	:	4.22 Å(reported)

This is a Full wwPDB EM 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/EMValidationReportHelp with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

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

:	0.0.1. dev 70
:	4.02b-467
:	20191225.v01 (using entries in the PDB archive December 25th 2019)
:	1.9.13
:	Engh & Huber (2001)
:	Parkinson et al. (1996)
:	2.36
	: : : : :

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $ELECTRON\ MICROSCOPY$

The reported resolution of this entry is 4.22 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f EM} {f structures} \ (\#{f Entries})$
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion < 40%). The numeric value is given above the bar.

Mol	Chain	Length		Quality of chain					
1	А	1159	- 29%	23%	•		46%		
1	С	1159	29%	22%	·		47%		
2	В	533	•		30%	•	34%		
2	D	533	33%		32%	•	34%		
3	Е	50	24%			70%		•	
4	F	50	20%			76%			
5	G	61	46%			70%		•	
6	Н	61	38%			61%		•	



2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 19956 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 Recombination activating gene 1 - MBP chimera.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	А	621	Total 4999	C 3128	N 899	O 935	S 37	0	0
1	С	616	Total 4972	C 3113	N 893	O 929	S 37	0	0

Chain	Residue	Modelled	Actual	Comment	Reference
А	-127	MET	-	initiating methionine	UNP POAEX9
А	-126	GLY	-	expression tag	UNP POAEX9
А	-125	SER	-	expression tag	UNP POAEX9
А	-124	SER	-	expression tag	UNP POAEX9
А	-123	HIS	-	expression tag	UNP POAEX9
А	-122	HIS	-	expression tag	UNP POAEX9
А	-121	HIS	-	expression tag	UNP POAEX9
А	-120	HIS	-	expression tag	UNP POAEX9
A	-119	HIS	-	expression tag	UNP P0AEX9
А	-118	HIS	-	expression tag	UNP POAEX9
А	-117	GLY	-	expression tag	UNP P0AEX9
А	-116	THR	-	expression tag	UNP POAEX9
А	-115	LYS	-	expression tag	UNP POAEX9
А	-114	THR	-	expression tag	UNP P0AEX9
A	251	GLY	-	linker	UNP P0AEX9
А	252	THR	-	linker	UNP P0AEX9
А	253	ASP	-	linker	UNP POAEX9
А	254	TYR	-	linker	UNP POAEX9
А	255	ASP	-	linker	UNP POAEX9
А	256	ILE	-	linker	UNP POAEX9
A	257	PRO	-	linker	UNP P0AEX9
A	258	THR	-	linker	UNP POAEX9
A	259	THR	-	linker	UNP POAEX9
A	260	LEU	-	linker	UNP POAEX9
A	261	GLU	-	linker	UNP POAEX9
A	262	VAL	-	linker	UNP POAEX9

There are 68 discrepancies between the modelled and reference sequences:



Chain	Residue	Modelled	Actual	Comment	Reference
А	263	LEU	-	linker	UNP POAEX9
A	264	PHE	-	linker	UNP POAEX9
A	265	GLN	-	linker	UNP POAEX9
A	266	GLY	-	linker	UNP POAEX9
А	267	PRO	-	linker	UNP POAEX9
A	268	LEU	-	linker	UNP POAEX9
А	269	GLY	-	linker	UNP POAEX9
А	270	SER	-	linker	UNP POAEX9
С	-127	MET	-	initiating methionine	UNP P0AEX9
С	-126	GLY	-	expression tag	UNP POAEX9
С	-125	SER	-	expression tag	UNP POAEX9
С	-124	SER	-	expression tag	UNP POAEX9
С	-123	HIS	-	expression tag	UNP POAEX9
С	-122	HIS	-	expression tag	UNP POAEX9
С	-121	HIS	-	expression tag	UNP POAEX9
С	-120	HIS	-	expression tag	UNP POAEX9
С	-119	HIS	-	expression tag	UNP POAEX9
С	-118	HIS	-	expression tag	UNP POAEX9
С	-117	GLY	-	expression tag	UNP POAEX9
С	-116	THR	-	expression tag	UNP POAEX9
С	-115	LYS	-	expression tag	UNP POAEX9
С	-114	THR	-	expression tag	UNP POAEX9
С	251	GLY	-	linker	UNP POAEX9
С	252	THR	-	linker	UNP POAEX9
С	253	ASP	-	linker	UNP P0AEX9
С	254	TYR	-	linker	UNP POAEX9
С	255	ASP	-	linker	UNP P0AEX9
С	256	ILE	-	linker	UNP P0AEX9
С	257	PRO	-	linker	UNP P0AEX9
C	258	THR	-	linker	UNP P0AEX9
С	259	THR	-	linker	UNP P0AEX9
C	260	LEU	-	linker	UNP P0AEX9
С	261	GLU	-	linker	UNP P0AEX9
С	262	VAL	-	linker	UNP P0AEX9
C	263	LEU	-	linker	UNP POAEX9
C	264	PHE	-	linker	UNP P0AEX9
С	265	GLN	-	linker	UNP POAEX9
C	266	GLY	-	linker	UNP POAEX9
C	267	PRO	-	linker	UNP POAEX9
С	268	LEU	-	linker	UNP POAEX9
C	269	GLY	-	linker	UNP POAEX9
С	270	SER	-	linker	UNP POAEX9



• Molecule 2 is a protein called Recombination activating gene 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	В	351	Total	С	Ν	0	\mathbf{S}	0	0
	551	2714	1716	470	509	19	0	0	
9	Л	351	Total	С	Ν	0	S	0	0
		001	2714	1716	470	509	19	U	U

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference	
В	-2	GLY	-	expression tag	UNP Q1RLW7	
В	-1	GLY	-	expression tag	UNP Q1RLW7	
В	0	SER	-	expression tag	UNP Q1RLW7	
D	-2	GLY	-	expression tag	UNP Q1RLW7	
D	-1	GLY	-	expression tag	UNP Q1RLW7	
D	0	SER	-	expression tag	UNP Q1RLW7	

• Molecule 3 is a DNA chain called Molecule name: Forward strand of 12-RSS substrate DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	Е	50	Total 1023	C 486	N 192	O 295	Р 50	0	0

• Molecule 4 is a DNA chain called Molecule name: Reverse strand of 12-RSS substrate DNA.

Mol	Chain	Residues		A	toms	AltConf	Trace		
4	F	50	Total 1027	C 489	N 183	O 305	Р 50	0	0

• Molecule 5 is a DNA chain called Molecule name: Forward strand of 23-RSS substrate DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	G	61	Total 1245	C 593	N 223	O 368	Р 61	0	0

• Molecule 6 is a DNA chain called Reverse strand of 23-RSS substrate DNA.

Mol	Chain	Residues	Atoms				AltConf	Trace	
6	Н	61	Total 1256	C 596	N 235	0 364	Р 61	0	0

• Molecule 7 is ZINC ION (three-letter code: ZN) (formula: Zn).



Mol	Chain	Residues	Atoms	AltConf
7	А	1	Total Zn 1 1	0
7	С	1	Total Zn 1 1	0

• Molecule 8 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	AltConf
8	А	2	Total Ca 2 2	0
8	С	2	Total Ca 2 2	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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 atom inclusion in map 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 diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). 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: Recombination activating gene 1 - MBP chimera









• Molecule 2: Recombination activating gene 2





• Molecule 6: Reverse strand of 23-RSS substrate DNA







4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	48002	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE	Depositor
	CORRECTION	
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose $(e^-/\text{\AA}^2)$	47	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.182	Depositor
Minimum map value	-0.085	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.009	Depositor
Recommended contour level	0.03	Depositor
Map size (Å)	237.69601, 237.69601, 237.69601	wwPDB
Map dimensions	192, 192, 192	wwPDB
Map angles $(^{\circ})$	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.238, 1.238, 1.238	Depositor



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: CA, ZN

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).

Mal	Chain	Bo	ond lengths	B	Bond angles		
	Ullalli	RMSZ	# Z > 5	RMSZ	# Z > 5		
1	А	0.32	0/5097	0.59	4/6856~(0.1%)		
1	С	0.33	0/5067	0.66	7/6812~(0.1%)		
2	В	0.34	0/2784	0.63	5/3784~(0.1%)		
2	D	0.35	1/2784~(0.0%)	0.61	3/3784~(0.1%)		
3	Е	0.61	0/1148	0.95	1/1768~(0.1%)		
4	F	0.62	0/1150	0.99	0/1774		
5	G	0.62	1/1394~(0.1%)	0.91	0/2148		
6	Н	0.65	0/1410	1.00	1/2175~(0.0%)		
All	All	0.42	2/20834~(0.0%)	0.73	$21/29101 \ (0.1\%)$		

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	А	0	1
1	С	0	1
2	В	0	3
2	D	0	1
All	All	0	6

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
2	D	83	GLN	CG-CD	-5.47	1.38	1.51
5	G	35	DC	C1'-N1	5.12	1.55	1.49

All (21) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	С	432	ASN	CB-CG-OD1	24.80	171.21	121.60
1	С	432	ASN	CB-CG-ND2	-20.88	66.60	116.70
1	А	423	GLN	CG-CD-OE1	17.99	157.58	121.60
1	А	423	GLN	CG-CD-NE2	-15.65	79.13	116.70
2	В	344	HIS	N-CA-CB	-11.29	90.27	110.60
1	С	816	LEU	C-N-CA	-9.46	98.06	121.70
2	D	83	GLN	CA-CB-CG	8.94	133.07	113.40
2	В	344	HIS	CB-CA-C	8.28	126.95	110.40
2	D	83	GLN	CG-CD-NE2	-8.03	97.44	116.70
2	В	10	ASN	CB-CA-C	-7.97	94.46	110.40
2	D	83	GLN	CB-CG-CD	7.04	129.90	111.60
6	Н	45	DG	O4'-C1'-N9	6.75	112.72	108.00
1	С	432	ASN	OD1-CG-ND2	-6.58	106.78	121.90
2	В	339	PRO	C-N-CD	-6.03	107.34	120.60
1	А	423	GLN	CA-CB-CG	5.94	126.46	113.40
1	А	423	GLN	OE1-CD-NE2	-5.90	108.33	121.90
1	С	817	HIS	CB-CG-CD2	-5.61	113.41	130.80
1	С	817	HIS	N-CA-CB	5.46	120.42	110.60
2	В	9	VAL	C-N-CA	5.27	134.87	121.70
3	Е	29	DG	O4'-C1'-N9	5.17	111.62	108.00
1	С	516	PHE	C-N-CA	-5.02	109.14	121.70

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	423	GLN	Sidechain
2	В	10	ASN	Sidechain, Mainchain
2	В	344	HIS	Sidechain
1	С	817	HIS	Sidechain
2	D	83	GLN	Sidechain

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	А	4999	0	4954	263	0
1	С	4972	0	4937	245	0



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	В	2714	0	2665	160	0
2	D	2714	0	2665	173	0
3	Е	1023	0	561	69	0
4	F	1027	0	566	67	0
5	G	1245	0	687	99	0
6	Н	1256	0	685	50	0
7	А	1	0	0	0	0
7	С	1	0	0	0	0
8	А	2	0	0	0	0
8	C	2	0	0	0	0
All	All	19956	0	17720	1020	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 27.

All (1020) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom 1	Atom 2	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
5:G:12:DT:O4	6:H:50:DA:N6	1.60	1.34
2:D:123:ARG:NH2	2:D:125:GLU:OE2	1.70	1.25
2:D:232:ARG:HH11	2:D:234:ILE:CG1	1.49	1.23
1:C:435:LYS:NZ	1:C:439:GLU:OE2	1.81	1.12
1:A:789:GLU:OE2	1:A:797:ARG:NH1	1.84	1.10
2:D:137:ARG:NH1	2:D:178:CYS:SG	2.25	1.09
2:D:338:THR:HG22	2:D:340:PRO:HA	1.30	1.08
1:A:447:LYS:NZ	1:C:463:GLU:OE2	1.88	1.05
1:A:505:ARG:NH1	1:C:1029:LYS:O	1.89	1.05
1:A:740:LEU:HD11	1:A:770:ARG:HH12	1.24	1.00
2:D:232:ARG:HH11	2:D:234:ILE:HG12	1.27	0.98
2:B:99:ASN:OD1	2:B:101:GLU:OE2	1.81	0.97
2:D:232:ARG:HH11	2:D:234:ILE:HG13	1.30	0.96
2:B:97:THR:OG1	2:B:101:GLU:OE2	1.86	0.94
5:G:10:DT:C4	5:G:11:DG:C6	2.57	0.93
1:C:523:ARG:NH1	3:E:24:DC:OP2	2.02	0.92
2:D:232:ARG:NH1	2:D:234:ILE:CG1	2.31	0.92
5:G:32:DC:N4	6:H:30:DG:O6	2.02	0.92
5:G:30:DC:H2"	5:G:31:DT:H71	1.51	0.92
2:B:76:ALA:HB3	2:B:93:HIS:O	1.71	0.91
3:E:31:DC:O2	4:F:20:DG:N2	2.05	0.90
2:D:58:ARG:NH1	5:G:7:DG:OP1	2.04	0.90
2:D:232:ARG:NH1	2:D:234:ILE:HG13	1.87	0.90



	i i i i i i i i i i i i i i i i i i i	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:C:598:ARG:NH1	1:C:604:ASP:OD2	2.04	0.90
1:A:546:SER:O	1:A:577:ARG:NH1	2.06	0.89
2:B:97:THR:OG1	2:B:99:ASN:OD1	1.89	0.89
2:B:137:ARG:NH1	2:B:178:CYS:SG	2.46	0.89
1:C:877:ARG:HH12	1:C:916:ARG:NH1	1.73	0.87
2:B:331:ILE:O	2:B:343:TYR:HA	1.76	0.86
1:C:590:GLU:OE2	1:C:714:GLY:N	2.08	0.85
5:G:9:DC:N3	6:H:53:DG:N1	2.24	0.85
1:A:413:GLN:O	1:A:426:ARG:NH1	2.10	0.85
2:B:10:ASN:HD22	2:B:11:CYS:CB	1.88	0.84
2:B:43:THR:HG22	2:B:45:ILE:H	1.43	0.84
2:D:147:SER:OG	2:D:240:LEU:HD22	1.77	0.84
1:A:617:GLU:OE2	1:A:701:ARG:NH1	2.10	0.83
1:A:624:ASP:HA	1:A:637:GLU:HB3	1.59	0.83
1:C:577:ARG:NH2	1:C:579:ASP:OD2	2.11	0.83
3:E:31:DC:N3	4:F:20:DG:N1	2.27	0.82
1:A:652:ARG:NH1	1:A:656:GLU:O	2.12	0.82
1:C:819:ASP:OD2	1:C:916:ARG:NH2	2.12	0.82
2:D:28:VAL:HG13	2:D:48:VAL:HB	1.61	0.82
1:C:590:GLU:HG2	1:C:713:VAL:HG23	1.62	0.81
1:C:575:ARG:NH1	1:C:576:PHE:O	2.14	0.81
2:B:10:ASN:ND2	2:B:10:ASN:C	2.34	0.80
1:A:486:CYS:HG	1:A:500:TYR:HH	1.28	0.80
1:C:760:SER:HB3	1:C:954:ILE:HD11	1.61	0.80
1:A:686:ASP:OD1	1:A:688:GLU:N	2.14	0.80
1:A:774:GLU:HG3	1:A:778:ARG:HH12	1.46	0.80
1:A:523:ARG:NH2	5:G:24:DG:OP2	2.14	0.79
1:A:738:GLU:OE2	1:A:775:ASN:ND2	2.16	0.79
1:A:481:LEU:O	1:A:517:GLN:NE2	2.16	0.79
1:C:538:TRP:CZ3	1:C:704:MET:CE	2.65	0.79
1:C:486:CYS:HG	1:C:500:TYR:HH	1.17	0.78
2:B:204:GLN:HB3	2:B:223:ILE:HG12	1.64	0.78
5:G:9:DC:O2	6:H:53:DG:N2	2.11	0.78
5:G:12:DT:O2	6:H:50:DA:H2	1.66	0.78
1:C:713:VAL:HG13	1:C:718:ARG:HD2	1.63	0.78
2:D:148:ARG:HH12	2:D:239:GLU:HB3	1.48	0.77
2:D:86:LYS:NZ	2:D:87:PRO:O	2.17	0.77
1:C:709:LEU:HB3	1:C:720:PHE:HB2	1.67	0.77
1:C:538:TRP:CZ3	1:C:704:MET:HE2	2.20	0.77
1:A:598:ARG:NH1	1:A:604:ASP:OD2	2.16	0.77
2:B:6:LEU:HD11	2:B:347:GLN:HB2	1.66	0.76



	la page	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:C:814:ASP:OD2	1:C:817:HIS:N	2.19	0.76
2:D:232:ARG:HE	2:D:234:ILE:HD11	1.50	0.76
3:E:28:DA:H2"	3:E:29:DG:OP2	1.86	0.76
6:H:39:DC:H2'	6:H:40:DA:C8	2.21	0.76
1:C:731:GLU:HG3	1:C:735:ARG:HD2	1.67	0.75
2:D:140:HIS:HD2	2:D:155:LEU:HD11	1.50	0.75
2:B:65:ASN:O	2:B:123:ARG:NH1	2.18	0.75
2:D:148:ARG:O	2:D:150:LYS:NZ	2.19	0.75
1:A:991:ARG:NH1	4:F:34:DG:O3'	2.19	0.74
2:D:258:ILE:HD13	2:D:284:ARG:HD2	1.67	0.74
4:F:36:DA:H2"	4:F:37:DA:OP2	1.85	0.74
1:A:858:GLN:O	1:A:862:LYS:NZ	2.16	0.74
2:B:104:SER:HB2	2:B:136:ALA:HB2	1.70	0.74
2:B:49:ARG:HH22	2:B:58:ARG:HD2	1.52	0.74
3:E:7:DG:H2"	3:E:8:DC:OP2	1.87	0.74
2:B:10:ASN:HD22	2:B:11:CYS:HB2	1.50	0.73
2:D:58:ARG:HH22	5:G:7:DG:H5"	1.53	0.73
2:D:56:LYS:HZ3	5:G:8:DC:H5"	1.54	0.73
5:G:12:DT:C4	6:H:50:DA:N6	2.54	0.73
2:D:49:ARG:NH1	2:D:58:ARG:HH21	1.87	0.73
2:B:270:GLU:OE2	2:B:270:GLU:N	2.22	0.73
5:G:12:DT:O2	6:H:50:DA:C2	2.41	0.72
2:B:148:ARG:HH12	2:B:239:GLU:HB3	1.53	0.72
2:B:49:ARG:HH22	2:B:58:ARG:HH11	1.35	0.72
2:B:283:LYS:HB2	2:B:317:TRP:HE1	1.54	0.72
1:C:617:GLU:OE2	1:C:729:TYR:OH	2.08	0.72
2:D:123:ARG:NH2	2:D:125:GLU:CD	2.41	0.72
1:C:825:GLU:OE2	1:C:950:TYR:OH	2.07	0.71
1:C:1002:LYS:HA	1:C:1005:GLU:HB3	1.72	0.71
1:C:836:GLU:HB3	1:C:839:GLN:HB2	1.72	0.71
1:A:741:GLU:N	1:A:741:GLU:OE2	2.22	0.71
1:C:738:GLU:OE2	1:C:775:ASN:ND2	2.24	0.71
1:A:688:GLU:OE2	2:B:73:ARG:NE	2.24	0.71
1:A:735:ARG:NH2	1:A:748:ILE:O	2.23	0.71
1:A:415:LEU:HD21	1:A:427:LEU:HD11	1.73	0.71
1:C:784:LYS:NZ	2:D:126:GLU:OE1	2.23	0.71
4:F:37:DA:OP2	4:F:37:DA:H8	1.74	0.71
1:A:421:ARG:NH2	6:H:14:DG:O6	2.24	0.71
1:A:493:THR:OG1	1:C:499:GLN:OE1	2.09	0.70
1:A:412:ARG:NH2	6:H:11:DT:OP1	2.24	0.70
1:A:774:GLU:HG3	1:A:778:ARG:NH1	2.07	0.70



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
5:G:11:DG:H1'	5:G:12:DT:H5'	1.73	0.70
2:B:48:VAL:HG23	2:B:55:LEU:HD11	1.74	0.69
5:G:10:DT:N3	5:G:11:DG:C5	2.60	0.69
2:B:49:ARG:HH12	2:B:58:ARG:NH1	1.90	0.69
1:A:575:ARG:NH1	1:A:576:PHE:O	2.26	0.69
5:G:10:DT:C6	5:G:11:DG:N7	2.60	0.69
4:F:22:DC:H2"	4:F:23:DT:OP2	1.92	0.69
1:A:598:ARG:HH12	1:A:604:ASP:CG	1.96	0.69
1:A:892:VAL:HG13	1:A:898:ARG:HG2	1.74	0.69
2:B:230:PRO:HB2	2:B:232:ARG:HG2	1.75	0.69
2:B:270:GLU:HB3	2:B:289:TYR:HE1	1.57	0.69
2:D:96:ARG:NH1	2:D:161:TYR:OH	2.25	0.69
5:G:12:DT:O4	6:H:50:DA:C6	2.46	0.69
3:E:18:DA:H2'	3:E:19:DC:C6	2.28	0.69
5:G:30:DC:H2"	5:G:31:DT:C7	2.23	0.69
1:C:641:ARG:NH1	1:C:987:ASN:OD1	2.24	0.68
4:F:12:DT:C6	4:F:13:DT:H72	2.29	0.68
1:C:861:LYS:HE3	1:C:862:LYS:HG3	1.75	0.68
4:F:42:DG:OP2	4:F:42:DG:H2'	1.93	0.68
2:D:338:THR:HG22	2:D:340:PRO:CA	2.18	0.68
1:A:740:LEU:HD11	1:A:770:ARG:NH1	2.04	0.68
5:G:49:DA:H2	6:H:13:DT:H3	1.38	0.68
1:A:749:CYS:SG	1:A:959:HIS:NE2	2.66	0.68
2:B:277:TYR:HA	2:B:284:ARG:H	1.58	0.68
1:C:988:LYS:O	1:C:991:ARG:HB2	1.93	0.68
5:G:31:DT:O2	6:H:32:DG:N2	2.26	0.68
1:C:418:LEU:HD11	1:C:426:ARG:NH1	2.10	0.67
2:D:284:ARG:NE	2:D:286:GLU:OE2	2.28	0.67
5:G:10:DT:C2	5:G:11:DG:C5	2.82	0.67
1:A:872:ASN:ND2	3:E:18:DA:OP1	2.27	0.67
2:B:108:MET:HE3	2:B:125:GLU:HB2	1.76	0.67
2:D:56:LYS:NZ	5:G:8:DC:H5"	2.09	0.67
1:A:771:SER:N	1:A:774:GLU:OE2	2.27	0.67
1:A:885:VAL:HG21	1:A:905:MET:HG2	1.77	0.67
2:B:28:VAL:HG13	2:B:48:VAL:HG13	1.76	0.67
3:E:41:DA:H8	3:E:41:DA:OP2	1.76	0.67
2:D:232:ARG:NE	2:D:234:ILE:HD11	2.10	0.67
1:C:771:SER:O	1:C:775:ASN:ND2	2.28	0.67
1:A:426:ARG:NH2	1:C:442:GLU:OE1	2.28	0.66
5:G:10:DT:C4	5:G:11:DG:O6	2.48	0.66
2:D:82:ALA:HB2	2:D:88:GLU:HB2	1.77	0.66



	l as page	Interatomic	Clash
Atom-1	Atom-2	distance $(Å)$	overlap (Å)
5:G:10:DT:H2"	5:G:11:DG:H8	1.61	0.66
2:D:210:LEU:HD13	2:D:292:LEU:HD21	1.78	0.66
2:B:62:PHE:HE1	2:B:122:LEU:HB2	1.59	0.65
2:D:168:THR:OG1	2:D:171:ASN:HB2	1.96	0.65
3:E:27:DC:OP2	3:E:27:DC:H2'	1.96	0.65
1:A:566:ASP:OD2	2:B:138:TYR:OH	2.14	0.65
1:A:861:LYS:NZ	1:A:862:LYS:NZ	2.44	0.65
1:A:755:THR:N	1:A:758:GLU:OE2	2.21	0.65
2:D:290:VAL:HG13	2:D:297:VAL:HG23	1.77	0.65
4:F:31:DT:H2"	4:F:32:DG:OP2	1.96	0.65
3:E:21:DG:H2'	3:E:22:DT:H71	1.77	0.65
1:A:506:THR:HG21	1:C:489:ILE:HG12	1.79	0.65
1:A:622:MET:HG2	1:A:625:VAL:HG12	1.77	0.65
1:A:664:GLN:O	1:A:666:GLN:NE2	2.30	0.65
2:B:212:ARG:NH1	2:B:213:GLN:OE1	2.30	0.65
1:C:418:LEU:HB2	1:C:423:GLN:HG3	1.78	0.65
1:C:735:ARG:NH1	1:C:743:SER:HA	2.12	0.65
4:F:36:DA:H8	4:F:36:DA:OP2	1.80	0.65
1:C:735:ARG:NH1	1:C:742:ALA:HA	2.12	0.65
3:E:18:DA:H2'	3:E:19:DC:H6	1.62	0.65
1:A:618:SER:OG	1:A:728:GLY:O	2.14	0.64
2:D:80:PHE:HD2	2:D:89:CYS:HB2	1.60	0.64
1:A:641:ARG:NH1	1:A:987:ASN:OD1	2.30	0.64
2:D:249:CYS:SG	2:D:250:THR:N	2.70	0.64
5:G:30:DC:C2'	5:G:31:DT:H71	2.24	0.64
1:A:590:GLU:HB2	1:A:713:VAL:HG23	1.79	0.64
1:A:988:LYS:HZ2	3:E:20:DA:H3'	1.61	0.64
1:A:682:VAL:HG12	1:A:683:ASP:H	1.61	0.64
1:A:861:LYS:HZ2	1:A:862:LYS:HZ3	1.44	0.64
4:F:32:DG:H2"	4:F:33:DT:OP2	1.97	0.64
1:A:671:GLU:OE1	1:A:671:GLU:N	2.30	0.64
1:A:620:ASP:OD1	1:A:621:GLY:N	2.31	0.64
1:C:832:ASP:HB3	1:C:837:VAL:HG11	1.79	0.64
1:A:482:HIS:HB3	1:A:485:VAL:HG23	1.80	0.63
1:A:553:ILE:HG12	1:A:576:PHE:HE1	1.62	0.63
2:B:10:ASN:ND2	2:B:10:ASN:O	2.30	0.63
5:G:10:DT:C4	5:G:11:DG:C5	2.86	0.63
1:A:623:GLY:HA2	1:A:638:LYS:HD3	1.81	0.63
1:A:817:HIS:O	1:A:819:ASP:N	2.32	0.63
2:B:229:ARG:NH2	2:B:280:GLU:OE1	2.27	0.63
2:B:277:TYR:OH	2:B:316:THR:OG1	2.14	0.62



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
2:B:316:THR:HG23	2:B:332:PRO:HG3	1.80	0.62
1:C:588:ASP:OD2	1:C:1018:SER:OG	2.18	0.62
2:B:10:ASN:HD22	2:B:10:ASN:C	2.02	0.62
4:F:7:DG:H2'	4:F:8:DG:H8	1.65	0.62
1:C:609:SER:HB3	1:C:653:LEU:HD11	1.81	0.62
5:G:10:DT:N1	5:G:11:DG:C8	2.68	0.62
2:B:199:GLU:OE2	2:B:199:GLU:N	2.33	0.62
2:D:328:LEU:HD11	2:D:345:PHE:HD2	1.63	0.62
1:A:496:SER:OG	1:A:499:GLN:N	2.29	0.61
1:C:741:GLU:N	1:C:741:GLU:OE1	2.32	0.61
2:D:1:MET:SD	2:D:302:ARG:NH1	2.73	0.61
2:D:49:ARG:NH1	2:D:58:ARG:HD2	2.14	0.61
5:G:32:DC:H2"	5:G:33:DC:H5	1.64	0.61
5:G:9:DC:N4	6:H:53:DG:O6	2.26	0.61
2:B:333:SER:OG	2:B:344:HIS:ND1	2.01	0.61
2:D:140:HIS:CD2	2:D:155:LEU:HD11	2.32	0.61
1:A:531:PRO:HG3	1:A:576:PHE:CE2	2.35	0.61
2:B:148:ARG:HH11	2:B:241:LEU:HD13	1.65	0.61
3:E:35:DA:H2	4:F:17:DC:H42	1.49	0.61
1:A:578:TYR:HH	1:A:724:PHE:HE2	1.46	0.61
1:A:754:SER:HB2	1:A:758:GLU:HB2	1.81	0.61
1:A:818:CYS:SG	1:A:915:TRP:NE1	2.72	0.61
2:B:217:TYR:HB3	2:B:233:LEU:HD21	1.82	0.61
1:C:795:ARG:NH1	2:D:39:ARG:NH1	2.49	0.61
1:C:840:LYS:NZ	1:C:843:PRO:HA	2.16	0.61
2:B:1:MET:HA	2:B:349:SER:O	2.01	0.61
4:F:28:DC:H2"	4:F:29:DA:C8	2.36	0.61
1:A:861:LYS:NZ	1:A:862:LYS:HZ3	1.99	0.61
2:B:49:ARG:HH22	2:B:58:ARG:NH1	1.99	0.61
1:C:871:MET:HG3	1:C:875:TYR:HB3	1.81	0.61
5:G:32:DC:H2"	5:G:33:DC:OP2	2.00	0.61
5:G:36:DT:H3	6:H:26:DA:H61	1.47	0.61
1:A:613:VAL:HG23	1:A:649:ILE:HD13	1.82	0.60
1:C:538:TRP:CE3	1:C:704:MET:HE1	2.36	0.60
6:H:26:DA:H2"	6:H:27:DG:H5'	1.82	0.60
3:E:26:DA:H2"	3:E:27:DC:OP2	2.01	0.60
2:B:7:THR:OG1	2:B:54:GLU:OE1	2.20	0.60
1:C:594:MET:SD	1:C:718:ARG:NH1	2.67	0.60
1:C:870:ARG:HD3	5:G:17:DC:N4	2.17	0.60
1:A:760:SER:OG	1:A:950:TYR:O	2.19	0.60
1:A:420:ARG:NE	5:G:43:DC:OP1	2.24	0.60



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:740:LEU:CD1	1:A:770:ARG:HH12	2.07	0.60
4:F:29:DA:H8	4:F:29:DA:OP2	1.84	0.60
1:C:622:MET:SD	1:C:987:ASN:ND2	2.74	0.60
2:B:10:ASN:ND2	2:B:11:CYS:SG	2.75	0.60
1:C:748:ILE:HG21	1:C:756:ARG:HE	1.66	0.60
2:D:28:VAL:CG1	2:D:48:VAL:HB	2.32	0.60
2:D:257:THR:HG23	2:D:284:ARG:HH22	1.66	0.59
1:A:675:ARG:NE	1:A:1017:THR:O	2.36	0.59
1:C:795:ARG:HH12	2:D:39:ARG:NH1	2.00	0.59
1:C:538:TRP:CZ3	1:C:704:MET:HE1	2.36	0.59
2:B:82:ALA:HB2	2:B:88:GLU:HB3	1.84	0.59
2:B:269:HIS:N	2:B:270:GLU:OE2	2.36	0.59
2:D:8:ALA:HA	2:D:55:LEU:HB3	1.84	0.59
2:D:232:ARG:NH1	2:D:234:ILE:HG12	2.10	0.59
3:E:14:DT:H2'	3:E:14:DT:OP2	2.01	0.59
1:A:915:TRP:HB2	1:A:975:ILE:HD12	1.83	0.59
2:B:2:SER:O	2:B:348:VAL:HA	2.02	0.59
1:C:951:ASP:OD1	1:C:951:ASP:N	2.35	0.59
1:C:652:ARG:NH2	1:C:657:ASP:O	2.36	0.59
2:B:62:PHE:CE1	2:B:122:LEU:HB2	2.38	0.59
3:E:6:DG:H8	3:E:6:DG:OP2	1.86	0.59
1:A:753:ASP:OD1	1:A:753:ASP:N	2.36	0.59
1:A:412:ARG:HE	6:H:11:DT:H5"	1.66	0.59
2:B:239:GLU:HB2	2:B:246:VAL:HG13	1.85	0.59
2:D:33:GLN:NE2	2:D:37:PRO:HA	2.17	0.59
1:A:771:SER:O	1:A:775:ASN:ND2	2.36	0.58
2:D:134:PRO:HG3	2:D:155:LEU:HD23	1.85	0.58
4:F:21:DT:H2"	4:F:22:DC:OP2	2.02	0.58
1:A:627:GLU:HG3	1:A:994:ARG:NH1	2.18	0.58
2:B:10:ASN:ND2	2:B:11:CYS:CB	2.65	0.58
1:C:814:ASP:OD1	1:C:815:ALA:N	2.36	0.58
2:D:79:HIS:NE2	2:D:88:GLU:OE2	2.37	0.58
3:E:21:DG:H4'	3:E:22:DT:OP1	2.03	0.58
5:G:41:DG:H2"	5:G:42:DG:C8	2.39	0.58
2:B:10:ASN:ND2	2:B:11:CYS:HB2	2.17	0.58
1:A:442:GLU:OE1	1:C:426:ARG:NH2	2.36	0.58
3:E:29:DG:OP2	3:E:29:DG:H2'	2.03	0.58
4:F:45:DC:H2"	4:F:46:DA:OP2	2.03	0.58
1:A:574:ARG:NH2	1:A:1003:THR:O	2.37	0.58
2:D:226:SER:HG	2:D:228:CYS:HG	1.44	0.58
1:A:420:ARG:NH1	1:C:459:ARG:HD3	2.19	0.57



		Interatomic	Clash
Atom-1	Atom-2	distance $(Å)$	overlap (Å)
2:B:274:PHE:CE2	2:B:348:VAL:HG21	2.39	0.57
6:H:28:DT:H2'	6:H:28:DT:OP2	2.04	0.57
2:B:168:THR:HG22	2:B:170:GLN:H	1.69	0.57
2:D:306:GLN:OE1	2:D:306:GLN:N	2.31	0.57
2:D:322:LEU:H	2:D:327:ALA:HA	1.68	0.57
3:E:22:DT:H2"	3:E:23:DG:C8	2.39	0.57
6:H:39:DC:H2'	6:H:40:DA:H8	1.66	0.57
1:A:435:LYS:HD2	1:A:446:VAL:HG21	1.86	0.57
2:D:307:TRP:HB3	2:D:311:ILE:HG23	1.84	0.57
4:F:27:DG:C6	4:F:28:DC:N4	2.73	0.57
2:B:73:ARG:HH12	2:B:96:ARG:NH1	2.03	0.57
5:G:10:DT:C2	5:G:11:DG:C8	2.92	0.57
1:A:986:GLY:O	1:A:989:LEU:HB2	2.04	0.57
2:D:49:ARG:HH12	2:D:58:ARG:HH21	1.51	0.57
2:D:310:GLU:OE1	2:D:346:TYR:OH	2.17	0.57
5:G:10:DT:C1'	5:G:11:DG:C8	2.88	0.57
1:A:817:HIS:C	1:A:819:ASP:H	2.07	0.57
1:A:903:LYS:HD3	1:A:941:LEU:HD22	1.86	0.57
2:B:148:ARG:NH1	2:B:241:LEU:HD13	2.19	0.57
2:D:6:LEU:HD11	2:D:347:GLN:HB2	1.87	0.57
2:D:152:ALA:HB2	2:D:240:LEU:HD21	1.86	0.57
3:E:37:DC:H1'	4:F:14:DG:H22	1.69	0.57
5:G:35:DC:H2'	5:G:36:DT:O4'	2.05	0.57
2:B:289:TYR:HD2	2:B:300:GLU:HB3	1.70	0.57
4:F:23:DT:OP2	4:F:23:DT:H2'	2.05	0.57
2:D:147:SER:OG	2:D:240:LEU:HD13	2.05	0.56
1:A:538:TRP:CH2	1:A:579:ASP:HB2	2.39	0.56
1:A:590:GLU:HB2	1:A:713:VAL:CG2	2.35	0.56
1:A:869:MET:HE3	1:A:870:ARG:HG2	1.86	0.56
1:C:557:LEU:O	2:D:173:ASN:ND2	2.25	0.56
1:A:763:MET:SD	1:A:943:SER:OG	2.61	0.56
2:B:212:ARG:HG3	2:B:269:HIS:NE2	2.20	0.56
2:D:71:PRO:HG2	2:D:98:PRO:HG3	1.86	0.56
3:E:37:DC:H2"	3:E:38:DA:C8	2.41	0.56
1:C:619:CYS:HB2	1:C:642:PHE:CD2	2.41	0.56
1:A:470:LEU:HD13	1:C:470:LEU:HD13	1.86	0.56
1:A:530:LEU:HD22	1:A:531:PRO:HD2	1.86	0.56
2:B:333:SER:CB	2:B:344:HIS:ND1	2.68	0.56
2:B:235:ARG:HB3	2:B:250:THR:HG23	1.87	0.56
2:B:338:THR:HG22	2:B:340:PRO:HA	1.87	0.56
1:C:558:SER:OG	2:D:172:TRP:N	2.39	0.56



Atom 1	Atom 2	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
5:G:51:DA:H2"	5:G:52:DA:OP2	2.06	0.56
6:H:34:DA:H2"	6:H:35:DC:OP2	2.06	0.56
2:D:331:ILE:O	2:D:343:TYR:HA	2.06	0.55
5:G:52:DA:OP2	5:G:52:DA:H2'	2.06	0.55
1:C:566:ASP:CG	2:D:159:ARG:HH12	2.09	0.55
5:G:30:DC:C2'	5:G:31:DT:C7	2.84	0.55
2:B:10:ASN:HD22	2:B:11:CYS:N	2.05	0.55
1:C:903:LYS:HZ3	1:C:907:LEU:HG	1.71	0.55
5:G:16:DA:H2"	5:G:17:DC:H5'	1.88	0.55
1:A:817:HIS:ND1	1:A:957:TYR:OH	2.35	0.55
1:C:779:TYR:CZ	1:C:804:LYS:HB2	2.42	0.55
2:D:307:TRP:HB2	2:D:312:SER:HB2	1.87	0.55
3:E:20:DA:H2"	3:E:21:DG:C8	2.41	0.55
1:A:505:ARG:HH12	1:C:1029:LYS:C	2.02	0.55
2:B:212:ARG:HG3	2:B:269:HIS:CE1	2.41	0.55
5:G:32:DC:C2	5:G:33:DC:N4	2.75	0.55
1:A:492:ASN:O	1:A:492:ASN:ND2	2.32	0.55
1:A:536:PHE:HB3	1:A:711:ILE:HD13	1.89	0.55
1:C:536:PHE:CE2	1:C:549:TRP:HB2	2.42	0.55
2:D:148:ARG:N	2:D:149:GLY:HA2	2.22	0.55
1:A:872:ASN:OD1	1:A:875:TYR:N	2.26	0.55
1:C:409:GLY:HA2	4:F:9:DT:H1'	1.87	0.55
1:C:675:ARG:NH2	1:C:1018:SER:HA	2.21	0.55
2:D:39:ARG:NH2	6:H:51:DC:OP2	2.40	0.55
5:G:10:DT:N3	5:G:11:DG:C6	2.75	0.55
1:A:912:LYS:HA	1:A:915:TRP:CZ3	2.42	0.55
5:G:44:DT:N3	5:G:45:DG:O6	2.39	0.55
1:A:553:ILE:HG12	1:A:576:PHE:CE1	2.40	0.55
1:A:869:MET:CE	1:A:870:ARG:HG2	2.36	0.54
1:A:487:LEU:HD22	1:A:526:GLU:HB2	1.88	0.54
1:C:708:ARG:HG2	1:C:721:ARG:HG2	1.88	0.54
1:C:735:ARG:HG2	1:C:740:LEU:HD12	1.88	0.54
2:D:277:TYR:HA	2:D:283:LYS:HA	1.87	0.54
5:G:10:DT:O4	5:G:11:DG:O6	2.25	0.54
6:H:35:DC:OP2	6:H:35:DC:H2'	2.08	0.54
1:A:731:GLU:N	1:A:731:GLU:OE1	2.40	0.54
1:A:748:ILE:HD13	1:A:756:ARG:HD2	1.89	0.54
1:C:735:ARG:HB3	1:C:740:LEU:HB2	1.89	0.54
2:D:1:MET:N	2:D:349:SER:O	2.37	0.54
1:A:883:GLU:OE1	1:A:883:GLU:N	2.33	0.54
1:A:959:HIS:HE2	1:A:964:HIS:HE1	1.56	0.54



	A de page	Interatomic	Clash
Atom-1	Atom-2	distance (\AA)	overlap (Å)
1:A:871:MET:SD	1:A:871:MET:N	2.80	0.54
2:D:284:ARG:HB2	2:D:284:ARG:NH1	2.22	0.54
3:E:31:DC:C2	4:F:20:DG:N2	2.72	0.54
4:F:20:DG:H2'	4:F:21:DT:C6	2.42	0.54
1:A:429:ASP:HB2	1:C:437:PHE:HE2	1.73	0.54
1:A:640:VAL:HB	1:A:680:MET:HE3	1.90	0.54
2:D:229:ARG:HE	2:D:259:THR:HG21	1.72	0.54
5:G:33:DC:H2'	5:G:34:DA:C4	2.42	0.54
1:A:956:ASN:HD21	1:A:960:LYS:NZ	2.05	0.54
2:D:6:LEU:HD22	2:D:53:GLY:HA2	1.88	0.54
2:D:159:ARG:HA	2:D:178:CYS:H	1.73	0.54
1:A:687:HIS:O	1:A:691:THR:OG1	2.22	0.54
1:C:628:LYS:HG2	1:C:994:ARG:NH1	2.23	0.54
5:G:15:DT:H2'	5:G:16:DA:H8	1.73	0.54
1:C:865:LEU:HD21	1:C:878:ARG:HG2	1.89	0.54
1:A:840:LYS:NZ	1:A:847:GLU:OE2	2.32	0.53
2:B:339:PRO:HB2	2:B:341:GLU:H	1.73	0.53
1:C:675:ARG:NH1	1:C:675:ARG:HG3	2.23	0.53
3:E:21:DG:H2"	3:E:22:DT:O5'	2.08	0.53
5:G:10:DT:C2'	5:G:11:DG:C8	2.91	0.53
5:G:32:DC:H1'	5:G:33:DC:C5	2.43	0.53
1:A:452:THR:O	1:A:456:LEU:HG	2.08	0.53
1:A:779:TYR:CZ	1:A:804:LYS:HB2	2.43	0.53
2:B:270:GLU:HB3	2:B:289:TYR:CE1	2.41	0.53
1:A:753:ASP:HB2	1:A:799:LYS:HG2	1.91	0.53
1:A:825:GLU:OE2	1:A:950:TYR:OH	2.24	0.53
2:D:58:ARG:NH2	5:G:7:DG:H5"	2.21	0.53
4:F:27:DG:H2'	4:F:28:DC:C6	2.43	0.53
6:H:33:DT:OP2	6:H:33:DT:H2'	2.08	0.53
2:B:108:MET:HE2	2:B:127:LYS:HD2	1.91	0.53
2:D:212:ARG:HG3	2:D:269:HIS:NE2	2.22	0.53
5:G:10:DT:N1	5:G:11:DG:N7	2.56	0.53
6:H:15:DT:H2"	6:H:16:DA:H5"	1.90	0.53
1:C:984:GLU:O	1:C:987:ASN:HB2	2.09	0.53
5:G:31:DT:C2	6:H:32:DG:N2	2.76	0.53
1:C:788:SER:HB2	2:D:65:ASN:HA	1.90	0.53
5:G:28:DT:H2"	5:G:29:DA:N7	2.24	0.53
1:A:534:HIS:CE1	1:A:587:LYS:HD2	2.43	0.53
1:A:922:ARG:HB3	1:A:922:ARG:NH1	2.24	0.53
1:C:534:HIS:CG	1:C:587:LYS:HZ2	2.27	0.53
1:C:885:VAL:HG21	1:C:905:MET:HG2	1.90	0.53



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:C:942:LEU:HD22	1:C:946:PHE:HB2	1.91	0.53
2:D:123:ARG:NH2	2:D:125:GLU:CG	2.71	0.53
4:F:7:DG:H2'	4:F:8:DG:C8	2.43	0.53
1:A:543:LYS:O	1:A:545:VAL:HG23	2.09	0.53
1:C:578:TYR:HB2	1:C:700:GLU:OE2	2.09	0.53
1:C:608:THR:HB	1:C:718:ARG:HG2	1.91	0.53
1:A:539:GLN:NE2	1:A:710:ILE:HG13	2.24	0.53
1:A:730:ASP:N	1:A:730:ASP:OD1	2.41	0.53
1:C:523:ARG:HH22	3:E:23:DG:H2'	1.74	0.53
2:B:49:ARG:NH2	2:B:58:ARG:HD2	2.21	0.52
1:C:410:ARG:HD2	3:E:41:DA:H2	1.74	0.52
2:D:188:GLU:OE2	2:D:189:PHE:HA	2.09	0.52
1:A:578:TYR:HA	1:A:677:LEU:HD21	1.91	0.52
2:D:212:ARG:NH1	2:D:269:HIS:HD2	2.07	0.52
5:G:10:DT:C2	5:G:11:DG:C4	2.98	0.52
1:C:773:ASP:OD1	1:C:773:ASP:N	2.40	0.52
1:A:613:VAL:HG22	1:A:649:ILE:HB	1.91	0.52
2:B:299:MET:SD	2:B:300:GLU:N	2.83	0.52
1:C:641:ARG:HH11	1:C:987:ASN:HA	1.73	0.52
3:E:31:DC:OP2	3:E:31:DC:H6	1.93	0.52
5:G:48:DC:H2"	5:G:49:DA:C8	2.44	0.52
2:D:266:ILE:HD13	2:D:350:PHE:CD2	2.44	0.52
5:G:34:DA:H8	5:G:34:DA:OP2	1.93	0.52
2:B:256:LEU:HD12	2:B:257:THR:H	1.74	0.52
1:C:999:ARG:N	1:C:1000:GLN:HA	2.25	0.52
6:H:12:DT:H2"	6:H:13:DT:H5'	1.92	0.52
1:A:470:LEU:HB2	1:C:473:MET:HE1	1.91	0.52
1:A:566:ASP:OD1	1:A:567:VAL:N	2.39	0.52
1:C:675:ARG:HG3	1:C:675:ARG:HH11	1.74	0.52
1:C:991:ARG:O	1:C:994:ARG:HB3	2.10	0.52
1:A:696:PRO:HG3	2:B:172:TRP:HB3	1.91	0.52
1:A:828:LYS:HB3	1:A:949:ARG:NH2	2.25	0.52
2:B:277:TYR:OH	2:B:317:TRP:O	2.14	0.52
1:C:563:SER:OG	1:C:566:ASP:OD2	2.26	0.52
2:D:60:ILE:HG22	2:D:61:SER:H	1.75	0.52
2:D:73:ARG:NH2	2:D:100:ASN:OD1	2.43	0.51
2:D:217:TYR:HD2	2:D:233:LEU:HD21	1.75	0.51
1:A:870:ARG:HH22	1:A:872:ASN:HA	1.74	0.51
2:B:148:ARG:N	2:B:149:GLY:HA2	2.25	0.51
1:C:889:CYS:O	1:C:898:ARG:NE	2.43	0.51
2:D:311:ILE:HB	2:D:346:TYR:HE2	1.75	0.51



	las page	Interatomic	Clash
Atom-1	Atom-2	distance $(Å)$	overlap (Å)
5:G:27:DG:H2"	5:G:28:DT:OP2	2.10	0.51
5:G:28:DT:OP2	5:G:28:DT:H2'	2.09	0.51
1:A:832:ASP:OD2	1:A:949:ARG:NH1	2.42	0.51
1:A:865:LEU:HD11	1:A:878:ARG:HB3	1.92	0.51
1:A:959:HIS:O	1:A:963:ALA:CB	2.59	0.51
2:D:141:THR:HG21	2:D:209:ALA:HB3	1.91	0.51
3:E:41:DA:C8	3:E:41:DA:H5'	2.45	0.51
2:B:21:LEU:HD12	2:B:320:GLY:HA3	1.92	0.51
5:G:52:DA:H2"	5:G:53:DA:C8	2.46	0.51
1:C:752:CYS:HG	1:C:766:HIS:CG	2.28	0.51
3:E:31:DC:N4	4:F:20:DG:O6	2.43	0.51
1:A:641:ARG:HH11	1:A:987:ASN:HA	1.76	0.51
2:B:249:CYS:SG	2:B:250:THR:N	2.84	0.51
1:C:650:SER:HA	1:C:662:ILE:HG13	1.93	0.51
1:C:709:LEU:O	1:C:719:SER:HA	2.10	0.51
1:C:863:MET:HE1	1:C:883:GLU:OE2	2.11	0.51
2:B:289:TYR:HB3	2:B:300:GLU:OE1	2.11	0.51
2:D:145:ILE:HB	2:D:214:ASP:HA	1.92	0.51
2:B:23:ASP:OD1	2:B:24:LEU:N	2.43	0.51
2:B:218:PHE:HB2	2:B:234:ILE:HB	1.91	0.51
1:C:926:ASP:OD1	1:C:926:ASP:N	2.39	0.51
2:D:78:ALA:HB1	2:D:144:VAL:HG23	1.93	0.51
4:F:36:DA:OP2	4:F:36:DA:C8	2.61	0.51
6:H:55:DC:H2"	6:H:56:DC:C6	2.46	0.51
1:A:479:PHE:HA	1:C:513:ARG:NH2	2.26	0.51
1:A:622:MET:HE1	1:A:990:PHE:HB3	1.93	0.51
1:A:959:HIS:NE2	1:A:964:HIS:HE1	2.09	0.51
2:B:8:ALA:HB2	2:B:55:LEU:HD23	1.92	0.51
3:E:9:DC:H2"	3:E:10:DT:H71	1.92	0.51
1:A:506:THR:O	1:A:510:THR:OG1	2.23	0.51
2:B:94:GLY:H	2:B:140:HIS:CE1	2.29	0.51
5:G:55:DC:H4'	5:G:56:DC:OP1	2.11	0.51
1:A:908:TYR:O	1:A:911:MET:HG2	2.11	0.50
1:C:415:LEU:HD21	1:C:427:LEU:HD11	1.92	0.50
2:D:229:ARG:HH22	2:D:280:GLU:HA	1.76	0.50
3:E:17:DC:H2"	3:E:18:DA:C8	2.46	0.50
3:E:20:DA:H2'	3:E:20:DA:OP2	2.11	0.50
5:G:13:DC:C5	5:G:14:DT:H73	2.46	0.50
1:A:650:SER:OG	1:A:651:ILE:N	2.44	0.50
2:B:147:SER:HB3	2:B:240:LEU:H	1.76	0.50
2:B:300:GLU:OE2	2:B:302:ARG:HD3	2.11	0.50



	At arra 2	Interatomic	Clash
Atom-1	Atom-2	distance (\AA)	overlap (Å)
1:C:641:ARG:NH1	1:C:987:ASN:HA	2.26	0.50
2:D:229:ARG:HH12	2:D:280:GLU:HG2	1.75	0.50
4:F:36:DA:OP2	4:F:36:DA:H2'	2.12	0.50
2:B:49:ARG:NH2	2:B:58:ARG:HH11	2.06	0.50
2:B:49:ARG:NH1	2:B:58:ARG:NH1	2.59	0.50
2:D:135:SER:N	2:D:137:ARG:HH21	2.09	0.50
3:E:8:DC:OP2	3:E:8:DC:H2'	2.12	0.50
6:H:32:DG:H2"	6:H:33:DT:OP2	2.11	0.50
1:C:534:HIS:CG	1:C:587:LYS:NZ	2.80	0.50
2:D:43:THR:HG23	2:D:62:PHE:CE1	2.47	0.50
2:D:202:ASP:OD2	2:D:223:ILE:HD13	2.10	0.50
1:C:540:PRO:HD2	1:C:707:SER:HA	1.94	0.50
1:A:849:ARG:HH11	1:A:849:ARG:HB2	1.77	0.50
1:C:778:ARG:NH2	1:C:799:LYS:HB2	2.27	0.50
2:D:269:HIS:HB3	2:D:292:LEU:HB2	1.94	0.50
1:A:861:LYS:HZ1	1:A:862:LYS:NZ	2.09	0.50
1:C:452:THR:HA	1:C:455:LEU:HD12	1.94	0.50
3:E:21:DG:H2'	3:E:22:DT:C6	2.46	0.50
1:A:859:LEU:O	1:A:864:LYS:N	2.44	0.49
2:B:140:HIS:HB3	2:B:158:GLY:HA3	1.94	0.49
1:C:896:GLU:OE1	1:C:896:GLU:N	2.36	0.49
1:C:992:ARG:HB2	5:G:21:DG:OP1	2.12	0.49
1:C:860:ARG:HG2	1:C:866:LYS:HE3	1.94	0.49
6:H:1:DC:H3'	6:H:2:DT:H71	1.92	0.49
1:A:788:SER:HB2	2:B:65:ASN:HA	1.94	0.49
1:A:1013:HIS:O	1:A:1017:THR:HG22	2.12	0.49
2:D:49:ARG:CZ	2:D:58:ARG:HD2	2.42	0.49
3:E:29:DG:H2"	3:E:30:DA:OP2	2.12	0.49
3:E:44:DC:H4'	3:E:45:DC:OP1	2.11	0.49
1:C:515:ILE:O	1:C:517:GLN:NE2	2.44	0.49
1:C:877:ARG:NH1	1:C:916:ARG:NH1	2.51	0.49
1:C:538:TRP:CZ3	1:C:709:LEU:HB2	2.48	0.49
1:C:641:ARG:HH11	1:C:987:ASN:CG	2.14	0.49
4:F:41:DA:H2"	4:F:42:DG:OP2	2.12	0.49
1:A:418:LEU:HB3	1:A:422:ALA:HB3	1.95	0.49
2:B:86:LYS:NZ	2:B:123:ARG:HD3	2.28	0.49
2:B:164:PRO:HA	2:B:167:ARG:HB2	1.94	0.49
1:C:737:MET:HE3	1:C:807:MET:HB3	1.93	0.49
1:C:790:SER:OG	1:C:791:ALA:N	2.45	0.49
1:A:869:MET:HB3	3:E:16:DA:H4'	1.95	0.49
1:A:999:ARG:NH1	1:A:1004:PHE:CG	2.81	0.49



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
6:H:29:DG:H2"	6:H:30:DG:OP2	2.12	0.49
1:A:493:THR:O	1:C:499:GLN:NE2	2.41	0.49
1:A:988:LYS:NZ	3:E:20:DA:P	2.85	0.49
1:C:740:LEU:HD21	1:C:770:ARG:HH12	1.78	0.49
2:D:130:VAL:HB	2:D:191:CYS:HA	1.93	0.49
2:D:222:HIS:HB2	2:D:259:THR:HG22	1.94	0.49
6:H:30:DG:H2"	6:H:31:DA:C8	2.48	0.49
4:F:40:DC:H1'	4:F:41:DA:C8	2.48	0.49
1:A:687:HIS:CD2	2:B:36:TRP:HE1	2.31	0.49
5:G:5:DT:H2'	5:G:5:DT:OP2	2.13	0.49
1:A:538:TRP:CD2	1:A:542:LEU:HD21	2.46	0.48
1:A:590:GLU:O	1:A:593:ILE:HG12	2.13	0.48
1:A:926:ASP:OD1	1:A:926:ASP:N	2.39	0.48
1:C:619:CYS:SG	1:C:620:ASP:N	2.85	0.48
2:D:47:GLY:N	2:D:60:ILE:HD11	2.27	0.48
2:D:212:ARG:NH1	2:D:293:ASP:HA	2.27	0.48
2:D:235:ARG:HB3	2:D:250:THR:HG23	1.94	0.48
4:F:10:DT:C6	4:F:11:DT:H72	2.48	0.48
6:H:35:DC:H2"	6:H:36:DT:C6	2.48	0.48
1:A:737:MET:HG2	1:A:806:PHE:CE1	2.48	0.48
1:C:415:LEU:O	1:C:423:GLN:NE2	2.47	0.48
1:C:555:ASP:N	1:C:555:ASP:OD1	2.46	0.48
1:C:638:LYS:HD2	1:C:683:ASP:OD2	2.13	0.48
1:C:749:CYS:HB3	1:C:752:CYS:O	2.13	0.48
1:A:429:ASP:HB2	1:C:437:PHE:CE2	2.47	0.48
1:A:911:MET:HB2	1:A:931:TYR:CE2	2.49	0.48
2:B:89:CYS:SG	2:B:110:SER:HB2	2.53	0.48
2:B:226:SER:OG	2:B:228:CYS:SG	2.56	0.48
2:D:80:PHE:CE1	2:D:151:THR:HG21	2.49	0.48
2:D:265:PRO:HB3	2:D:271:TYR:CZ	2.49	0.48
4:F:27:DG:H2"	4:F:28:DC:H5'	1.96	0.48
5:G:10:DT:H2"	5:G:11:DG:C8	2.45	0.48
1:A:420:ARG:HH11	1:C:459:ARG:HD3	1.78	0.48
5:G:9:DC:H2"	5:G:10:DT:OP2	2.12	0.48
1:A:968:ILE:HG23	1:A:971:ARG:HD3	1.96	0.48
4:F:32:DG:OP2	4:F:32:DG:H8	1.94	0.48
5:G:32:DC:C2'	5:G:33:DC:H5	2.27	0.48
1:A:858:GLN:NE2	1:A:890:GLU:OE1	2.47	0.48
3:E:45:DC:H42	4:F:6:DG:H22	1.61	0.48
1:A:410:ARG:HB2	5:G:52:DA:H2	1.78	0.48
1:A:694:LEU:O	1:A:698:VAL:HG23	2.13	0.48



		Interatomic	Clash
Atom-1	Atom-2	distance $(Å)$	overlap (Å)
1:A:771:SER:HB3	1:A:774:GLU:OE2	2.14	0.48
1:C:796:ASP:O	1:C:799:LYS:HD2	2.12	0.48
2:B:93:HIS:HA	2:B:94:GLY:HA2	1.62	0.48
4:F:37:DA:OP2	4:F:37:DA:C8	2.62	0.48
5:G:32:DC:C1'	5:G:33:DC:C5	2.97	0.48
1:A:490:ARG:O	1:A:493:THR:HG22	2.14	0.47
1:A:829:ILE:O	1:A:833:GLU:HB2	2.14	0.47
1:A:839:GLN:N	1:A:839:GLN:OE1	2.46	0.47
1:A:849:ARG:HB2	1:A:849:ARG:NH1	2.29	0.47
2:B:16:GLN:HB3	2:B:34:LYS:HB3	1.95	0.47
2:B:229:ARG:O	2:B:257:THR:OG1	2.20	0.47
2:B:316:THR:OG1	2:B:317:TRP:N	2.47	0.47
1:C:889:CYS:HB2	1:C:898:ARG:HG3	1.95	0.47
2:D:113:SER:OG	2:D:114:ARG:N	2.47	0.47
5:G:10:DT:O4	5:G:11:DG:C6	2.67	0.47
1:A:589:LEU:O	1:A:593:ILE:HG23	2.14	0.47
1:A:737:MET:HG2	1:A:806:PHE:HE1	1.79	0.47
2:B:5:PRO:HA	2:B:346:TYR:HD1	1.79	0.47
2:B:87:PRO:HA	2:B:88:GLU:HA	1.50	0.47
1:C:845:ARG:O	1:C:849:ARG:NE	2.35	0.47
5:G:44:DT:H2'	5:G:45:DG:C8	2.49	0.47
1:A:534:HIS:CG	1:A:587:LYS:HD2	2.49	0.47
2:B:214:ASP:OD2	2:B:237:HIS:HE1	1.96	0.47
1:C:993:PHE:HE2	1:C:1013:HIS:HA	1.79	0.47
2:D:158:GLY:N	2:D:204:GLN:O	2.45	0.47
4:F:17:DC:H2"	4:F:18:DC:N1	2.29	0.47
1:A:683:ASP:OD1	1:A:685:SER:N	2.29	0.47
1:A:834:ILE:HD13	1:A:891:LEU:HD12	1.97	0.47
1:C:651:ILE:HB	1:C:662:ILE:HD11	1.94	0.47
2:D:158:GLY:O	2:D:178:CYS:HB3	2.14	0.47
3:E:5:DT:O2	4:F:47:DG:N2	2.48	0.47
1:A:610:GLY:O	1:A:651:ILE:HD12	2.14	0.47
2:B:85:GLY:O	2:B:87:PRO:HD3	2.15	0.47
1:A:626:SER:OG	4:F:34:DG:OP1	2.26	0.47
1:A:700:GLU:OE2	2:B:169:THR:HG21	2.15	0.47
1:A:713:VAL:HG13	1:A:718:ARG:NE	2.30	0.47
2:B:66:SER:HB3	2:B:123:ARG:HB2	1.96	0.47
1:C:489:ILE:O	1:C:493:THR:HG22	2.14	0.47
1:C:538:TRP:CE3	1:C:704:MET:CE	2.96	0.47
1:C:835:GLY:HA3	1:C:851:TRP:CE2	2.49	0.47
2:D:27:ASP:OD1	2:D:28:VAL:N	2.47	0.47



	the page	Interatomic	Clash
Atom-1	Atom-2	distance (\AA)	overlap (Å)
2:D:145:ILE:HD11	2:D:152:ALA:HB3	1.96	0.47
1:A:641:ARG:NH1	1:A:987:ASN:N	2.62	0.47
1:A:738:GLU:HG3	1:A:738:GLU:O	2.14	0.47
1:A:817:HIS:C	1:A:819:ASP:N	2.67	0.47
1:A:959:HIS:O	1:A:963:ALA:HB3	2.15	0.47
1:A:985:SER:O	1:A:988:LYS:HB3	2.15	0.47
1:A:988:LYS:NZ	3:E:20:DA:H3'	2.30	0.47
1:C:628:LYS:N	1:C:994:ARG:HH12	2.13	0.47
2:D:41:CYS:SG	2:D:43:THR:HB	2.54	0.47
4:F:28:DC:H2'	4:F:28:DC:OP2	2.14	0.47
5:G:32:DC:H2"	5:G:33:DC:C5	2.45	0.47
2:B:333:SER:HG	2:B:344:HIS:CG	2.18	0.47
2:B:339:PRO:HD2	2:B:341:GLU:O	2.14	0.47
1:C:779:TYR:CE2	1:C:804:LYS:HB2	2.49	0.47
4:F:17:DC:H2"	4:F:18:DC:C6	2.50	0.47
1:A:410:ARG:HB2	5:G:52:DA:C2	2.50	0.47
1:C:961:THR:HA	1:C:965:VAL:HG23	1.97	0.47
4:F:33:DT:H1'	4:F:34:DG:H5'	1.97	0.47
2:B:67:SER:OG	2:B:124:CYS:O	2.24	0.47
2:B:78:ALA:HB3	2:B:91:LEU:HB2	1.97	0.47
2:B:131:GLY:HA3	2:B:132:ASP:C	2.35	0.47
1:C:512:GLY:HA2	1:C:513:ARG:HB3	1.97	0.47
1:C:582:LEU:HD23	1:C:582:LEU:HA	1.55	0.47
2:D:268:TYR:O	2:D:269:HIS:ND1	2.46	0.47
2:B:302:ARG:HH11	2:B:351:GLN:HG3	1.79	0.46
1:C:486:CYS:SG	1:C:522:LEU:HD11	2.56	0.46
1:C:708:ARG:HG2	1:C:721:ARG:HH11	1.79	0.46
4:F:15:DT:H2"	4:F:16:DT:C5	2.49	0.46
5:G:22:DT:H2"	5:G:23:DG:H8	1.80	0.46
1:A:956:ASN:HD21	1:A:960:LYS:HZ2	1.63	0.46
2:B:93:HIS:HB2	2:B:106:LEU:HA	1.97	0.46
1:C:487:LEU:HB2	1:C:522:LEU:HD13	1.97	0.46
1:C:752:CYS:SG	1:C:766:HIS:ND1	2.84	0.46
1:C:814:ASP:OD1	1:C:816:LEU:N	2.47	0.46
2:D:80:PHE:CD2	2:D:89:CYS:HB2	2.46	0.46
5:G:22:DT:H2"	5:G:23:DG:C8	2.51	0.46
1:A:908:TYR:HB2	1:A:938:PHE:HE2	1.80	0.46
1:C:543:LYS:O	1:C:545:VAL:HG23	2.15	0.46
2:D:97:THR:HB	2:D:99:ASN:O	2.15	0.46
2:D:233:LEU:HB3	2:D:252:LEU:HD23	1.98	0.46
5:G:56:DC:H2'	5:G:57:DT:H71	1.97	0.46



	bus page	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:565:ASP:OD1	1:A:566:ASP:N	2.48	0.46
1:A:649:ILE:HG23	1:A:663:PHE:HB3	1.97	0.46
1:A:942:LEU:HD21	1:A:950:TYR:CE2	2.50	0.46
1:C:742:ALA:HA	1:C:743:SER:HA	1.52	0.46
3:E:8:DC:H2"	3:E:9:DC:C6	2.50	0.46
3:E:34:DG:N1	4:F:17:DC:C2	2.76	0.46
5:G:10:DT:C2	5:G:11:DG:N7	2.83	0.46
5:G:32:DC:N4	6:H:30:DG:C6	2.72	0.46
1:A:558:SER:OG	2:B:173:ASN:N	2.46	0.46
1:A:485:VAL:HG13	1:A:1024:PHE:HB3	1.98	0.46
1:C:528:GLU:HA	1:C:533:PHE:CD2	2.50	0.46
1:C:576:PHE:CE1	1:C:1010:LEU:HD21	2.51	0.46
2:D:212:ARG:NH1	2:D:269:HIS:CD2	2.83	0.46
1:C:961:THR:HG23	1:C:962:LEU:HG	1.98	0.46
1:C:975:ILE:C	1:C:977:ALA:H	2.19	0.46
1:C:1019:LYS:O	1:C:1022:GLN:N	2.49	0.46
4:F:18:DC:H2'	4:F:19:DA:C8	2.51	0.46
1:A:409:GLY:HA2	6:H:9:DT:O2	2.15	0.46
1:C:487:LEU:O	1:C:491:VAL:HG13	2.16	0.46
4:F:26:DA:H2"	4:F:27:DG:C8	2.50	0.46
1:A:505:ARG:NH1	1:C:1029:LYS:C	2.65	0.46
2:B:306:GLN:OE1	2:B:306:GLN:N	2.40	0.46
2:D:72:LEU:HB3	2:D:95:GLY:HA3	1.97	0.46
2:D:229:ARG:NH2	2:D:280:GLU:OE2	2.49	0.46
1:A:416:LEU:HD21	1:C:448:SER:HB3	1.98	0.46
1:A:422:ALA:HB1	6:H:12:DT:H5"	1.97	0.46
1:A:693:ILE:O	1:A:696:PRO:HD2	2.16	0.46
2:B:99:ASN:CG	2:B:101:GLU:OE2	2.53	0.46
1:C:725:ARG:HE	1:C:810:GLN:NE2	2.14	0.46
1:C:795:ARG:NH2	2:D:39:ARG:NH1	2.63	0.46
2:D:261:ALA:HB1	2:D:273:ILE:HG23	1.97	0.46
1:A:534:HIS:ND1	1:A:587:LYS:HD2	2.31	0.45
1:A:544:ASN:O	1:A:544:ASN:ND2	2.49	0.45
1:C:914:VAL:HG21	1:C:931:TYR:HD2	1.80	0.45
1:C:1023:LYS:HA	1:C:1023:LYS:HD3	1.77	0.45
2:D:49:ARG:NH1	2:D:58:ARG:HB3	2.31	0.45
2:D:88:GLU:HG3	2:D:89:CYS:N	2.31	0.45
1:A:1021:LEU:HD12	1:A:1021:LEU:H	1.82	0.45
2:D:49:ARG:HH11	2:D:58:ARG:HB3	1.80	0.45
2:B:277:TYR:HD1	2:B:283:LYS:HA	1.80	0.45
2:B:302:ARG:HE	2:B:350:PHE:HE1	1.64	0.45



	bus puge	Interatomic	Clash
Atom-1	Atom-2	distance $(Å)$	overlap (Å)
2:D:96:ARG:HD2	2:D:138:TYR:CD2	2.51	0.45
3:E:15:DT:C2	3:E:16:DA:C6	3.04	0.45
5:G:30:DC:C2	5:G:31:DT:C4	3.04	0.45
1:A:716:LEU:HB2	1:A:718:ARG:HE	1.81	0.45
1:A:919:CYS:HB3	1:A:922:ARG:NH1	2.31	0.45
2:B:10:ASN:HD22	2:B:11:CYS:CA	2.29	0.45
2:B:204:GLN:HB3	2:B:223:ILE:CG1	2.40	0.45
2:D:333:SER:HB2	2:D:342:ALA:HA	1.99	0.45
1:A:486:CYS:SG	1:A:500:TYR:OH	2.48	0.45
1:A:621:GLY:HA2	1:A:640:VAL:HA	1.99	0.45
2:B:49:ARG:HH12	2:B:58:ARG:HH12	1.62	0.45
2:B:318:PHE:CE1	2:B:330:ALA:HB3	2.51	0.45
1:C:482:HIS:HB3	1:C:485:VAL:HG23	1.98	0.45
1:C:795:ARG:HH12	2:D:39:ARG:HH12	1.62	0.45
2:D:87:PRO:HA	2:D:88:GLU:HA	1.71	0.45
4:F:13:DT:C2	4:F:14:DG:C8	3.05	0.45
1:C:651:ILE:HD11	1:C:653:LEU:HB2	1.98	0.45
2:D:242:LEU:HA	2:D:243:GLY:HA2	1.53	0.45
1:A:624:ASP:HA	1:A:637:GLU:CB	2.39	0.45
3:E:20:DA:H2"	3:E:21:DG:H8	1.82	0.45
4:F:41:DA:H2"	4:F:42:DG:C8	2.51	0.45
5:G:12:DT:H2"	5:G:13:DC:C6	2.52	0.45
5:G:34:DA:H2"	5:G:35:DC:O4'	2.17	0.45
1:C:431:LYS:NZ	3:E:34:DG:OP1	2.42	0.45
1:C:859:LEU:O	1:C:865:LEU:N	2.41	0.45
2:D:310:GLU:OE2	2:D:344:HIS:ND1	2.50	0.45
4:F:13:DT:C2	4:F:14:DG:N7	2.84	0.45
1:A:459:ARG:HH21	1:C:420:ARG:NH1	2.14	0.45
1:C:669:ASN:O	1:C:982:GLY:HA2	2.17	0.45
1:A:450:CYS:SG	1:C:454:PHE:HA	2.57	0.45
1:A:587:LYS:HZ1	1:A:711:ILE:HG23	1.82	0.45
2:B:65:ASN:C	2:B:123:ARG:HH11	2.14	0.45
1:C:880:MET:HG3	1:C:909:LEU:HD21	1.99	0.45
2:D:86:LYS:HA	2:D:86:LYS:HD2	1.78	0.45
2:D:234:ILE:HA	2:D:234:ILE:HD13	1.63	0.45
2:D:274:PHE:CZ	2:D:348:VAL:HG21	2.52	0.45
2:B:208:VAL:HG21	2:B:261:ALA:HB3	1.98	0.44
1:C:613:VAL:HG12	1:C:615:VAL:HG23	2.00	0.44
1:A:609:SER:HB3	1:A:653:LEU:HD11	1.99	0.44
1:A:692:ALA:HA	2:B:100:ASN:ND2	2.31	0.44
1:A:919:CYS:HB3	1:A:922:ARG:HH12	1.83	0.44



	t i cas pagem	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
2:B:282:GLN:NE2	2:B:283:LYS:O	2.51	0.44
1:C:688:GLU:OE2	2:D:73:ARG:HD3	2.18	0.44
1:C:816:LEU:HD22	1:C:981:GLU:HB2	1.98	0.44
2:D:16:GLN:HG3	2:D:18:GLY:O	2.18	0.44
2:D:148:ARG:HH12	2:D:239:GLU:CB	2.22	0.44
4:F:9:DT:H2'	4:F:10:DT:C6	2.53	0.44
5:G:33:DC:H2'	5:G:34:DA:C5	2.52	0.44
1:A:536:PHE:HB3	1:A:711:ILE:CD1	2.47	0.44
1:A:832:ASP:CG	1:A:949:ARG:HH12	2.21	0.44
1:A:974:SER:OG	1:A:975:ILE:N	2.51	0.44
2:D:1:MET:CA	2:D:349:SER:O	2.65	0.44
5:G:10:DT:C2'	5:G:11:DG:H8	2.26	0.44
6:H:7:DG:O5'	6:H:7:DG:H8	2.00	0.44
1:A:526:GLU:O	1:A:529:LEU:N	2.47	0.44
1:A:529:LEU:HD13	1:A:1021:LEU:HD11	2.00	0.44
1:A:797:ARG:O	1:A:799:LYS:NZ	2.40	0.44
1:C:752:CYS:HB2	1:C:964:HIS:CE1	2.52	0.44
1:C:861:LYS:HE2	1:C:861:LYS:HB3	1.87	0.44
1:C:975:ILE:HG12	1:C:976:GLY:N	2.32	0.44
2:D:39:ARG:HB2	6:H:50:DA:OP1	2.18	0.44
2:D:283:LYS:NZ	2:D:311:ILE:O	2.35	0.44
3:E:27:DC:H2"	3:E:28:DA:C8	2.52	0.44
1:A:893:PRO:O	1:A:898:ARG:NH2	2.46	0.44
1:A:950:TYR:HB3	1:A:954:ILE:HD11	2.00	0.44
2:D:339:PRO:HB2	2:D:341:GLU:H	1.82	0.44
2:B:135:SER:OG	2:B:137:ARG:NH2	2.50	0.44
2:B:274:PHE:CZ	2:B:348:VAL:CG2	3.01	0.44
2:D:19:PHE:HE2	2:D:21:LEU:HD21	1.83	0.44
3:E:1:DG:H2"	3:E:2:DA:C8	2.53	0.44
4:F:31:DT:OP2	4:F:31:DT:H6	1.99	0.44
2:B:141:THR:OG1	2:B:156:PHE:O	2.29	0.44
1:C:709:LEU:HD12	1:C:710:ILE:H	1.81	0.44
1:C:778:ARG:HH21	1:C:799:LYS:HB2	1.83	0.44
1:C:903:LYS:NZ	1:C:907:LEU:HG	2.32	0.44
1:C:975:ILE:HG12	1:C:976:GLY:H	1.82	0.44
2:D:219:LEU:HD11	2:D:256:LEU:N	2.32	0.44
5:G:24:DG:H1'	5:G:25:DT:H5'	2.00	0.44
1:A:611:PHE:HB3	1:A:649:ILE:HD11	1.99	0.44
1:A:658:ASP:OD1	1:A:659:GLY:N	2.50	0.44
1:A:950:TYR:HB3	1:A:954:ILE:CD1	2.47	0.44
2:B:277:TYR:HA	2:B:284:ARG:N	2.30	0.44



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:C:725:ARG:HE	1:C:810:GLN:HE21	1.65	0.44
1:C:985:SER:O	1:C:987:ASN:N	2.50	0.44
1:C:1011:LYS:NZ	3:E:24:DC:OP1	2.51	0.44
2:D:257:THR:O	2:D:284:ARG:NH1	2.51	0.44
1:C:586:LEU:HD23	1:C:586:LEU:HA	1.75	0.44
1:C:669:ASN:OD1	1:C:979:ALA:HB3	2.17	0.44
2:D:43:THR:HG22	2:D:45:ILE:H	1.82	0.44
5:G:11:DG:C4	5:G:12:DT:C5	3.05	0.44
1:A:649:ILE:HG22	1:A:663:PHE:HD2	1.83	0.43
1:C:691:THR:HG22	1:C:806:PHE:CZ	2.53	0.43
3:E:33:DG:H5'	3:E:33:DG:C8	2.52	0.43
2:D:197:LEU:HA	2:D:197:LEU:HD23	1.69	0.43
5:G:30:DC:H2"	5:G:31:DT:OP2	2.17	0.43
2:B:217:TYR:HE1	2:B:235:ARG:HG3	1.83	0.43
1:C:731:GLU:OE2	1:C:960:LYS:NZ	2.36	0.43
1:C:946:PHE:HA	1:C:948:TYR:CE1	2.53	0.43
6:H:8:DG:H2'	6:H:9:DT:C6	2.53	0.43
1:A:511:SER:O	1:A:513:ARG:N	2.51	0.43
2:B:8:ALA:HA	2:B:55:LEU:HB3	2.00	0.43
2:D:129:LEU:HB3	2:D:192:CYS:SG	2.58	0.43
2:D:325:GLY:O	2:D:350:PHE:N	2.46	0.43
3:E:15:DT:H1'	3:E:16:DA:N7	2.33	0.43
5:G:5:DT:OP2	5:G:5:DT:H6	2.02	0.43
1:A:755:THR:HG23	1:A:758:GLU:OE2	2.18	0.43
1:A:778:ARG:HD3	1:A:798:VAL:O	2.18	0.43
2:B:49:ARG:NH2	2:B:58:ARG:NH1	2.66	0.43
2:B:197:LEU:HD23	2:B:197:LEU:HA	1.88	0.43
1:C:522:LEU:HA	1:C:525:ALA:HB3	2.01	0.43
1:C:616:LYS:HB2	1:C:647:MET:HG3	1.99	0.43
1:C:686:ASP:OD1	1:C:686:ASP:N	2.51	0.43
2:D:96:ARG:HG2	2:D:102:LEU:CD2	2.48	0.43
4:F:27:DG:O5'	4:F:27:DG:H8	2.01	0.43
1:A:893:PRO:C	1:A:898:ARG:HH21	2.22	0.43
1:C:859:LEU:HB3	1:C:865:LEU:HB2	2.01	0.43
1:C:860:ARG:HH12	1:C:864:LYS:HE2	1.83	0.43
2:D:228:CYS:SG	2:D:230:PRO:HD3	2.59	0.43
2:D:317:TRP:HB3	2:D:331:ILE:HD13	2.00	0.43
5:G:33:DC:N3	6:H:30:DG:N2	2.66	0.43
1:A:836:GLU:O	1:A:840:LYS:HB3	2.18	0.43
1:A:882:ARG:HA	1:A:885:VAL:HB	2.00	0.43
1:C:972:ASP:OD1	1:C:972:ASP:N	2.52	0.43



	bus page	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
3:E:26:DA:C2'	3:E:27:DC:OP2	2.66	0.43
5:G:11:DG:C5	5:G:12:DT:C7	3.01	0.43
1:C:621:GLY:HA2	1:C:640:VAL:HG22	2.01	0.43
1:C:716:LEU:HD12	1:C:718:ARG:HH21	1.84	0.43
1:C:840:LYS:HZ2	1:C:843:PRO:HA	1.81	0.43
3:E:5:DT:C2	4:F:47:DG:N2	2.86	0.43
3:E:45:DC:N4	4:F:6:DG:H22	2.17	0.43
1:A:870:ARG:NH2	1:A:872:ASN:HA	2.34	0.43
2:B:137:ARG:HB2	2:B:159:ARG:O	2.19	0.43
1:C:735:ARG:HH12	1:C:742:ALA:HA	1.81	0.43
1:C:795:ARG:CZ	2:D:39:ARG:NH1	2.82	0.43
3:E:16:DA:H5'	3:E:16:DA:C8	2.54	0.43
3:E:40:DA:H2"	3:E:41:DA:OP2	2.19	0.43
4:F:43:DG:H4'	4:F:44:DC:OP1	2.18	0.43
1:A:810:GLN:HG3	1:A:971:ARG:HH11	1.84	0.43
2:B:323:GLY:O	2:B:326:THR:OG1	2.30	0.43
1:C:587:LYS:O	1:C:590:GLU:HB2	2.19	0.43
1:C:675:ARG:HD3	1:C:1017:THR:HB	1.99	0.43
1:C:811:PRO:HG2	1:C:971:ARG:HH12	1.84	0.43
6:H:4:DC:H2"	6:H:5:DA:C8	2.54	0.43
6:H:29:DG:H1'	6:H:30:DG:H5'	2.00	0.43
1:A:787:PHE:CG	1:A:797:ARG:NH1	2.87	0.42
2:B:22:LEU:HD22	2:B:90:TYR:CD1	2.54	0.42
2:B:109:LEU:HD21	2:B:122:LEU:HD13	2.01	0.42
2:B:167:ARG:NH1	2:B:172:TRP:CE2	2.87	0.42
1:C:545:VAL:HA	2:D:169:THR:HG23	2.01	0.42
1:C:681:PHE:HZ	1:C:994:ARG:HH21	1.67	0.42
1:C:904:LEU:HD12	1:C:904:LEU:O	2.19	0.42
1:A:577:ARG:NH2	1:A:579:ASP:OD2	2.49	0.42
1:A:997:ASN:HD22	1:A:1012:HIS:CE1	2.37	0.42
1:C:667:LYS:HE3	1:C:670:SER:OG	2.18	0.42
3:E:8:DC:H2"	3:E:9:DC:C5	2.54	0.42
6:H:37:DA:H2"	6:H:38:DC:H6	1.84	0.42
1:A:683:ASP:OD1	1:A:684:GLU:N	2.52	0.42
2:B:9:VAL:HG11	2:B:56:LYS:HD2	2.01	0.42
2:B:66:SER:HB3	2:B:123:ARG:HA	2.01	0.42
1:C:729:TYR:HB2	1:C:734:VAL:CG2	2.49	0.42
1:C:813:LEU:HD23	1:C:813:LEU:H	1.84	0.42
3:E:31:DC:OP2	3:E:31:DC:C6	2.73	0.42
4:F:46:DA:O5'	4:F:46:DA:C8	2.72	0.42
2:B:83:GLN:N	2:B:87:PRO:HB3	2.34	0.42



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
2:B:207:HIS:ND1	2:B:207:HIS:O	2.53	0.42
2:B:268:TYR:N	2:B:270:GLU:OE2	2.53	0.42
2:B:341:GLU:OE1	2:B:343:TYR:N	2.51	0.42
1:C:579:ASP:OD1	1:C:579:ASP:N	2.53	0.42
1:C:680:MET:HG2	1:C:693:ILE:HG21	2.01	0.42
1:C:810:GLN:OE1	1:C:811:PRO:HD2	2.20	0.42
1:C:828:LYS:HE2	1:C:949:ARG:HH22	1.84	0.42
2:D:22:LEU:HD13	2:D:90:TYR:CG	2.54	0.42
1:A:487:LEU:HA	1:A:522:LEU:HD11	2.01	0.42
2:B:49:ARG:HH12	2:B:58:ARG:HH11	1.67	0.42
2:B:217:TYR:CE1	2:B:235:ARG:HG3	2.53	0.42
2:B:233:LEU:HD23	2:B:234:ILE:N	2.35	0.42
1:C:573:SER:O	1:C:574:ARG:NH1	2.45	0.42
1:C:828:LYS:HE2	1:C:828:LYS:HB3	1.80	0.42
5:G:2:DA:H1'	5:G:3:DT:H5'	2.02	0.42
1:A:610:GLY:O	1:A:652:ARG:N	2.53	0.42
2:B:22:LEU:HD11	2:B:92:ILE:HD11	2.01	0.42
2:B:162:MET:HE1	2:B:171:ASN:HB3	2.02	0.42
2:B:294:ASP:N	2:B:294:ASP:OD1	2.50	0.42
1:C:818:CYS:SG	1:C:819:ASP:N	2.93	0.42
1:A:681:PHE:CE1	1:A:1006:LEU:HD21	2.54	0.42
1:A:778:ARG:HB3	1:A:798:VAL:HG22	2.01	0.42
2:B:135:SER:N	2:B:137:ARG:HH21	2.17	0.42
2:B:332:PRO:HA	2:B:342:ALA:O	2.20	0.42
1:C:732:LYS:HE3	1:C:732:LYS:HB2	1.89	0.42
1:C:860:ARG:NH1	1:C:864:LYS:HE2	2.35	0.42
1:C:971:ARG:NH1	1:C:972:ASP:OD2	2.52	0.42
2:D:21:LEU:O	2:D:22:LEU:HD23	2.20	0.42
1:A:594:MET:O	1:A:598:ARG:HG2	2.19	0.42
1:C:834:ILE:HD13	1:C:834:ILE:HA	1.94	0.42
1:C:956:ASN:O	1:C:959:HIS:HB3	2.19	0.42
5:G:49:DA:N3	6:H:14:DG:N2	2.68	0.42
1:A:421:ARG:HH12	1:A:425:HIS:HD2	1.68	0.42
1:A:771:SER:O	1:A:774:GLU:HG2	2.19	0.42
1:A:821:GLY:O	1:A:825:GLU:HB2	2.20	0.42
2:B:184:LEU:HG	2:B:195:HIS:CD2	2.55	0.42
2:B:201:THR:HB	2:B:232:ARG:HH12	1.85	0.42
1:C:883:GLU:OE1	1:C:883:GLU:N	2.48	0.42
2:D:1:MET:HA	2:D:349:SER:O	2.20	0.42
4:F:35:DT:H2"	4:F:36:DA:C8	2.55	0.42
5:G:17:DC:H6	5:G:17:DC:H2'	1.75	0.42



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:487:LEU:HD11	1:A:1015:LEU:HD13	2.02	0.42
1:A:580:VAL:O	1:A:583:VAL:HG22	2.20	0.42
1:A:709:LEU:O	1:A:719:SER:HA	2.20	0.42
2:B:73:ARG:NH1	2:B:96:ARG:NH1	2.66	0.42
1:C:991:ARG:HA	1:C:991:ARG:HD2	1.87	0.42
5:G:54:DC:H2"	5:G:55:DC:C6	2.54	0.42
1:A:892:VAL:CG1	1:A:898:ARG:HG2	2.44	0.41
2:B:21:LEU:O	2:B:22:LEU:HD23	2.20	0.41
1:C:1006:LEU:HD23	1:C:1006:LEU:HA	1.58	0.41
3:E:6:DG:OP2	3:E:6:DG:C8	2.70	0.41
5:G:11:DG:N1	5:G:12:DT:C4	2.88	0.41
6:H:42:DT:H3'	6:H:43:DG:C8	2.55	0.41
1:A:982:GLY:O	1:A:985:SER:HB3	2.19	0.41
2:B:177:ASP:N	2:B:177:ASP:OD1	2.52	0.41
1:C:512:GLY:HA2	1:C:513:ARG:C	2.40	0.41
1:C:588:ASP:OD1	1:C:1019:LYS:HB2	2.20	0.41
2:D:93:HIS:HA	2:D:94:GLY:HA2	1.66	0.41
2:D:283:LYS:CD	2:D:317:TRP:HE1	2.33	0.41
2:D:328:LEU:HD11	2:D:345:PHE:CD2	2.51	0.41
3:E:33:DG:H2"	3:E:34:DG:C8	2.55	0.41
1:A:642:PHE:HB3	1:A:678:CYS:HB3	2.01	0.41
2:B:180:PRO:HB3	2:B:203:GLY:N	2.36	0.41
2:D:43:THR:HG22	2:D:45:ILE:N	2.34	0.41
2:D:315:ARG:HG3	2:D:316:THR:HG23	2.01	0.41
2:D:321:SER:O	2:D:321:SER:OG	2.36	0.41
3:E:18:DA:H2"	3:E:19:DC:O5'	2.20	0.41
1:C:519:LEU:HD23	1:C:519:LEU:HA	1.71	0.41
1:C:705:MET:H	1:C:705:MET:HG2	1.67	0.41
3:E:34:DG:C8	3:E:35:DA:C6	3.08	0.41
5:G:42:DG:O5'	5:G:42:DG:H8	2.03	0.41
6:H:5:DA:H2"	6:H:6:DG:N7	2.35	0.41
1:A:851:TRP:HA	1:A:854:THR:HG22	2.02	0.41
1:A:895:GLU:OE2	1:A:898:ARG:NH2	2.53	0.41
1:A:991:ARG:HH11	4:F:34:DG:C3'	2.32	0.41
1:C:864:LYS:HD3	1:C:864:LYS:HA	1.95	0.41
2:D:24:LEU:CD2	2:D:29:TYR:CD2	3.03	0.41
2:D:202:ASP:OD2	2:D:223:ILE:HG21	2.21	0.41
2:D:287:CYS:SG	2:D:304:PRO:HA	2.61	0.41
4:F:13:DT:H6	4:F:13:DT:H2'	1.66	0.41
1:A:822:ASN:OD1	1:A:822:ASN:N	2.52	0.41
1:C:610:GLY:C	1:C:651:ILE:HD12	2.40	0.41



		Interatomic	Clash
Atom-1	Atom-2	distance $(Å)$	overlap (Å)
1:C:895:GLU:OE1	1:C:898:ARG:NH1	2.54	0.41
3:E:29:DG:H1'	3:E:30:DA:C8	2.55	0.41
4:F:47:DG:H2"	4:F:48:DA:OP2	2.20	0.41
6:H:42:DT:H3'	6:H:43:DG:N7	2.35	0.41
2:B:142:LEU:HD12	2:B:154:VAL:O	2.21	0.41
1:C:498:SER:O	1:C:502:LYS:HG3	2.20	0.41
1:C:534:HIS:CE1	1:C:587:LYS:NZ	2.89	0.41
1:C:628:LYS:HB2	4:F:32:DG:H4'	2.01	0.41
2:D:217:TYR:OH	2:D:235:ARG:NH2	2.52	0.41
4:F:13:DT:H4'	4:F:14:DG:OP1	2.21	0.41
6:H:7:DG:H2'	6:H:8:DG:H8	1.85	0.41
6:H:45:DG:H2'	6:H:45:DG:N3	2.36	0.41
1:A:489:ILE:HD11	1:C:503:MET:HA	2.02	0.41
1:A:647:MET:HE1	1:A:977:ALA:O	2.21	0.41
1:A:740:LEU:HD21	1:A:770:ARG:HH22	1.85	0.41
1:A:903:LYS:HE2	1:A:907:LEU:HD11	2.02	0.41
1:C:452:THR:O	1:C:456:LEU:HG	2.21	0.41
1:C:738:GLU:HA	1:C:805:PRO:HB3	2.01	0.41
1:C:838:TYR:OH	1:C:897:ARG:HG3	2.21	0.41
1:C:1029:LYS:HE2	1:C:1029:LYS:HB3	1.78	0.41
2:D:286:GLU:HA	2:D:304:PRO:HB3	2.03	0.41
5:G:11:DG:C5	5:G:12:DT:H73	2.55	0.41
1:A:430:LEU:HB2	1:C:437:PHE:CG	2.56	0.41
1:A:820:ILE:O	1:A:824:THR:HG22	2.20	0.41
1:A:860:ARG:HH21	1:A:867:PRO:HD2	1.85	0.41
1:A:968:ILE:O	1:A:971:ARG:HG2	2.20	0.41
1:A:1019:LYS:HB2	1:A:1019:LYS:HE2	1.74	0.41
1:A:1024:PHE:HD1	1:A:1024:PHE:HA	1.73	0.41
2:B:76:ALA:CB	2:B:93:HIS:O	2.57	0.41
2:B:212:ARG:HG2	2:B:213:GLN:OE1	2.20	0.41
1:C:485:VAL:HG22	1:C:1024:PHE:CG	2.56	0.41
1:C:789:GLU:CD	1:C:797:ARG:HE	2.24	0.41
1:C:795:ARG:HH22	2:D:39:ARG:NH1	2.18	0.41
2:D:130:VAL:O	2:D:192:CYS:N	2.37	0.41
2:D:256:LEU:CD2	2:D:288:THR:HG21	2.51	0.41
2:D:290:VAL:CG1	2:D:297:VAL:HG23	2.49	0.41
4:F:14:DG:C5	4:F:15:DT:C4	3.09	0.41
4:F:15:DT:H2"	4:F:16:DT:C4	2.55	0.41
1:A:503:MET:O	1:A:507:VAL:HG12	2.21	0.41
1:A:578:TYR:OH	1:A:724:PHE:HE2	2.01	0.41
2:B:180:PRO:O	2:B:200:LEU:HD23	2.21	0.41



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:C:481:LEU:HD23	1:C:481:LEU:HA	1.78	0.41
1:C:492:ASN:OD1	1:C:1025:MET:HG2	2.20	0.41
1:C:798:VAL:O	1:C:801:VAL:HG23	2.20	0.41
1:C:840:LYS:HD3	1:C:843:PRO:HB3	2.03	0.41
2:D:259:THR:OG1	2:D:277:TYR:O	2.26	0.41
2:D:275:GLY:HA3	2:D:317:TRP:HH2	1.86	0.41
3:E:33:DG:C2	3:E:34:DG:C6	3.09	0.41
6:H:51:DC:H2"	6:H:52:DA:C8	2.56	0.41
1:A:682:VAL:HG12	1:A:683:ASP:N	2.33	0.40
1:A:897:ARG:HH12	1:A:945:MET:HG3	1.86	0.40
1:A:988:LYS:HZ2	3:E:20:DA:P	2.44	0.40
1:A:993:PHE:HE2	1:A:1013:HIS:HA	1.86	0.40
2:B:210:LEU:HD23	2:B:210:LEU:HA	1.88	0.40
2:D:147:SER:OG	2:D:240:LEU:CD2	2.61	0.40
5:G:11:DG:N2	5:G:12:DT:C2	2.90	0.40
1:A:816:LEU:HD13	1:A:981:GLU:HG2	2.03	0.40
2:B:273:ILE:HB	2:B:288:THR:OG1	2.21	0.40
1:C:665:GLU:HG2	1:C:668:PRO:HB3	2.03	0.40
1:C:778:ARG:NH2	1:C:798:VAL:O	2.49	0.40
2:D:274:PHE:HZ	2:D:348:VAL:HG11	1.86	0.40
4:F:31:DT:OP2	4:F:31:DT:C6	2.74	0.40
1:A:593:ILE:HD11	1:A:713:VAL:HG21	2.03	0.40
1:A:862:LYS:HE2	1:A:887:ALA:HB1	2.02	0.40
1:A:894:SER:O	1:A:898:ARG:HG3	2.22	0.40
2:B:48:VAL:HG23	2:B:55:LEU:CD1	2.48	0.40
2:B:197:LEU:HD12	2:B:249:CYS:HB3	2.03	0.40
2:B:269:HIS:O	2:B:292:LEU:HD23	2.22	0.40
1:C:833:GLU:OE2	1:C:948:TYR:OH	2.35	0.40
1:C:989:LEU:HA	1:C:989:LEU:HD23	1.74	0.40
3:E:45:DC:H2"	3:E:46:DT:OP2	2.20	0.40
4:F:27:DG:C2'	4:F:28:DC:H5'	2.51	0.40
5:G:34:DA:OP2	5:G:34:DA:C8	2.73	0.40
1:A:529:LEU:HD23	1:A:529:LEU:HA	1.88	0.40
1:A:778:ARG:HE	1:A:799:LYS:CE	2.34	0.40
1:C:577:ARG:HB3	1:C:579:ASP:OD1	2.22	0.40
1:C:597:LEU:HD21	1:C:603:ASP:O	2.22	0.40
1:C:795:ARG:NH1	2:D:39:ARG:HH11	2.18	0.40
2:D:96:ARG:HG2	2:D:102:LEU:HD22	2.03	0.40
2:D:150:LYS:H	2:D:150:LYS:HG2	1.61	0.40
2:D:233:LEU:O	2:D:234:ILE:HD13	2.21	0.40
2:D:234:ILE:HG23	2:D:234:ILE:HD12	1.49	0.40



continuea from prees	cae page		1
Atom 1	Atom 2	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
2:D:238:VAL:HG12	2:D:247:LEU:HG	2.04	0.40
3:E:19:DC:H2"	3:E:20:DA:C8	2.57	0.40
5:G:4:DC:H2"	5:G:5:DT:OP2	2.21	0.40
5:G:32:DC:N4	6:H:30:DG:H1	2.19	0.40
1:A:441:GLU:OE1	1:C:412:ARG:NH1	2.55	0.40
1:A:495:LEU:HD22	1:A:499:GLN:HB2	2.04	0.40
1:A:557:LEU:HD12	1:A:557:LEU:H	1.87	0.40
2:D:131:GLY:HA2	2:D:132:ASP:HB2	2.03	0.40
2:D:159:ARG:HB3	2:D:176:VAL:O	2.21	0.40
3:E:16:DA:H2"	3:E:17:DC:OP2	2.21	0.40
3:E:35:DA:H2"	3:E:36:DA:OP2	2.22	0.40
4:F:17:DC:H6	4:F:17:DC:H2'	1.59	0.40
4:F:20:DG:C2'	4:F:21:DT:H5'	2.52	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 PDB entries followed by that with respect to all EM entries.

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	Perce	ntiles
1	А	619/1159~(53%)	558~(90%)	59 (10%)	2(0%)	41	76
1	С	612/1159~(53%)	551 (90%)	61 (10%)	0	100	100
2	В	349/533~(66%)	306~(88%)	43 (12%)	0	100	100
2	D	349/533~(66%)	312 (89%)	37~(11%)	0	100	100
All	All	1929/3384~(57%)	1727 (90%)	200 (10%)	2(0%)	54	85

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	818	CYS
1	А	512	GLY



5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	А	547/1000~(55%)	521~(95%)	26~(5%)	25 52
1	С	545/1000~(54%)	515~(94%)	30~(6%)	21 49
2	В	303/465~(65%)	290~(96%)	13 (4%)	29 55
2	D	303/465~(65%)	291~(96%)	12 (4%)	31 56
All	All	1698/2930~(58%)	1617~(95%)	81 (5%)	29 52

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	479	PHE
1	А	492	ASN
1	А	497	CYS
1	А	503	MET
1	А	505	ARG
1	А	538	TRP
1	А	590	GLU
1	А	619	CYS
1	А	627	GLU
1	А	629	HIS
1	А	656	GLU
1	А	675	ARG
1	А	678	CYS
1	А	680	MET
1	А	719	SER
1	А	777	GLU
1	А	790	SER
1	А	822	ASN
1	А	844	SER
1	А	865	LEU
1	А	866	LYS
1	А	869	MET
1	А	878	ARG
1	А	933	TYR



Mol	Chain	Res	Type
1	А	945	MET
1	А	971	ARG
2	В	3	LEU
2	В	4	GLN
2	В	24	LEU
2	В	41	CYS
2	В	58	ARG
2	В	62	PHE
2	В	63	SER
2	В	177	ASP
2	В	189	PHE
2	В	284	ARG
2	В	289	TYR
2	В	292	LEU
2	В	293	ASP
1	С	479	PHE
1	С	486	CYS
1	С	490	ARG
1	С	492	ASN
1	С	496	SER
1	С	560	TRP
1	С	573	SER
1	С	578	TYR
1	С	586	LEU
1	С	603	ASP
1	С	619	CYS
1	С	663	PHE
1	С	671	GLU
1	С	677	LEU
1	С	678	CYS
1	С	712	SER
1	C	719	SER
1	С	818	CYS
1	С	849	ARG
1	C	902	LEU
1	С	917	SER
1	C	933	TYR
1	С	972	ASP
1	C	984	GLU
1	С	988	LYS
1	C	994	ARG
1	С	999	ARG



Mol	Chain	Res	Type
1	С	1000	GLN
1	С	1024	PHE
1	С	1029	LYS
2	D	1	MET
2	D	24	LEU
2	D	90	TYR
2	D	103	SER
2	D	167	ARG
2	D	178	CYS
2	D	187	LEU
2	D	205	SER
2	D	212	ARG
2	D	232	ARG
2	D	249	CYS
2	D	279	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	775	ASN
1	А	956	ASN
1	А	964	HIS
2	В	10	ASN
1	С	775	ASN
1	С	1012	HIS
2	D	33	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 monosaccharides in this entry.



5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.
There are no bond length 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 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-7847. These allow visual inspection of the internal detail of the map and identification of artifacts.

No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections (i)

6.1.1 Primary map



The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

6.2.1 Primary map



X Index: 96







The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

6.3.1 Primary map



X Index: 92

Y Index: 82

Z Index: 83

The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

6.4.1 Primary map



The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.



6.5 Orthogonal surface views (i)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.03. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.6 Mask visualisation (i)

This section was not generated. No masks/segmentation were deposited.



7 Map analysis (i)

This section contains the results of statistical analysis of the map.

7.1 Map-value distribution (i)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.



7.2 Volume estimate (i)



The volume at the recommended contour level is 193 $\rm nm^3;$ this corresponds to an approximate mass of 175 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.



7.3 Rotationally averaged power spectrum (i)



*Reported resolution corresponds to spatial frequency of 0.237 ${\rm \AA^{-1}}$



8 Fourier-Shell correlation (i)

This section was not generated. No FSC curve or half-maps provided.



9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-7847 and PDB model 6DBQ. Per-residue inclusion information can be found in section 3 on page 7.

9.1 Map-model overlay (i)



The images above show the 3D surface view of the map at the recommended contour level 0.03 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.



9.2 Q-score mapped to coordinate model (i)



The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model (i)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.03).



9.4 Atom inclusion (i)



At the recommended contour level, 89% of all backbone atoms, 76% of all non-hydrogen atoms, are inside the map.



9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.03) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7580	0.2960
А	0.7870	0.3220
В	0.8450	0.3320
С	0.8000	0.3280
D	0.8280	0.3380
Е	0.6870	0.2030
F	0.7400	0.2100
G	0.4890	0.1660
Н	0.4850	0.1640

