



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 05:57 PM UTC

PDB ID : 1MA7 / pdb_00001ma7
Title : Crystal structure of Cre site-specific recombinase complexed with a mutant DNA substrate, LoxP-A8/T27
Authors : Martin, S.S.; Chu, V.C.; Baldwin, E.P.
Deposited on : 2002-08-01
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

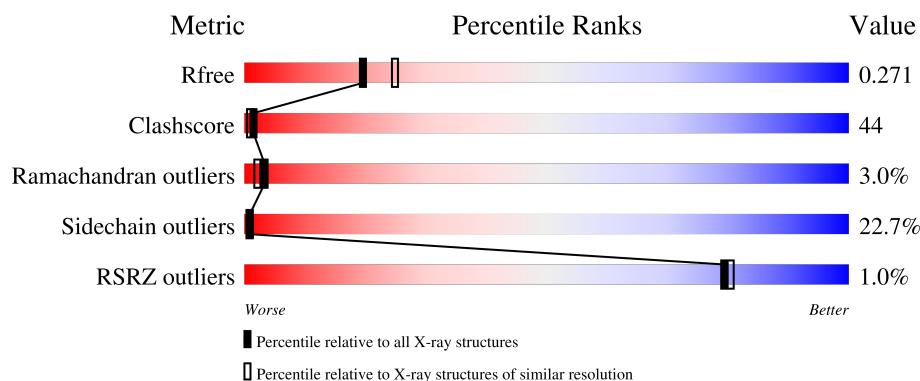
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	C	34	
2	D	34	
3	A	349	
3	B	349	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6699 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 DNA chain called LOXP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	32	Total	C	N	O	P	0	0	0
			654	317	115	191	31			

- Molecule 2 is a DNA chain called LOXP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	34	Total	C	N	O	P	0	0	0
			693	336	123	201	33			

- Molecule 3 is a protein called CRE RECOMBINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	323	Total	C	N	O	S	45	0	0
			2557	1588	487	467	15			
3	B	317	Total	C	N	O	S	11	0	0
			2512	1562	481	454	15			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	HIS	-	expression tag	UNP P06956
A	-3	HIS	-	expression tag	UNP P06956
A	-2	HIS	-	expression tag	UNP P06956
A	-1	HIS	-	expression tag	UNP P06956
A	0	HIS	-	expression tag	UNP P06956
A	1	HIS	-	expression tag	UNP P06956
B	-4	HIS	-	expression tag	UNP P06956
B	-3	HIS	-	expression tag	UNP P06956
B	-2	HIS	-	expression tag	UNP P06956
B	-1	HIS	-	expression tag	UNP P06956
B	0	HIS	-	expression tag	UNP P06956
B	1	HIS	-	expression tag	UNP P06956

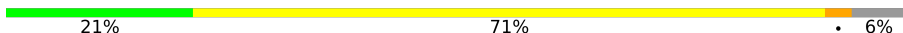
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	38	Total 38	O 38	0	0
4	D	41	Total 41	O 41	0	0
4	A	60	Total 60	O 60	0	0
4	B	144	Total 144	O 144	0	0

3 Residue-property plots [i](#)

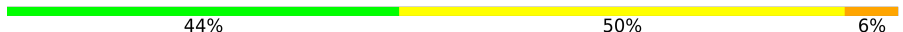
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: LOXP

Chain C: 

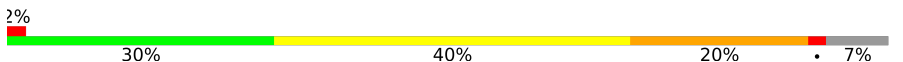


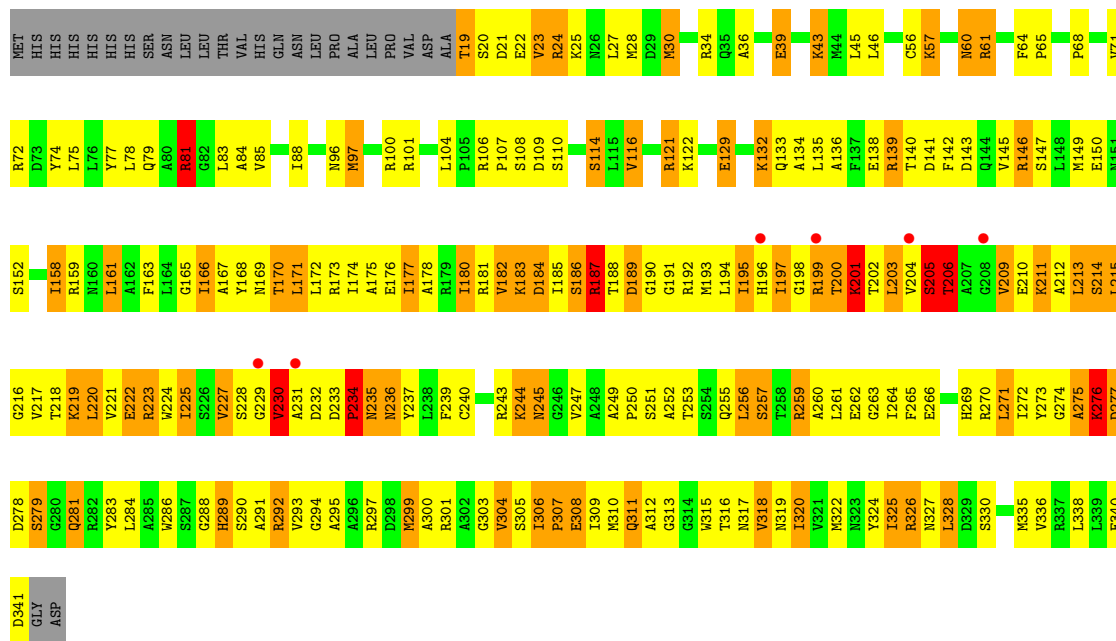
• Molecule 2: LOXP

Chain D: 



• Molecule 3: CRE RECOMBINASE

Chain A: 



• Molecule 3: CRE RECOMBINASE

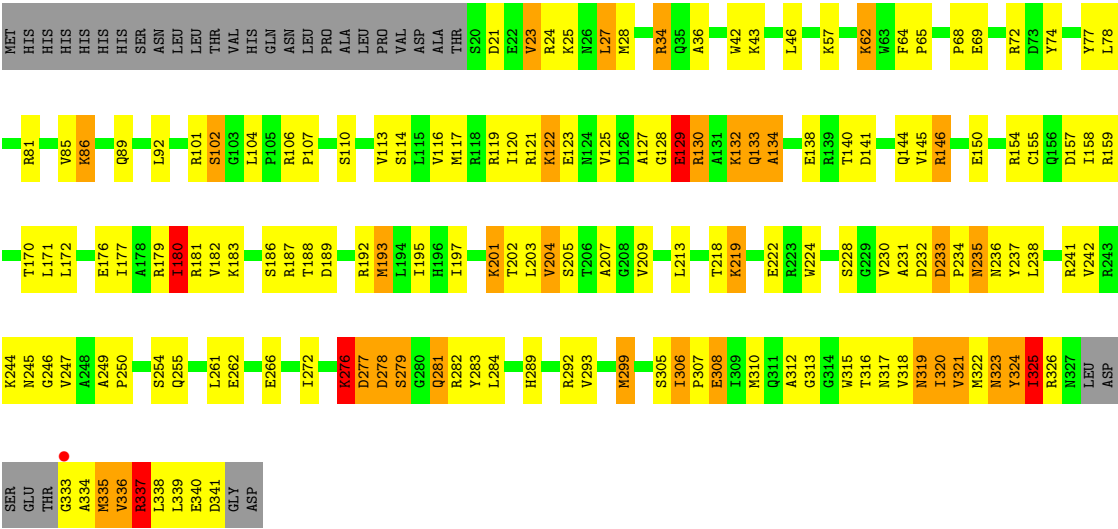
Chain B:

46%

34%

9%

9%



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	107.54Å 122.24Å 180.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	5.00 – 2.30 5.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.0 (5.00-2.30) 87.6 (5.00-2.30)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	10.55 (at 2.24Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.239 , 0.290 0.241 , 0.271	Depositor DCC
R_{free} test set	2787 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	45.1	Xtriage
Anisotropy	0.167	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 93.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6699	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.71% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	C	0.22	0/732	0.99	2/1126 (0.2%)
2	D	0.26	0/777	1.04	2/1197 (0.2%)
3	A	0.62	0/2598	1.15	18/3503 (0.5%)
3	B	0.52	0/2552	0.98	8/3438 (0.2%)
All	All	0.52	0/6659	1.06	30/9264 (0.3%)

There are no bond length outliers.

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	235	ASN	N-CA-C	-12.16	97.72	111.71
3	A	182	VAL	N-CA-C	10.31	121.04	111.45
3	A	281	GLN	N-CA-C	10.13	124.33	110.35
3	B	233	ASP	CA-CB-CG	9.96	122.56	112.60
3	A	304	VAL	N-CA-C	8.70	120.81	108.36
3	A	81	ARG	N-CA-C	-7.24	104.44	113.28
3	A	206	THR	N-CA-C	6.95	117.44	108.34
3	A	307	PRO	N-CA-C	-6.94	103.40	113.75
3	B	320	ILE	N-CA-C	-6.94	103.76	110.42
3	A	278	ASP	N-CA-C	6.75	120.45	110.46
3	A	232	ASP	N-CA-C	-6.55	104.14	111.28
3	B	36	ALA	N-CA-C	-6.41	105.42	113.18
3	A	215	LEU	CB-CA-C	6.36	119.11	111.22
1	C	20	DG	C4-N9-C1'	-6.17	117.74	127.00
3	A	199	ARG	N-CA-C	5.73	119.53	112.03
3	B	276	LYS	N-CA-C	5.71	119.55	109.96
3	B	323	ASN	N-CA-C	5.63	117.09	111.07
3	B	102	SER	N-CA-C	-5.42	106.51	113.23
3	A	220	LEU	N-CA-C	-5.42	105.37	111.28
3	A	227	VAL	N-CA-C	5.36	118.14	112.83
3	A	271	LEU	N-CA-C	-5.36	105.43	111.28
3	B	34	ARG	CB-CA-C	-5.34	102.26	110.81
1	C	20	DG	C8-N9-C1'	5.27	134.91	127.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	304	VAL	CB-CA-C	5.22	117.94	110.84
3	A	211	LYS	N-CA-C	5.20	117.97	110.28
2	D	18	DA	O4'-C1'-N9	-5.07	100.79	108.40
3	A	187	ARG	N-CA-C	-5.06	107.23	113.41
3	A	170	THR	N-CA-C	5.02	118.56	112.23
2	D	15	DC	O4'-C1'-N1	5.01	115.92	108.40
3	B	134	ALA	N-CA-C	5.01	117.31	110.24

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	C	654	0	368	49	0
2	D	693	0	389	38	0
3	A	2557	0	2578	300	0
3	B	2512	0	2537	187	0
4	A	60	0	0	8	0
4	B	144	0	0	2	0
4	C	38	0	0	2	0
4	D	41	0	0	1	0
All	All	6699	0	5872	523	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 44.

All (523) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:145:VAL:HG13	3:A:149:MET:HE3	1.23	1.20
3:A:235:ASN:HB2	3:A:252:ALA:HB1	1.15	1.14
2:D:23:DA:H5'	3:B:201:LYS:HB2	1.31	1.08
3:B:306:ILE:HG23	3:B:310:MET:HE3	1.36	1.06
3:A:169:ASN:HB2	3:A:217:VAL:HG21	1.40	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:182:VAL:HB	3:A:231:ALA:HB2	1.40	1.03
3:A:202:THR:HB	3:A:205:SER:HB2	1.33	1.03
2:D:21:DT:H2''	2:D:22:DT:H5'	1.45	0.98
1:C:16:DG:H5'	3:B:202:THR:HB	1.46	0.96
3:A:213:LEU:HB3	3:A:217:VAL:HB	1.48	0.95
3:A:225:ILE:HG13	3:A:230:VAL:HG23	1.53	0.91
2:D:14:DG:H2''	2:D:15:DC:H6	1.36	0.90
3:A:182:VAL:CB	3:A:231:ALA:HB2	2.02	0.89
3:A:235:ASN:HB2	3:A:252:ALA:CB	2.00	0.88
3:B:299:MET:HE1	3:B:312:ALA:HB2	1.52	0.88
3:A:23:VAL:CG2	3:A:104:LEU:HD21	2.03	0.88
1:C:1:DA:C8	1:C:1:DA:H5'	2.09	0.88
2:D:16:DA:H2''	2:D:17:DT:C7	2.04	0.87
3:A:193:MET:HB2	3:A:218:THR:CG2	2.04	0.87
3:A:236:ASN:N	3:A:252:ALA:HA	1.88	0.86
3:A:202:THR:HG22	3:A:203:LEU:H	1.38	0.86
3:A:65:PRO:HG3	3:A:104:LEU:HD13	1.57	0.86
3:B:299:MET:HE1	3:B:312:ALA:CB	2.06	0.85
3:A:213:LEU:HB2	3:A:218:THR:HG23	1.56	0.85
3:B:130:ARG:HG2	3:B:130:ARG:HH11	1.41	0.85
3:A:200:THR:HG21	3:A:206:THR:HG23	1.58	0.85
3:B:310:MET:HE1	3:B:318:VAL:HG22	1.57	0.84
3:B:203:LEU:HD23	3:B:204:VAL:N	1.93	0.83
3:A:213:LEU:CB	3:A:218:THR:HG23	2.08	0.83
3:A:199:ARG:HH22	3:A:292:ARG:NH2	1.77	0.83
3:A:213:LEU:CB	3:A:217:VAL:HB	2.09	0.83
3:B:133:GLN:HE22	3:B:324:TYR:HA	1.42	0.82
1:C:19:DT:H2''	1:C:20:DG:C8	2.16	0.81
3:B:340:GLU:O	3:B:341:ASP:C	2.23	0.81
3:A:197:ILE:CG2	3:A:209:VAL:HG22	2.11	0.81
3:B:180:ILE:HD13	3:B:195:ILE:HG21	1.61	0.81
3:A:187:ARG:NH2	3:A:193:MET:HE2	1.96	0.80
3:A:275:ALA:C	3:A:276:LYS:HG2	2.05	0.80
1:C:25:DA:C8	1:C:25:DA:H5'	2.17	0.80
3:A:304:VAL:HG12	3:A:308:GLU:HB3	1.62	0.80
3:A:193:MET:HE3	3:A:222:GLU:OE2	1.81	0.80
3:A:299:MET:HE1	3:B:335:MET:HB2	1.62	0.79
3:A:177:ILE:O	3:A:180:ILE:HG23	1.82	0.79
1:C:14:DA:N7	3:B:86:LYS:NZ	2.31	0.79
3:A:231:ALA:O	3:A:234:PRO:HD3	1.83	0.79
3:A:193:MET:HB2	3:A:218:THR:HG22	1.64	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:199:ARG:NH2	3:A:292:ARG:HH21	1.80	0.78
2:D:23:DA:C5'	3:B:201:LYS:HB2	2.12	0.78
3:A:163:PHE:CE2	3:A:261:LEU:HD22	2.19	0.78
3:A:310:MET:HE2	3:A:318:VAL:HG13	1.65	0.77
3:A:213:LEU:HD12	3:A:218:THR:CA	2.14	0.77
3:A:182:VAL:HG23	3:A:236:ASN:O	1.83	0.77
3:A:213:LEU:HD12	3:A:218:THR:HA	1.65	0.77
3:A:245:ASN:N	3:A:245:ASN:HD22	1.79	0.77
3:B:306:ILE:HG22	3:B:307:PRO:HD3	1.66	0.77
1:C:16:DG:H5'	3:B:202:THR:CB	2.15	0.77
3:A:169:ASN:CB	3:A:217:VAL:HG21	2.14	0.77
3:B:306:ILE:CG2	3:B:310:MET:HE3	2.13	0.77
1:C:1:DA:H2''	1:C:2:DT:C5'	2.15	0.76
3:B:85:VAL:O	3:B:89:GLN:HG3	1.86	0.76
3:B:310:MET:CE	3:B:318:VAL:HG13	2.16	0.76
3:A:310:MET:HE1	3:A:318:VAL:HG22	1.68	0.75
1:C:1:DA:H2''	1:C:2:DT:H5''	1.68	0.75
3:B:310:MET:CE	3:B:318:VAL:HG22	2.18	0.74
3:A:101:ARG:HH12	3:B:114:SER:HB3	1.53	0.74
1:C:1:DA:H5'	1:C:1:DA:H8	1.50	0.74
3:A:182:VAL:CG1	3:A:231:ALA:HB2	2.17	0.74
3:A:133:GLN:HE21	3:A:324:TYR:HE1	1.35	0.74
1:C:16:DG:C5'	3:B:202:THR:HB	2.16	0.73
2:D:14:DG:H2''	2:D:15:DC:C6	2.23	0.73
3:A:189:ASP:OD2	3:A:194:LEU:HG	1.89	0.73
3:A:204:VAL:HG23	3:B:323:ASN:HD21	1.54	0.73
3:A:201:LYS:HA	3:B:130:ARG:NH2	2.03	0.72
3:B:146:ARG:O	3:B:150:GLU:HB2	1.89	0.72
3:B:23:VAL:HG21	3:B:104:LEU:HD21	1.70	0.72
3:B:133:GLN:NE2	3:B:324:TYR:HA	2.05	0.72
3:A:166:ILE:HA	3:A:217:VAL:HG11	1.71	0.72
3:B:133:GLN:HE21	3:B:324:TYR:HD1	1.35	0.72
1:C:2:DT:H2''	1:C:3:DA:N7	2.05	0.71
3:A:235:ASN:CB	3:A:252:ALA:HB1	2.09	0.71
3:A:145:VAL:CG1	3:A:149:MET:HE3	2.13	0.71
3:A:193:MET:HB2	3:A:218:THR:HB	1.71	0.71
3:B:24:ARG:NH1	3:B:24:ARG:HB3	2.05	0.71
3:A:202:THR:CB	3:A:205:SER:HB2	2.18	0.71
3:A:165:GLY:C	3:A:217:VAL:HG13	2.15	0.70
3:A:245:ASN:H	3:A:245:ASN:ND2	1.89	0.70
3:A:275:ALA:O	3:A:276:LYS:HG2	1.92	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:304:VAL:CG1	3:A:308:GLU:HB3	2.21	0.70
3:A:310:MET:CE	3:A:318:VAL:HG22	2.22	0.69
3:B:116:VAL:O	3:B:120:ILE:HG13	1.91	0.69
3:A:303:GLY:HA2	4:A:396:HOH:O	1.93	0.69
3:A:259:ARG:HA	3:A:259:ARG:NE	2.07	0.69
1:C:32:DT:H2''	1:C:33:DA:C8	2.27	0.69
2:D:16:DA:H4'	2:D:17:DT:OP1	1.93	0.68
3:A:262:GLU:O	3:A:266:GLU:HG3	1.93	0.68
2:D:16:DA:H2''	2:D:17:DT:H73	1.75	0.68
3:A:297:ARG:HA	3:A:325:ILE:HG22	1.75	0.68
3:A:85:VAL:HG23	3:A:129:GLU:OE2	1.94	0.68
3:A:201:LYS:HA	3:B:130:ARG:HH21	1.58	0.68
1:C:25:DA:H5'	1:C:25:DA:H8	1.59	0.68
3:A:166:ILE:HD11	3:A:177:ILE:HG12	1.75	0.68
3:B:92:LEU:HD23	3:B:117:MET:HG2	1.75	0.67
3:A:213:LEU:HD13	3:A:217:VAL:HG12	1.76	0.67
3:A:245:ASN:HD22	3:A:245:ASN:H	1.39	0.67
3:A:235:ASN:C	3:A:252:ALA:HA	2.19	0.67
3:A:163:PHE:HE2	3:A:261:LEU:HD22	1.60	0.67
3:B:325:ILE:HD12	3:B:325:ILE:N	2.10	0.67
3:B:119:ARG:O	3:B:123:GLU:HG3	1.95	0.66
2:D:18:DA:OP2	3:B:121:ARG:NE	2.28	0.66
3:B:133:GLN:NE2	3:B:324:TYR:HD1	1.93	0.66
2:D:20:DA:OP2	3:B:106:ARG:NH1	2.28	0.66
3:A:193:MET:HB2	3:A:218:THR:CB	2.25	0.66
3:A:272:ILE:HG22	3:A:273:TYR:CD1	2.31	0.66
3:A:197:ILE:HG21	3:A:209:VAL:HG22	1.76	0.65
3:A:235:ASN:HB3	3:A:252:ALA:O	1.96	0.65
3:B:323:ASN:O	3:B:325:ILE:N	2.29	0.65
1:C:28:DA:H2''	1:C:29:DA:H8	1.59	0.65
3:A:24:ARG:HH11	3:A:24:ARG:HG2	1.59	0.65
3:A:313:GLY:HA3	3:A:315:TRP:CZ3	2.31	0.65
1:C:1:DA:C2'	1:C:2:DT:H5''	2.26	0.65
3:B:62:LYS:HB3	4:B:452:HOH:O	1.97	0.65
2:D:11:DA:OP1	3:A:81:ARG:NH2	2.29	0.65
3:A:202:THR:HB	3:A:205:SER:CB	2.18	0.65
3:B:128:GLY:O	3:B:129:GLU:C	2.40	0.65
3:A:306:ILE:N	3:A:307:PRO:HD2	2.12	0.65
3:B:27:LEU:HD11	3:B:102:SER:HB3	1.78	0.65
3:A:158:ILE:HG21	3:A:227:VAL:CG2	2.27	0.65
3:A:158:ILE:HG21	3:A:227:VAL:HG21	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:236:ASN:CA	3:A:252:ALA:HA	2.26	0.65
1:C:28:DA:H2''	1:C:29:DA:C8	2.31	0.64
3:A:197:ILE:HD13	3:A:198:GLY:N	2.12	0.64
3:A:204:VAL:HG23	3:B:323:ASN:ND2	2.11	0.64
3:A:214:SER:C	3:A:216:GLY:H	2.06	0.64
3:B:187:ARG:NH2	3:B:222:GLU:OE2	2.28	0.64
3:A:209:VAL:HG23	3:A:210:GLU:N	2.13	0.64
3:A:310:MET:CE	3:A:318:VAL:HG13	2.27	0.64
1:C:1:DA:H1'	1:C:2:DT:H5''	1.80	0.64
3:A:276:LYS:O	3:A:277:ASP:HB2	1.97	0.63
3:A:199:ARG:NH2	3:A:292:ARG:NH2	2.42	0.63
3:B:154:ARG:O	3:B:158:ILE:HG13	1.98	0.63
3:B:277:ASP:CB	3:B:284:LEU:HD13	2.28	0.63
3:A:184:ASP:HB3	3:A:196:HIS:O	1.98	0.63
3:A:60:ASN:C	3:A:61:ARG:HD3	2.23	0.63
3:A:212:ALA:O	3:A:213:LEU:HD23	1.99	0.63
3:A:159:ARG:HB2	3:A:224:TRP:CZ3	2.33	0.63
3:A:213:LEU:HD12	3:A:218:THR:CG2	2.29	0.63
1:C:13:DA:H5''	3:B:132:LYS:O	1.99	0.63
1:C:16:DG:OP1	3:B:315:TRP:HA	1.98	0.63
3:A:96:ASN:ND2	3:A:108:SER:OG	2.32	0.62
3:A:231:ALA:HB1	3:A:234:PRO:HG3	1.81	0.62
3:B:306:ILE:N	3:B:307:PRO:HD2	2.13	0.62
1:C:15:DT:H2''	1:C:16:DG:C8	2.35	0.62
3:A:21:ASP:OD1	3:A:24:ARG:NH2	2.33	0.62
3:A:236:ASN:HA	3:A:252:ALA:HA	1.80	0.62
3:B:319:ASN:OD1	3:B:319:ASN:N	2.31	0.62
3:B:89:GLN:HG2	3:B:117:MET:HE2	1.82	0.62
3:A:340:GLU:O	3:A:341:ASP:C	2.42	0.62
3:B:335:MET:HE3	3:B:335:MET:O	2.00	0.61
3:A:277:ASP:HB2	3:A:284:LEU:HB3	1.82	0.61
3:A:306:ILE:HD11	3:A:322:MET:HE1	1.81	0.61
3:A:187:ARG:O	3:A:189:ASP:N	2.32	0.61
3:A:166:ILE:HA	3:A:217:VAL:CG1	2.30	0.60
3:A:299:MET:HE3	3:B:335:MET:HA	1.82	0.60
3:A:305:SER:OG	3:A:307:PRO:HG2	2.01	0.60
3:A:297:ARG:HA	3:A:325:ILE:CG2	2.30	0.60
3:B:313:GLY:HA3	3:B:315:TRP:CZ3	2.36	0.60
3:A:295:ALA:O	3:A:299:MET:HB2	2.01	0.60
3:B:182:VAL:HB	3:B:234:PRO:HA	1.84	0.60
3:A:132:LYS:HB3	3:A:283:TYR:HD2	1.67	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:277:ASP:HB2	3:B:284:LEU:HD13	1.83	0.60
1:C:34:DT:H1'	3:A:244:LYS:HD3	1.83	0.59
3:A:293:VAL:O	3:A:297:ARG:HG3	2.02	0.59
1:C:11:DA:OP1	3:B:81:ARG:NH2	2.29	0.59
1:C:15:DT:H2''	1:C:16:DG:N7	2.16	0.59
3:A:184:ASP:O	3:A:196:HIS:N	2.33	0.59
3:A:313:GLY:HA3	3:A:315:TRP:CE3	2.38	0.59
2:D:16:DA:H2''	2:D:17:DT:C5	2.36	0.59
3:A:183:LYS:C	3:A:185:ILE:H	2.10	0.59
3:A:202:THR:HG22	3:A:203:LEU:N	2.14	0.59
3:A:228:SER:O	3:A:250:PRO:HG3	2.03	0.59
2:D:7:DT:O4	3:A:259:ARG:HG3	2.02	0.59
3:A:310:MET:HE1	3:A:318:VAL:N	2.17	0.59
3:B:310:MET:HE1	3:B:318:VAL:CG2	2.30	0.59
3:A:139:ARG:HG3	3:A:139:ARG:HH11	1.68	0.58
3:A:201:LYS:CA	3:B:130:ARG:HH21	2.15	0.58
3:A:178:ALA:HB2	3:A:261:LEU:HD11	1.86	0.58
3:B:128:GLY:O	3:B:130:ARG:NH1	2.36	0.58
2:D:30:DG:N7	4:D:56:HOH:O	2.32	0.58
3:A:189:ASP:H	3:A:194:LEU:CD1	2.16	0.58
3:B:334:ALA:O	3:B:337:ARG:HB2	2.04	0.58
3:A:310:MET:HE1	3:A:317:ASN:C	2.29	0.58
3:A:104:LEU:N	3:A:104:LEU:HD23	2.18	0.58
3:B:305:SER:O	3:B:308:GLU:N	2.34	0.58
2:D:23:DA:H2'	2:D:24:DT:C7	2.34	0.58
1:C:14:DA:N3	3:B:201:LYS:NZ	2.52	0.57
3:A:72:ARG:HG3	3:A:116:VAL:HB	1.85	0.57
3:A:259:ARG:NH2	3:A:262:GLU:HB3	2.18	0.57
3:A:217:VAL:O	3:A:221:VAL:HG23	2.05	0.57
3:B:262:GLU:O	3:B:266:GLU:HG3	2.03	0.57
3:A:174:ILE:HG23	4:A:394:HOH:O	2.04	0.57
3:A:308:GLU:HA	3:A:311:GLN:NE2	2.20	0.57
3:A:132:LYS:HB3	3:A:283:TYR:CD2	2.38	0.57
3:B:24:ARG:HB3	3:B:24:ARG:CZ	2.35	0.57
3:A:197:ILE:HD13	3:A:198:GLY:C	2.29	0.57
3:B:306:ILE:HG22	3:B:307:PRO:CD	2.34	0.57
3:A:263:GLY:HA2	3:A:266:GLU:OE1	2.04	0.57
1:C:2:DT:H2''	1:C:3:DA:C8	2.40	0.57
1:C:16:DG:H5''	3:B:316:THR:HG23	1.87	0.57
2:D:23:DA:H2'	2:D:24:DT:H72	1.85	0.57
3:A:310:MET:HE1	3:A:318:VAL:CA	2.35	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:107:PRO:O	3:A:110:SER:HB3	2.04	0.57
3:A:110:SER:O	3:A:114:SER:HB2	2.05	0.57
3:A:299:MET:HE1	3:B:335:MET:CB	2.34	0.57
3:A:84:ALA:O	3:A:88:ILE:HG13	2.04	0.56
3:A:172:LEU:HD11	3:A:197:ILE:HG13	1.87	0.56
3:B:130:ARG:HH11	3:B:130:ARG:CG	2.17	0.56
3:B:197:ILE:HB	3:B:209:VAL:CG2	2.35	0.56
3:B:281:GLN:O	3:B:284:LEU:HG	2.04	0.56
3:B:333:GLY:C	3:B:336:VAL:HG23	2.29	0.56
3:A:23:VAL:HG21	3:A:104:LEU:HD21	1.87	0.56
3:B:64:PHE:HA	3:B:65:PRO:C	2.30	0.56
2:D:6:DT:OP2	3:A:257:SER:HB3	2.06	0.56
3:A:197:ILE:HG22	3:A:209:VAL:HG22	1.85	0.56
3:A:317:ASN:OD1	3:A:319:ASN:HB2	2.05	0.56
3:A:178:ALA:HB2	3:A:261:LEU:CD1	2.36	0.56
1:C:14:DA:N6	4:C:65:HOH:O	2.37	0.56
3:A:197:ILE:HG23	3:A:209:VAL:HG13	1.88	0.56
3:A:306:ILE:HD11	3:A:322:MET:CE	2.36	0.56
3:B:293:VAL:HG22	3:B:324:TYR:CE1	2.41	0.56
3:A:146:ARG:HG2	3:A:161:LEU:HD21	1.87	0.56
3:A:336:VAL:O	3:A:340:GLU:HG3	2.06	0.56
3:A:197:ILE:CG1	3:A:198:GLY:H	2.18	0.56
3:A:279:SER:OG	3:A:281:GLN:HG2	2.06	0.56
3:B:181:ARG:NH1	3:B:235:ASN:O	2.37	0.55
1:C:10:DT:H71	3:B:43:LYS:HD3	1.88	0.55
3:A:300:ALA:HB2	3:A:325:ILE:HG21	1.87	0.55
3:A:308:GLU:HA	3:A:311:GLN:HE21	1.70	0.55
3:B:318:VAL:O	3:B:322:MET:HG2	2.07	0.55
1:C:27:DT:H2''	1:C:28:DA:O5'	2.06	0.55
3:B:293:VAL:HG22	3:B:324:TYR:CD1	2.41	0.55
2:D:2:DT:H2''	2:D:3:DA:C8	2.42	0.55
3:A:45:LEU:HD22	3:A:97:MET:HE2	1.88	0.55
3:A:187:ARG:CZ	3:A:193:MET:HG2	2.37	0.55
3:A:189:ASP:N	3:A:194:LEU:HD11	2.22	0.55
3:A:306:ILE:CD1	3:A:322:MET:HE1	2.37	0.55
3:A:27:LEU:HD23	3:A:30:MET:HE3	1.89	0.55
3:A:74:TYR:O	3:A:77:TYR:HB3	2.07	0.55
2:D:26:DC:C5'	3:B:282:ARG:HG3	2.37	0.54
3:A:146:ARG:O	3:A:150:GLU:HG3	2.07	0.54
3:A:276:LYS:HA	4:A:389:HOH:O	2.06	0.54
3:A:159:ARG:HB2	3:A:224:TRP:CE3	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:270:ARG:O	3:A:274:GLY:HA2	2.07	0.54
2:D:2:DT:O2	3:A:244:LYS:NZ	2.32	0.54
3:A:101:ARG:HH12	3:B:114:SER:CB	2.20	0.54
3:A:243:ARG:HB2	3:A:245:ASN:ND2	2.23	0.54
3:B:323:ASN:C	3:B:325:ILE:H	2.16	0.54
3:B:203:LEU:HD23	3:B:203:LEU:C	2.31	0.54
3:A:43:LYS:HE3	3:A:43:LYS:HA	1.90	0.54
3:B:172:LEU:HD21	3:B:197:ILE:HG13	1.89	0.54
1:C:1:DA:C1'	1:C:2:DT:H5''	2.37	0.53
3:B:130:ARG:HG2	3:B:130:ARG:NH1	2.18	0.53
3:A:243:ARG:HG3	4:A:362:HOH:O	2.08	0.53
3:A:325:ILE:O	3:A:328:LEU:HB2	2.09	0.53
2:D:5:DC:C2'	2:D:6:DT:H5'	2.38	0.53
3:A:39:GLU:O	3:A:43:LYS:HG2	2.07	0.53
3:A:230:VAL:HG13	3:A:231:ALA:N	2.23	0.53
3:A:166:ILE:N	3:A:217:VAL:HG13	2.23	0.53
3:A:171:LEU:HD11	3:B:335:MET:SD	2.49	0.53
3:A:171:LEU:HD12	3:B:335:MET:HG2	1.91	0.53
3:B:27:LEU:CD1	3:B:102:SER:HB3	2.39	0.53
3:A:121:ARG:HH11	3:A:121:ARG:CG	2.21	0.53
3:A:218:THR:O	3:A:221:VAL:HB	2.09	0.53
3:A:219:LYS:O	3:A:222:GLU:HB2	2.09	0.53
3:A:24:ARG:HG2	3:A:24:ARG:NH1	2.24	0.53
3:A:141:ASP:OD2	3:A:286:TRP:HZ2	1.92	0.53
3:A:245:ASN:ND2	3:A:247:VAL:H	2.07	0.53
3:B:141:ASP:O	3:B:145:VAL:HG23	2.09	0.53
3:A:310:MET:HE1	3:A:318:VAL:CG2	2.39	0.52
3:B:128:GLY:O	3:B:129:GLU:O	2.27	0.52
1:C:1:DA:H2''	1:C:2:DT:H5'	1.90	0.52
3:A:23:VAL:HG22	3:A:104:LEU:HD21	1.89	0.52
3:A:197:ILE:HG12	3:A:198:GLY:H	1.75	0.52
3:B:92:LEU:CD2	3:B:117:MET:HG2	2.39	0.52
3:B:318:VAL:O	3:B:318:VAL:HG12	2.09	0.52
1:C:26:DC:OP1	3:A:283:TYR:N	2.41	0.52
3:B:134:ALA:HA	3:B:283:TYR:CD1	2.45	0.52
3:A:229:GLY:O	3:A:230:VAL:HG12	2.10	0.52
1:C:7:DT:H1'	1:C:8:DA:H5'	1.92	0.52
3:B:281:GLN:CD	3:B:284:LEU:HD21	2.34	0.52
1:C:33:DA:H2''	1:C:34:DT:OP2	2.09	0.52
3:A:317:ASN:OD1	3:A:319:ASN:N	2.43	0.52
1:C:16:DG:H5''	3:B:316:THR:CG2	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:299:MET:CE	3:B:335:MET:HA	2.40	0.51
3:A:65:PRO:HG3	3:A:104:LEU:CD1	2.33	0.51
3:A:213:LEU:HD12	3:A:218:THR:HG22	1.92	0.51
3:B:310:MET:HE2	3:B:318:VAL:HA	1.92	0.51
3:B:310:MET:HE1	3:B:318:VAL:HG13	1.89	0.51
3:B:281:GLN:HG2	3:B:284:LEU:HD21	1.92	0.51
2:D:15:DC:H2'	2:D:16:DA:C8	2.45	0.51
3:A:182:VAL:HB	3:A:231:ALA:CB	2.28	0.51
3:A:189:ASP:H	3:A:194:LEU:HD11	1.75	0.51
2:D:17:DT:H3'	3:B:121:ARG:HD3	1.91	0.51
3:B:159:ARG:HB2	3:B:224:TRP:CZ3	2.46	0.51
3:B:204:VAL:HG12	3:B:204:VAL:O	2.11	0.51
3:A:57:LYS:HB2	3:A:57:LYS:NZ	2.26	0.51
3:A:204:VAL:O	3:A:205:SER:O	2.28	0.51
3:B:277:ASP:HB3	3:B:284:LEU:HD13	1.93	0.51
3:A:72:ARG:HG3	3:A:116:VAL:CB	2.42	0.50
3:B:276:LYS:O	3:B:277:ASP:O	2.29	0.50
3:A:170:THR:O	3:A:211:LYS:HE3	2.10	0.50
3:A:71:VAL:O	3:A:75:LEU:HG	2.12	0.50
3:B:188:THR:OG1	3:B:192:ARG:HB2	2.11	0.50
3:A:21:ASP:CG	3:A:24:ARG:HH22	2.18	0.50
3:B:170:THR:O	3:B:171:LEU:HB2	2.10	0.50
3:A:139:ARG:HG3	3:A:139:ARG:NH1	2.27	0.50
3:A:72:ARG:HD3	4:A:381:HOH:O	2.12	0.50
3:B:132:LYS:HB3	3:B:283:TYR:CD2	2.46	0.50
3:B:192:ARG:HG2	3:B:213:LEU:O	2.12	0.50
2:D:14:DG:C4	2:D:15:DC:C6	3.00	0.49
3:A:168:TYR:HD2	3:B:339:LEU:HD11	1.78	0.49
2:D:14:DG:C2'	2:D:15:DC:H6	2.17	0.49
3:A:161:LEU:HD22	3:A:220:LEU:HD13	1.94	0.49
3:A:177:ILE:C	3:A:180:ILE:HG23	2.38	0.49
3:A:243:ARG:HB2	3:A:245:ASN:HD21	1.77	0.49
2:D:14:DG:H3'	3:A:320:ILE:HG13	1.94	0.49
3:A:181:ARG:HA	3:A:237:TYR:HA	1.95	0.49
3:A:56:CYS:O	3:A:60:ASN:N	2.46	0.49
3:A:224:TRP:CE3	3:A:228:SER:HB3	2.47	0.49
3:A:166:ILE:O	3:A:167:ALA:C	2.56	0.49
3:A:202:THR:CG2	3:A:203:LEU:H	2.11	0.49
3:B:68:PRO:HB3	3:B:110:SER:OG	2.12	0.49
3:B:245:ASN:ND2	3:B:247:VAL:HG23	2.28	0.49
3:B:140:THR:O	3:B:144:GLN:HG3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:212:ALA:C	3:A:213:LEU:HD23	2.37	0.48
3:A:299:MET:O	3:A:304:VAL:HB	2.12	0.48
3:A:204:VAL:CG2	3:B:323:ASN:ND2	2.76	0.48
3:B:154:ARG:HB2	3:B:157:ASP:OD2	2.12	0.48
3:B:177:ILE:O	3:B:180:ILE:HG13	2.13	0.48
3:B:183:LYS:HB3	3:B:234:PRO:HB2	1.95	0.48
3:A:190:GLY:O	3:A:191:GLY:C	2.54	0.48
3:B:182:VAL:HG23	3:B:236:ASN:O	2.13	0.48
3:A:168:TYR:CD2	3:B:339:LEU:HD11	2.48	0.48
3:A:306:ILE:O	3:A:307:PRO:C	2.57	0.48
3:B:219:LYS:HB2	3:B:219:LYS:HE2	1.59	0.48
3:B:310:MET:HE2	3:B:318:VAL:HG13	1.96	0.48
3:A:20:SER:O	3:A:24:ARG:HB3	2.13	0.48
3:A:182:VAL:HG11	3:A:231:ALA:HB2	1.93	0.48
3:B:315:TRP:CZ2	3:B:324:TYR:HE2	2.32	0.48
3:B:319:ASN:HA	3:B:322:MET:HB2	1.94	0.48
3:A:187:ARG:HA	3:A:187:ARG:HD3	1.47	0.48
3:A:292:ARG:NH2	3:A:313:GLY:O	2.46	0.48
3:A:260:ALA:O	3:A:264:ILE:HG13	2.14	0.48
3:B:182:VAL:HG11	3:B:231:ALA:HA	1.96	0.48
3:A:174:ILE:HG13	3:A:175:ALA:N	2.29	0.47
3:B:171:LEU:O	3:B:292:ARG:HG3	2.14	0.47
3:A:273:TYR:CD1	3:A:273:TYR:N	2.81	0.47
3:A:170:THR:HG22	3:A:172:LEU:H	1.80	0.47
3:B:282:ARG:O	3:B:283:TYR:HB2	2.15	0.47
2:D:26:DC:H5''	3:B:282:ARG:HG3	1.94	0.47
3:A:277:ASP:O	3:A:284:LEU:HD13	2.13	0.47
3:B:78:LEU:O	3:B:81:ARG:HB2	2.15	0.47
1:C:16:DG:H5'	3:B:202:THR:CG2	2.44	0.47
3:B:336:VAL:O	3:B:340:GLU:HG2	2.14	0.47
3:A:24:ARG:O	3:A:28:MET:HG3	2.14	0.47
3:A:209:VAL:HG23	3:A:210:GLU:H	1.76	0.47
3:A:276:LYS:O	3:A:277:ASP:CB	2.63	0.47
3:B:281:GLN:NE2	3:B:284:LEU:HD21	2.30	0.47
1:C:13:DA:C2	2:D:23:DA:C2	3.03	0.47
3:A:223:ARG:O	3:A:227:VAL:HG22	2.15	0.47
3:A:300:ALA:CB	3:A:325:ILE:HG21	2.44	0.47
3:B:219:LYS:HE2	3:B:219:LYS:CA	2.40	0.47
3:B:318:VAL:O	3:B:322:MET:HB2	2.15	0.47
2:D:21:DT:C2'	2:D:22:DT:H5'	2.31	0.46
3:A:121:ARG:CG	3:A:121:ARG:NH1	2.78	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:263:GLY:HA2	3:A:266:GLU:CD	2.40	0.46
3:A:186:SER:C	3:A:187:ARG:HH11	2.23	0.46
3:A:259:ARG:HA	3:A:259:ARG:HE	1.77	0.46
3:B:128:GLY:O	3:B:130:ARG:HG2	2.16	0.46
2:D:17:DT:H3'	3:B:121:ARG:CD	2.46	0.46
3:A:209:VAL:CG2	3:A:210:GLU:N	2.79	0.46
3:A:224:TRP:CZ3	3:A:228:SER:HB3	2.50	0.46
3:A:265:PHE:CZ	3:A:291:ALA:HB2	2.51	0.46
2:D:5:DC:H2''	2:D:6:DT:H5'	1.98	0.46
3:B:224:TRP:CZ3	3:B:228:SER:HB3	2.51	0.46
3:B:325:ILE:C	3:B:326:ARG:O	2.56	0.46
3:A:34:ARG:C	3:A:36:ALA:H	2.23	0.46
3:A:307:PRO:O	3:A:311:GLN:HB3	2.15	0.46
3:A:25:LYS:NZ	3:B:69:GLU:OE2	2.49	0.45
3:A:61:ARG:HH11	3:A:61:ARG:CG	2.28	0.45
3:A:142:PHE:O	3:A:146:ARG:HB2	2.16	0.45
3:A:169:ASN:OD1	3:A:213:LEU:HB3	2.16	0.45
3:A:205:SER:HB3	3:A:206:THR:H	1.64	0.45
3:A:272:ILE:HG22	3:A:273:TYR:CE1	2.51	0.45
1:C:23:DA:C2	2:D:13:DA:C2	3.03	0.45
3:A:318:VAL:O	3:A:322:MET:HG2	2.16	0.45
3:B:245:ASN:ND2	3:B:247:VAL:CG2	2.79	0.45
3:B:277:ASP:HB3	3:B:279:SER:OG	2.16	0.45
3:B:245:ASN:CG	3:B:247:VAL:HG23	2.42	0.45
3:B:276:LYS:O	3:B:277:ASP:C	2.60	0.45
2:D:17:DT:H2''	2:D:18:DA:C8	2.51	0.45
3:A:176:GLU:OE1	3:A:199:ARG:HG3	2.17	0.45
3:A:272:ILE:CG2	3:A:273:TYR:CE1	2.99	0.45
3:B:306:ILE:N	3:B:307:PRO:CD	2.79	0.45
3:B:305:SER:O	3:B:306:ILE:C	2.58	0.45
3:A:133:GLN:NE2	3:A:324:TYR:CE1	2.79	0.45
3:A:197:ILE:CD1	3:A:198:GLY:H	2.29	0.45
3:B:203:LEU:HD23	3:B:204:VAL:CA	2.46	0.45
1:C:23:DA:C8	1:C:24:DT:H72	2.52	0.45
3:A:61:ARG:HD3	3:A:61:ARG:N	2.29	0.45
3:B:231:ALA:O	3:B:232:ASP:C	2.60	0.45
3:A:19:THR:OG1	3:A:22:GLU:N	2.38	0.45
3:A:166:ILE:HG22	3:A:217:VAL:HG12	1.98	0.45
3:A:213:LEU:CG	3:A:218:THR:HG23	2.46	0.45
3:A:214:SER:O	3:A:218:THR:OG1	2.30	0.45
3:B:281:GLN:O	3:B:281:GLN:HG2	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:21:DT:H2'	2:D:22:DT:C6	2.52	0.44
3:A:106:ARG:O	3:A:109:ASP:HB2	2.16	0.44
3:A:169:ASN:OD1	3:A:213:LEU:HA	2.16	0.44
3:A:289:HIS:O	3:A:290:SER:C	2.59	0.44
3:A:184:ASP:O	3:A:195:ILE:HA	2.17	0.44
3:A:306:ILE:N	3:A:307:PRO:CD	2.79	0.44
1:C:32:DT:C2'	1:C:33:DA:C8	2.98	0.44
3:A:183:LYS:H	3:A:183:LYS:HG3	1.51	0.44
3:A:281:GLN:HG3	3:A:284:LEU:HD21	1.99	0.44
1:C:14:DA:O5'	1:C:14:DA:H2'	2.17	0.44
3:A:171:LEU:HD13	3:A:312:ALA:HB1	2.00	0.44
3:A:197:ILE:CG1	3:A:198:GLY:N	2.81	0.44
3:A:249:ALA:O	3:A:250:PRO:C	2.60	0.44
3:B:197:ILE:HB	3:B:209:VAL:HG22	1.98	0.44
3:B:261:LEU:HD23	3:B:261:LEU:HA	1.81	0.44
1:C:9:DG:H2''	1:C:10:DT:C6	2.53	0.44
2:D:17:DT:H2''	2:D:18:DA:N7	2.32	0.44
3:A:135:LEU:O	3:A:286:TRP:HD1	2.00	0.44
3:A:171:LEU:HD11	3:A:295:ALA:HB1	2.00	0.44
3:B:336:VAL:C	3:B:340:GLU:HG3	2.43	0.44
1:C:24:DT:H2''	1:C:25:DA:C8	2.52	0.44
3:A:183:LYS:HG3	3:A:234:PRO:HB3	2.00	0.44
3:A:233:ASP:O	3:A:236:ASN:HB2	2.18	0.44
3:A:234:PRO:C	3:A:236:ASN:N	2.73	0.44
3:A:275:ALA:O	3:A:276:LYS:CG	2.63	0.44
3:A:187:ARG:NH2	3:A:193:MET:CE	2.76	0.43
3:B:218:THR:HG22	3:B:222:GLU:OE2	2.17	0.43
3:A:187:ARG:HD3	3:A:192:ARG:O	2.18	0.43
3:A:221:VAL:O	3:A:222:GLU:C	2.61	0.43
3:A:259:ARG:NH2	3:A:262:GLU:CB	2.81	0.43
3:A:214:SER:C	3:A:216:GLY:N	2.74	0.43
3:B:125:VAL:C	3:B:127:ALA:N	2.76	0.43
3:A:193:MET:CB	3:A:218:THR:HB	2.45	0.43
3:B:155:CYS:HB3	3:B:242:VAL:HG11	2.00	0.43
1:C:6:DT:H1'	1:C:7:DT:H5'	2.00	0.43
3:A:174:ILE:HG21	3:A:288:GLY:HA3	2.01	0.43
3:A:213:LEU:HD13	3:A:217:VAL:C	2.43	0.43
3:B:72:ARG:HG2	3:B:72:ARG:HH11	1.83	0.43
3:B:107:PRO:O	3:B:113:VAL:HB	2.19	0.43
3:A:243:ARG:HD2	4:A:362:HOH:O	2.19	0.43
3:A:274:GLY:C	3:A:276:LYS:H	2.27	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:34:ARG:HG2	3:B:42:TRP:CE2	2.53	0.43
3:B:179:ARG:O	3:B:180:ILE:C	2.61	0.43
1:C:16:DG:O5'	3:B:316:THR:OG1	2.29	0.42
3:A:299:MET:HE1	3:B:335:MET:CA	2.49	0.42
2:D:16:DA:H2''	2:D:17:DT:H71	1.97	0.42
3:B:132:LYS:HB3	3:B:283:TYR:CE2	2.54	0.42
3:B:204:VAL:O	3:B:205:SER:HB3	2.18	0.42
3:B:64:PHE:CZ	3:B:104:LEU:HD12	2.54	0.42
3:A:68:PRO:HB3	3:A:110:SER:OG	2.19	0.42
3:A:251:SER:HB3	4:A:393:HOH:O	2.18	0.42
3:A:239:PHE:HB3	3:A:256:LEU:HD12	2.01	0.42
3:B:237:TYR:CE1	3:B:255:GLN:HG2	2.54	0.42
1:C:24:DT:OP1	3:A:173:ARG:HD3	2.19	0.42
3:A:78:LEU:HB3	3:A:83:LEU:HD12	2.01	0.42
3:A:233:ASP:O	3:A:235:ASN:N	2.53	0.42
3:A:139:ARG:HD3	3:A:143:ASP:CG	2.44	0.42
3:A:269:HIS:ND1	4:A:389:HOH:O	2.36	0.42
3:B:125:VAL:C	3:B:127:ALA:H	2.28	0.42
3:A:170:THR:HG22	3:A:172:LEU:HB2	2.02	0.42
3:A:183:LYS:C	3:A:185:ILE:N	2.77	0.42
3:A:81:ARG:HB3	3:A:83:LEU:HG	2.01	0.42
3:A:171:LEU:HD21	3:A:295:ALA:HB3	2.02	0.42
3:B:249:ALA:N	3:B:250:PRO:HD3	2.35	0.42
3:A:134:ALA:O	3:A:136:ALA:N	2.54	0.41
3:B:122:LYS:HE3	3:B:122:LYS:HB3	1.62	0.41
3:B:244:LYS:C	3:B:246:GLY:H	2.28	0.41
3:A:64:PHE:HA	3:A:65:PRO:C	2.45	0.41
3:A:304:VAL:CG1	3:A:308:GLU:CB	2.96	0.41
3:B:101:ARG:HD3	4:B:439:HOH:O	2.20	0.41
3:B:310:MET:HG2	3:B:321:VAL:HG21	2.01	0.41
3:A:85:VAL:N	3:A:129:GLU:OE2	2.52	0.41
3:B:130:ARG:NH1	3:B:130:ARG:CG	2.79	0.41
3:B:238:LEU:HD12	3:B:238:LEU:O	2.19	0.41
3:B:278:ASP:OD1	3:B:278:ASP:N	2.51	0.41
3:A:132:LYS:HB2	3:A:283:TYR:HE2	1.84	0.41
3:A:270:ARG:NE	3:A:276:LYS:HD3	2.35	0.41
3:A:326:ARG:HG3	3:A:327:ASN:N	2.35	0.41
3:B:188:THR:O	3:B:189:ASP:C	2.63	0.41
3:B:313:GLY:HA3	3:B:315:TRP:CE3	2.56	0.41
3:A:159:ARG:NH1	3:A:240:CYS:O	2.53	0.41
3:A:230:VAL:HG22	3:A:231:ALA:H	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:15:DT:H4'	3:B:202:THR:CG2	2.51	0.41
3:A:138:GLU:HG2	3:A:294:GLY:HA2	2.02	0.41
3:B:193:MET:HE3	3:B:222:GLU:HG3	2.03	0.41
2:D:2:DT:H2''	2:D:3:DA:N7	2.35	0.41
2:D:14:DG:C2'	2:D:15:DC:C6	2.99	0.41
3:B:27:LEU:CD1	3:B:102:SER:CB	2.99	0.41
3:B:132:LYS:HB2	3:B:132:LYS:NZ	2.36	0.41
3:B:133:GLN:NE2	3:B:324:TYR:CD1	2.79	0.41
3:B:323:ASN:C	3:B:325:ILE:N	2.76	0.41
3:B:325:ILE:HG22	3:B:326:ARG:N	2.36	0.41
3:A:177:ILE:C	3:A:180:ILE:CG2	2.94	0.41
3:A:197:ILE:CD1	3:A:199:ARG:HB2	2.50	0.41
1:C:24:DT:H2''	1:C:25:DA:OP2	2.20	0.40
3:A:299:MET:CE	3:B:335:MET:CA	2.99	0.40
3:B:233:ASP:HA	3:B:234:PRO:HD3	1.85	0.40
3:B:334:ALA:O	3:B:336:VAL:N	2.54	0.40
1:C:15:DT:H71	4:C:57:HOH:O	2.21	0.40
3:A:61:ARG:CG	3:A:61:ARG:NH1	2.84	0.40
3:B:281:GLN:CG	3:B:284:LEU:HD21	2.51	0.40
3:A:235:ASN:HB2	3:A:252:ALA:CA	2.50	0.40
3:B:23:VAL:HG21	3:B:104:LEU:CD2	2.48	0.40
3:A:222:GLU:O	3:A:225:ILE:HB	2.21	0.40
3:B:74:TYR:O	3:B:77:TYR:HB3	2.22	0.40
3:B:197:ILE:HB	3:B:209:VAL:HG23	2.04	0.40
3:B:237:TYR:CZ	3:B:255:GLN:HG2	2.57	0.40
3:B:310:MET:SD	3:B:317:ASN:O	2.79	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	
3	A	321/349 (92%)	273 (85%)	37 (12%)	11 (3%)	3	2
3	B	313/349 (90%)	283 (90%)	22 (7%)	8 (3%)	4	3
All	All	634/698 (91%)	556 (88%)	59 (9%)	19 (3%)	3	2

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	188	THR
3	A	201	LYS
3	A	205	SER
3	A	230	VAL
3	B	129	GLU
3	B	277	ASP
3	B	324	TYR
3	B	337	ARG
3	A	276	LYS
3	A	279	SER
3	B	207	ALA
3	A	234	PRO
3	A	277	ASP
3	B	335	MET
3	A	184	ASP
3	A	275	ALA
3	B	325	ILE
3	A	236	ASN
3	B	180	ILE

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	
3	A	270/293 (92%)	192 (71%)	78 (29%)	0	0
3	B	264/293 (90%)	221 (84%)	43 (16%)	2	2
All	All	534/586 (91%)	413 (77%)	121 (23%)	1	1

All (121) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	19	THR
3	A	23	VAL
3	A	24	ARG
3	A	30	MET
3	A	39	GLU
3	A	43	LYS
3	A	46	LEU
3	A	57	LYS
3	A	60	ASN
3	A	61	ARG
3	A	79	GLN
3	A	81	ARG
3	A	97	MET
3	A	100	ARG
3	A	114	SER
3	A	116	VAL
3	A	121	ARG
3	A	122	LYS
3	A	129	GLU
3	A	132	LYS
3	A	139	ARG
3	A	140	THR
3	A	146	ARG
3	A	147	SER
3	A	152	SER
3	A	158	ILE
3	A	161	LEU
3	A	166	ILE
3	A	171	LEU
3	A	177	ILE
3	A	180	ILE
3	A	183	LYS
3	A	186	SER
3	A	187	ARG
3	A	189	ASP
3	A	195	ILE
3	A	197	ILE
3	A	200	THR
3	A	201	LYS
3	A	203	LEU
3	A	205	SER
3	A	206	THR

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Mol	Chain	Res	Type
3	A	209	VAL
3	A	213	LEU
3	A	214	SER
3	A	215	LEU
3	A	219	LYS
3	A	222	GLU
3	A	223	ARG
3	A	225	ILE
3	A	230	VAL
3	A	234	PRO
3	A	244	LYS
3	A	245	ASN
3	A	253	THR
3	A	255	GLN
3	A	256	LEU
3	A	257	SER
3	A	259	ARG
3	A	271	LEU
3	A	276	LYS
3	A	289	HIS
3	A	292	ARG
3	A	299	MET
3	A	301	ARG
3	A	306	ILE
3	A	308	GLU
3	A	309	ILE
3	A	311	GLN
3	A	316	THR
3	A	318	VAL
3	A	320	ILE
3	A	325	ILE
3	A	326	ARG
3	A	328	LEU
3	A	330	SER
3	A	335	MET
3	A	338	LEU
3	B	21	ASP
3	B	23	VAL
3	B	25	LYS
3	B	27	LEU
3	B	28	MET
3	B	46	LEU

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Mol	Chain	Res	Type
3	B	57	LYS
3	B	62	LYS
3	B	86	LYS
3	B	122	LYS
3	B	129	GLU
3	B	130	ARG
3	B	132	LYS
3	B	133	GLN
3	B	138	GLU
3	B	146	ARG
3	B	176	GLU
3	B	180	ILE
3	B	186	SER
3	B	193	MET
3	B	201	LYS
3	B	204	VAL
3	B	219	LYS
3	B	230	VAL
3	B	235	ASN
3	B	241	ARG
3	B	254	SER
3	B	272	ILE
3	B	276	LYS
3	B	278	ASP
3	B	279	SER
3	B	281	GLN
3	B	289	HIS
3	B	299	MET
3	B	306	ILE
3	B	308	GLU
3	B	319	ASN
3	B	320	ILE
3	B	321	VAL
3	B	325	ILE
3	B	336	VAL
3	B	337	ARG
3	B	338	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	40	HIS

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Mol	Chain	Res	Type
3	A	89	GLN
3	A	96	ASN
3	A	144	GLN
3	A	245	ASN
3	A	255	GLN
3	A	269	HIS
3	A	319	ASN
3	B	26	ASN
3	B	133	GLN
3	B	269	HIS
3	B	323	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	C	32/34 (94%)	-1.11	0	100 100	33, 51, 69, 93	0
2	D	34/34 (100%)	-1.05	0	100 100	33, 48, 83, 98	0
3	A	320/349 (91%)	-0.23	6 (1%)	66 68	28, 65, 92, 99	7 (2%)
3	B	317/349 (90%)	-0.74	1 (0%)	90 90	24, 43, 85, 100	3 (0%)
All	All	703/766 (91%)	-0.54	7 (0%)	79 80	24, 53, 90, 100	10 (1%)

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	231	ALA	3.6
3	B	333	GLY	3.0
3	A	229	GLY	2.5
3	A	208	GLY	2.3
3	A	196	HIS	2.2
3	A	199	ARG	2.1
3	A	204	VAL	2.1

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.