



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:07 PM UTC

PDB ID : 8R7Y / pdb_00008r7y
Title : Deoxyribonucleoside regulator DeoR in complex with the DNA operator
Authors : Pachl, P.; Soltysova, M.; Rezacova, P.
Deposited on : 2023-11-27
Resolution : 3.70 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:

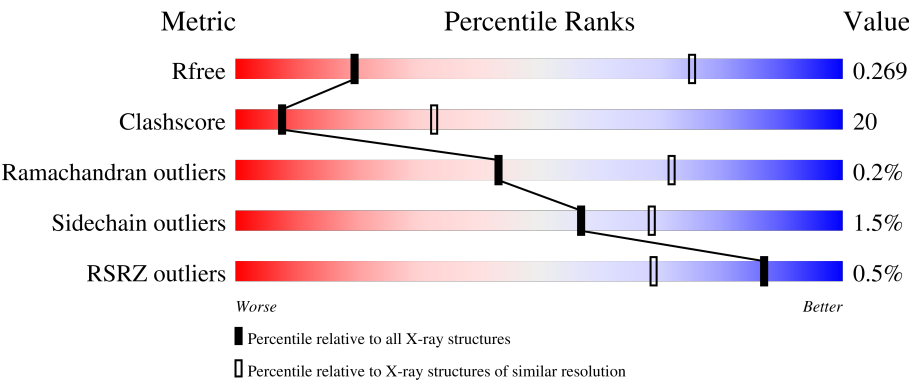
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.70 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.



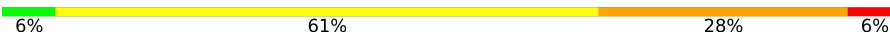
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	1131 (3.80-3.60)
Clashscore	190562	1171 (3.80-3.60)
Ramachandran outliers	187476	1129 (3.80-3.60)
Sidechain outliers	187428	1126 (3.80-3.60)
RSRZ outliers	180081	1130 (3.80-3.60)

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	A	318	% 64%32%..
1	B	318	% 73%24%..
1	C	318	% 62%34%..
1	D	318	66%31%..
2	E	18	17%39%44%

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Mol	Chain	Length	Quality of chain
2	H	18	
3	F	18	
3	G	18	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11264 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Deoxyribonucleoside regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	312	Total	C	N	O	S	0	0	0
			2454	1546	429	470	9			
1	B	311	Total	C	N	O	S	0	0	0
			2450	1547	428	466	9			
1	C	312	Total	C	N	O	S	0	0	0
			2435	1539	419	468	9			
1	D	312	Total	C	N	O	S	0	0	0
			2453	1547	428	470	8			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP P39140
A	-3	ILE	-	expression tag	UNP P39140
A	-2	ASP	-	expression tag	UNP P39140
A	-1	PRO	-	expression tag	UNP P39140
A	0	PHE	-	expression tag	UNP P39140
A	1	THR	-	expression tag	UNP P39140
B	-4	GLY	-	expression tag	UNP P39140
B	-3	ILE	-	expression tag	UNP P39140
B	-2	ASP	-	expression tag	UNP P39140
B	-1	PRO	-	expression tag	UNP P39140
B	0	PHE	-	expression tag	UNP P39140
B	1	THR	-	expression tag	UNP P39140
C	-4	GLY	-	expression tag	UNP P39140
C	-3	ILE	-	expression tag	UNP P39140
C	-2	ASP	-	expression tag	UNP P39140
C	-1	PRO	-	expression tag	UNP P39140
C	0	PHE	-	expression tag	UNP P39140
C	1	THR	-	expression tag	UNP P39140
D	-4	GLY	-	expression tag	UNP P39140
D	-3	ILE	-	expression tag	UNP P39140
D	-2	ASP	-	expression tag	UNP P39140

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	PRO	-	expression tag	UNP P39140
D	0	PHE	-	expression tag	UNP P39140
D	1	THR	-	expression tag	UNP P39140

- Molecule 2 is a DNA chain called OL18 DNA operator, strand 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	18	Total 366	C 179	N 61	O 109	P 17	0	0	0
2	H	18	Total 366	C 179	N 61	O 109	P 17	0	0	0

- Molecule 3 is a DNA chain called OL18 DNA operator, strand 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	18	Total 366	C 178	N 68	O 103	P 17	0	0	0
3	G	18	Total 366	C 178	N 68	O 103	P 17	0	0	0

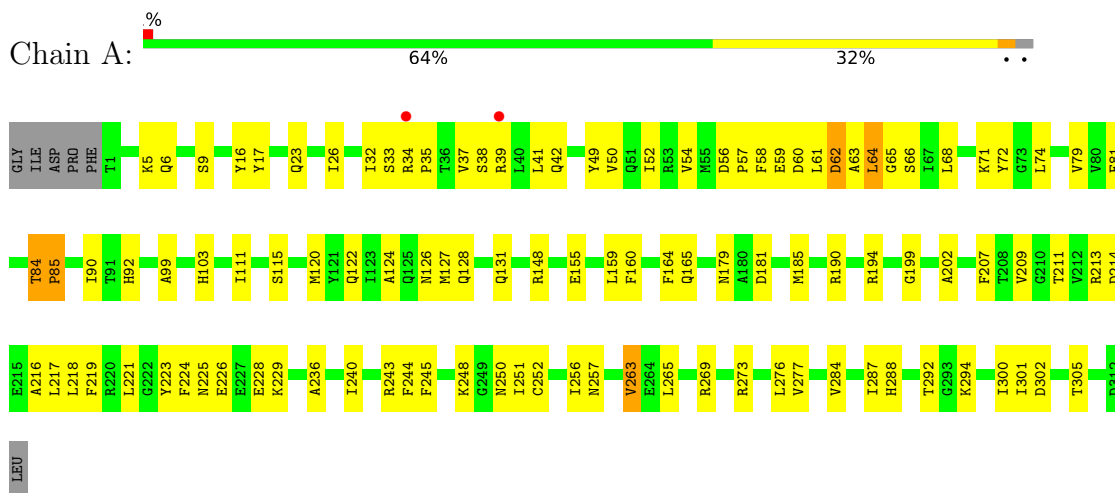
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total 2	O 2	0	0
4	B	1	Total 1	O 1	0	0
4	C	1	Total 1	O 1	0	0
4	E	1	Total 1	O 1	0	0
4	G	1	Total 1	O 1	0	0
4	H	2	Total 2	O 2	0	0

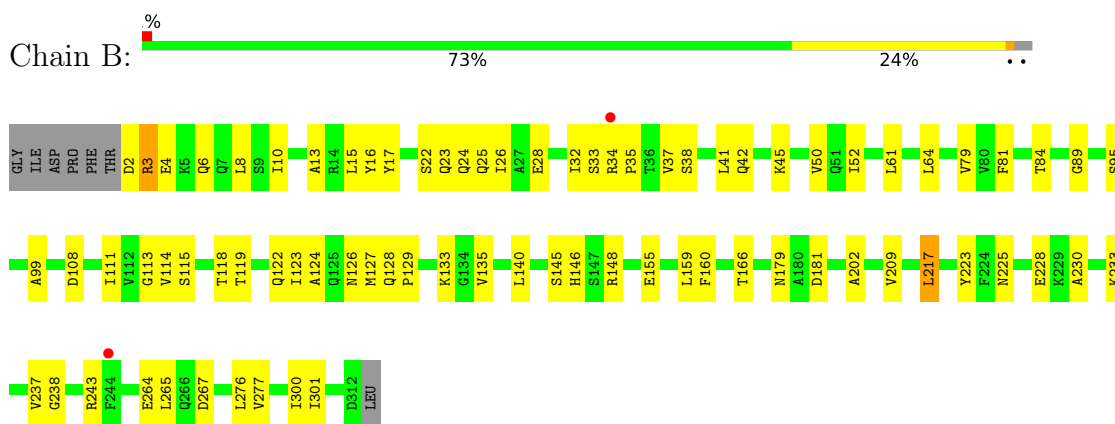
3 Residue-property plots [i](#)

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.

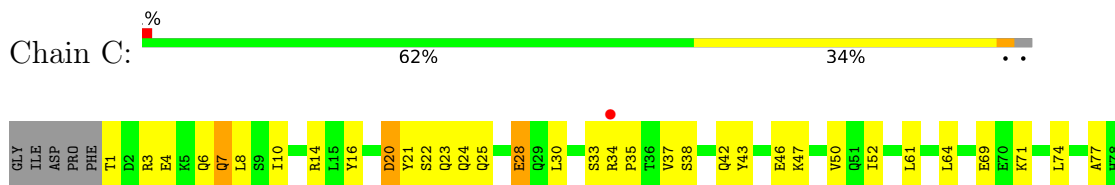
• Molecule 1: Deoxyribonucleoside regulator

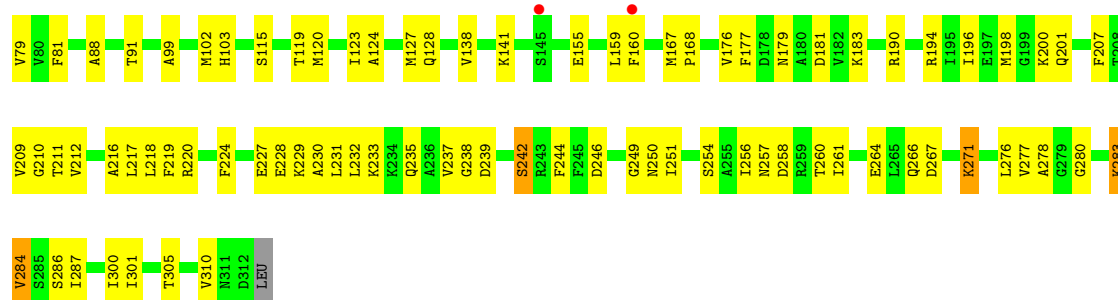


• Molecule 1: Deoxyribonucleoside regulator



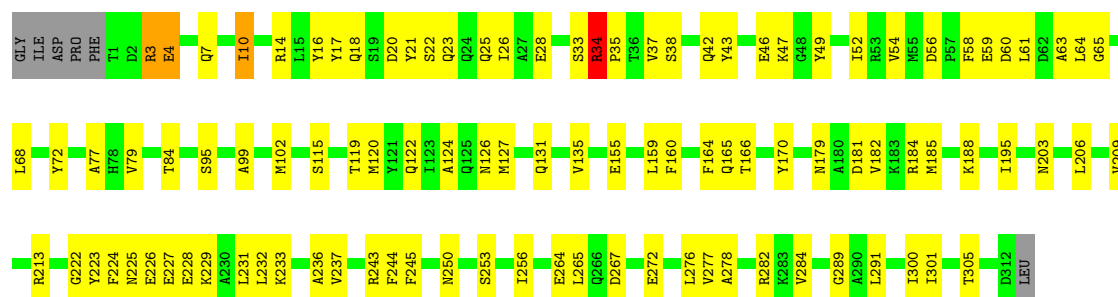
• Molecule 1: Deoxyribonucleoside regulator





- Molecule 1: Deoxyribonucleoside regulator

Chain D: 66% 31% ..



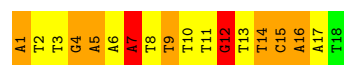
- Molecule 2: OL18 DNA operator, strand 1

Chain E: 17% 39% 44%



- Molecule 2: OL18 DNA operator, strand 1

Chain H: 6% 44% 39% 11%



- Molecule 3: OL18 DNA operator, strand 2

Chain F: 6% 61% 28% 6%



- Molecule 3: OL18 DNA operator, strand 2

Chain G: 11% 56% 33%



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	166.67Å 166.67Å 331.70Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.92 – 3.70 48.92 – 3.70	Depositor EDS
% Data completeness (in resolution range)	99.2 (48.92-3.70) 99.2 (48.92-3.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.01 (at 3.67Å)	Xtriage
Refinement program	REFMAC 5.8.0405	Depositor
R, R_{free}	0.266 , 0.333 (Not available) , 0.269	Depositor DCC
R_{free} test set	1482 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	149.4	Xtriage
Anisotropy	0.055	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 123.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11264	wwPDB-VP
Average B, all atoms (Å ²)	173.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.73% 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	A	0.44	0/2493	1.00	3/3367 (0.1%)
1	B	0.44	0/2490	1.01	1/3362 (0.0%)
1	C	0.44	0/2475	1.03	3/3347 (0.1%)
1	D	0.44	0/2493	1.02	1/3369 (0.0%)
2	E	0.41	0/409	1.24	0/630
2	H	0.41	0/409	1.35	6/630 (1.0%)
3	F	0.41	0/411	1.30	1/632 (0.2%)
3	G	0.42	0/411	1.30	1/632 (0.2%)
All	All	0.44	0/11591	1.07	16/15969 (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	D	0	1
2	E	0	8
2	H	0	7
3	F	0	10
3	G	0	10
All	All	0	36

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	17	DA	O3'-P-O5'	8.77	117.15	104.00
1	C	28	GLU	CA-C-N	7.88	130.68	120.44
1	C	28	GLU	C-N-CA	7.88	130.68	120.44
2	H	9	DT	P-O5'-C5'	-6.86	109.71	120.00
3	G	2	DT	C2'-C3'-O3'	-6.63	101.56	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	14	DT	P-O5'-C5'	-6.38	110.43	120.00
2	H	16	DA	P-O5'-C5'	-6.34	110.48	120.00
1	A	84	THR	CA-C-N	6.08	126.97	120.47
1	A	84	THR	C-N-CA	6.08	126.97	120.47
1	B	4	GLU	N-CA-C	-5.49	106.63	113.38
2	H	7	DA	O3'-P-O5'	-5.46	95.81	104.00
1	C	119	THR	CB-CA-C	-5.38	102.36	110.92
2	H	12	DG	O3'-P-O5'	-5.24	96.14	104.00
1	A	85	PRO	N-CA-C	5.19	123.57	114.75
1	D	119	THR	CB-CA-C	-5.11	102.79	110.92
2	H	15	DC	P-O3'-C3'	-5.11	112.54	120.20

There are no chirality outliers.

All (36) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	34	ARG	Sidechain
2	E	1	DA	Sidechain
2	E	12	DG	Sidechain
2	E	16	DA	Sidechain
2	E	17	DA	Sidechain
2	E	4	DG	Sidechain
2	E	5	DA	Sidechain
2	E	6	DA	Sidechain
2	E	7	DA	Sidechain
3	F	1	DA	Sidechain
3	F	10	DA	Sidechain
3	F	11	DA	Sidechain
3	F	16	DA	Sidechain
3	F	17	DA	Sidechain
3	F	4	DG	Sidechain
3	F	5	DA	Sidechain
3	F	6	DA	Sidechain
3	F	8	DA	Sidechain
3	F	9	DA	Sidechain
3	G	1	DA	Sidechain
3	G	10	DA	Sidechain
3	G	11	DA	Sidechain
3	G	16	DA	Sidechain
3	G	17	DA	Sidechain
3	G	4	DG	Sidechain
3	G	5	DA	Sidechain

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Mol	Chain	Res	Type	Group
3	G	6	DA	Sidechain
3	G	8	DA	Sidechain
3	G	9	DA	Sidechain
2	H	1	DA	Sidechain
2	H	12	DG	Sidechain
2	H	16	DA	Sidechain
2	H	4	DG	Sidechain
2	H	5	DA	Sidechain
2	H	6	DA	Sidechain
2	H	7	DA	Sidechain

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2454	0	2456	108	1
1	B	2450	0	2451	71	0
1	C	2435	0	2417	121	0
1	D	2453	0	2445	101	0
2	E	366	0	209	20	2
2	H	366	0	209	20	1
3	F	366	0	206	20	0
3	G	366	0	206	10	1
4	A	2	0	0	0	0
4	B	1	0	0	1	0
4	C	1	0	0	0	0
4	E	1	0	0	0	0
4	G	1	0	0	0	0
4	H	2	0	0	0	0
All	All	11264	0	10599	438	3

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 20.

All (438) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:23:GLN:OE1	1:C:38:SER:OG	1.69	1.08
1:B:111:ILE:HG22	1:B:202:ALA:HA	1.37	1.02
1:C:52:ILE:CG2	1:D:52:ILE:HD12	1.90	1.02
1:B:23:GLN:OE1	1:B:38:SER:OG	1.76	1.01
1:C:8:LEU:HD11	1:C:30:LEU:HD22	1.42	1.01
1:C:43:TYR:OH	1:C:47:LYS:HE2	1.61	1.00
1:B:111:ILE:CG2	1:B:202:ALA:HA	1.92	0.99
1:A:16:TYR:HB2	1:A:26:ILE:HD11	1.45	0.99
1:C:212:VAL:HG21	1:C:237:VAL:HA	1.42	0.98
1:D:209:VAL:HG13	1:D:276:LEU:HD11	1.49	0.94
1:A:251:ILE:HG23	1:A:257:ASN:HD22	1.28	0.94
1:C:14:ARG:HH22	1:D:60:ASP:HB2	1.33	0.93
1:A:217:LEU:O	1:A:221:LEU:HD13	1.68	0.92
1:D:72:TYR:CD2	1:D:291:LEU:HB3	2.05	0.92
2:E:2:DT:H2''	2:E:3:DT:C5'	2.01	0.91
1:C:228:GLU:O	1:C:231:LEU:HG	1.71	0.91
1:A:16:TYR:HB2	1:A:26:ILE:CD1	2.01	0.89
1:B:145:SER:HB3	1:B:223:TYR:OH	1.71	0.88
1:C:52:ILE:HG22	1:D:52:ILE:HD12	1.55	0.88
1:B:22:SER:HB3	1:B:25:GLN:HG3	1.55	0.87
1:C:23:GLN:CD	1:C:38:SER:OG	2.17	0.86
1:C:14:ARG:NH2	1:D:56:ASP:OD2	2.08	0.86
1:C:198:MET:HA	1:C:201:GLN:HB2	1.56	0.85
1:A:23:GLN:OE1	1:A:38:SER:OG	1.97	0.83
1:B:217:LEU:HD12	1:B:217:LEU:H	1.44	0.81
1:A:111:ILE:CG2	1:A:202:ALA:HA	2.10	0.81
2:E:2:DT:H2''	2:E:3:DT:H5'	1.61	0.81
1:C:251:ILE:HB	1:C:257:ASN:ND2	1.94	0.81
1:C:284:VAL:HG22	1:C:305:THR:HG23	1.61	0.81
1:D:72:TYR:HD2	1:D:291:LEU:HB3	1.46	0.79
1:A:23:GLN:CD	1:A:38:SER:OG	2.25	0.79
1:C:138:VAL:HB	1:C:198:MET:HE2	1.63	0.78
1:A:111:ILE:HG22	1:A:202:ALA:HA	1.65	0.78
1:D:253:SER:HB3	1:D:256:ILE:HG12	1.63	0.78
1:C:99:ALA:HA	1:C:127:MET:SD	2.23	0.78
1:C:1:THR:HB	1:C:4:GLU:HB2	1.63	0.77
1:B:111:ILE:HG22	1:B:202:ALA:CA	2.14	0.77
1:C:176:VAL:HG21	1:C:224:PHE:HE1	1.51	0.76
1:C:212:VAL:HG12	1:C:233:LYS:HA	1.68	0.76
1:A:251:ILE:HG23	1:A:257:ASN:ND2	2.01	0.74
2:E:2:DT:H2''	2:E:3:DT:H5''	1.67	0.74
1:C:7:GLN:HB2	1:D:58:PHE:CZ	2.22	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:16:TYR:O	1:D:20:ASP:HA	1.88	0.74
1:D:95:SER:HB3	1:D:122:GLN:HG3	1.68	0.74
1:D:95:SER:CB	1:D:122:GLN:HG3	2.18	0.74
1:C:211:THR:HG23	1:C:216:ALA:HB2	1.70	0.74
1:C:176:VAL:HG21	1:C:224:PHE:CE1	2.24	0.73
1:A:56:ASP:HB2	1:A:57:PRO:HD2	1.69	0.73
1:B:16:TYR:HB2	1:B:26:ILE:HD11	1.70	0.73
2:E:1:DA:H2"	2:E:2:DT:H5"	1.70	0.72
1:A:16:TYR:CB	1:A:26:ILE:HD11	2.18	0.72
1:C:212:VAL:CG2	1:C:237:VAL:HA	2.19	0.71
1:D:99:ALA:HA	1:D:127:MET:SD	2.29	0.71
1:A:68:LEU:HD21	1:A:72:TYR:HD2	1.54	0.71
1:A:288:HIS:O	1:A:292:THR:HG23	1.90	0.71
1:A:99:ALA:HA	1:A:127:MET:SD	2.30	0.71
1:C:50:VAL:HG12	1:D:54:VAL:HG22	1.73	0.70
1:B:155:GLU:O	1:B:159:LEU:HG	1.91	0.70
2:E:10:DT:C6	2:E:11:DT:H72	2.27	0.69
1:C:61:LEU:HD11	1:C:81:PHE:HB2	1.72	0.69
1:C:8:LEU:HD11	1:C:30:LEU:CD2	2.22	0.69
1:C:16:TYR:O	1:C:20:ASP:HA	1.93	0.69
1:B:145:SER:CB	1:B:223:TYR:OH	2.41	0.69
1:D:155:GLU:O	1:D:159:LEU:HG	1.92	0.69
1:B:217:LEU:HD12	1:B:217:LEU:N	2.08	0.68
1:B:24:GLN:O	1:B:28:GLU:HG2	1.93	0.68
1:C:231:LEU:O	1:C:235:GLN:HB2	1.94	0.68
1:D:61:LEU:H	1:D:64:LEU:HD13	1.57	0.68
1:B:230:ALA:HA	1:B:233:LYS:HE2	1.74	0.68
1:A:81:PHE:HZ	1:B:15:LEU:HD23	1.59	0.68
1:C:224:PHE:O	1:C:229:LYS:HE2	1.93	0.68
1:B:146:HIS:NE2	1:B:148:ARG:NH1	2.42	0.67
1:C:52:ILE:CG2	1:D:52:ILE:CD1	2.70	0.67
1:A:217:LEU:O	1:A:221:LEU:CD1	2.40	0.67
1:D:224:PHE:CZ	1:D:256:ILE:HB	2.29	0.67
1:A:84:THR:OG1	1:A:85:PRO:HD2	1.94	0.67
1:C:231:LEU:HD12	1:C:232:LEU:N	2.10	0.67
1:D:182:VAL:HA	1:D:185:MET:HE3	1.75	0.67
1:C:155:GLU:O	1:C:159:LEU:HG	1.93	0.67
1:B:217:LEU:H	1:B:217:LEU:CD1	2.08	0.66
1:C:230:ALA:O	1:C:233:LYS:HG2	1.95	0.66
1:A:6:GLN:OE1	1:A:49:TYR:OH	2.13	0.66
1:A:236:ALA:HA	1:A:245:PHE:O	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:14:ARG:NH2	1:D:60:ASP:HB2	2.09	0.66
1:D:47:LYS:HB2	1:D:49:TYR:CE2	2.31	0.65
3:G:1:DA:C8	3:G:2:DT:H72	2.31	0.65
1:B:84:THR:CG2	1:B:89:GLY:HA3	2.27	0.65
1:B:225:ASN:HB2	1:B:228:GLU:HB2	1.79	0.65
1:A:155:GLU:O	1:A:159:LEU:HG	1.95	0.65
1:C:52:ILE:HG22	1:D:52:ILE:CD1	2.27	0.65
3:F:4:DG:H2''	3:F:5:DA:C8	2.31	0.65
1:C:209:VAL:HA	1:C:239:ASP:O	1.97	0.65
3:F:10:DA:H5'	3:F:10:DA:H8	1.62	0.65
1:B:24:GLN:NE2	1:B:28:GLU:OE2	2.25	0.64
1:A:226:GLU:HA	1:A:229:LYS:CB	2.28	0.64
2:E:4:DG:H2''	2:E:5:DA:H8	1.61	0.64
1:C:278:ALA:HB1	1:C:283:LYS:HE3	1.80	0.64
3:F:15:DC:H2''	3:F:16:DA:H5''	1.79	0.64
1:A:209:VAL:HG22	1:A:240:ILE:HG12	1.79	0.64
1:C:52:ILE:HD13	1:D:17:TYR:HB3	1.80	0.64
1:C:251:ILE:HB	1:C:257:ASN:HD21	1.61	0.64
1:C:228:GLU:HA	1:C:231:LEU:CD2	2.28	0.64
1:A:54:VAL:HG12	1:B:50:VAL:HG22	1.80	0.63
1:B:23:GLN:CD	1:B:38:SER:OG	2.40	0.63
1:A:218:LEU:HG	1:A:224:PHE:CE2	2.34	0.63
3:G:14:DT:H1'	3:G:15:DC:H5'	1.79	0.63
1:A:52:ILE:HG12	1:B:52:ILE:HG22	1.81	0.62
1:C:7:GLN:HB2	1:D:58:PHE:CE2	2.34	0.62
1:D:22:SER:HB3	1:D:25:GLN:HG3	1.82	0.62
1:A:23:GLN:CD	1:A:38:SER:HG	2.05	0.62
1:B:95:SER:HA	1:B:123:ILE:HG13	1.82	0.62
1:D:224:PHE:CE2	1:D:256:ILE:HD12	2.35	0.62
3:F:12:DT:H2'	3:F:13:DT:H71	1.82	0.62
2:H:3:DT:H2''	2:H:4:DG:C8	2.35	0.62
3:F:10:DA:H5'	3:F:10:DA:C8	2.34	0.62
1:A:273:ARG:HG3	1:A:273:ARG:HH11	1.65	0.61
1:D:225:ASN:O	1:D:228:GLU:HB2	2.00	0.61
2:H:9:DT:H2'	2:H:10:DT:C7	2.30	0.61
1:C:211:THR:CG2	1:C:216:ALA:HB2	2.29	0.61
1:D:182:VAL:HA	1:D:185:MET:CE	2.30	0.61
2:E:4:DG:H2''	2:E:5:DA:C8	2.35	0.61
1:B:3:ARG:HD3	1:B:6:GLN:CD	2.25	0.61
1:A:218:LEU:HG	1:A:224:PHE:HE2	1.66	0.61
1:C:284:VAL:CG2	1:C:305:THR:HG23	2.29	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:GLU:O	1:A:64:LEU:HB2	2.01	0.61
1:A:207:PHE:CE1	1:A:240:ILE:HD13	2.35	0.60
2:E:1:DA:C4	2:E:2:DT:C5	2.89	0.60
3:F:11:DA:H2'	3:F:12:DT:H72	1.82	0.60
1:A:81:PHE:HZ	1:B:15:LEU:CD2	2.15	0.60
1:B:126:ASN:O	1:B:128:GLN:HG2	2.00	0.60
1:D:65:GLY:HA2	1:D:77:ALA:HB3	1.82	0.60
1:A:23:GLN:HE21	1:A:41:LEU:HD12	1.67	0.60
1:D:47:LYS:HB2	1:D:49:TYR:CD2	2.37	0.60
1:B:111:ILE:CG2	1:B:202:ALA:CA	2.75	0.60
1:A:216:ALA:O	1:A:219:PHE:HB2	2.02	0.60
2:H:4:DG:H2''	2:H:5:DA:C8	2.37	0.60
1:C:24:GLN:NE2	1:C:28:GLU:OE2	2.34	0.59
1:D:203:ASN:HB2	1:D:272:GLU:OE1	2.02	0.59
1:B:64:LEU:HD23	1:B:79:VAL:HG11	1.83	0.59
1:C:264:GLU:HB3	1:C:266:GLN:OE1	2.02	0.59
1:A:226:GLU:HA	1:A:229:LYS:HB3	1.82	0.59
1:C:283:LYS:O	1:C:287:ILE:HG13	2.02	0.59
1:D:170:TYR:O	1:D:195:ILE:HD13	2.02	0.59
1:C:61:LEU:CD1	1:C:81:PHE:HB2	2.32	0.58
3:F:16:DA:C2	3:F:17:DA:C4	2.91	0.58
1:C:52:ILE:HG21	1:D:52:ILE:HD12	1.83	0.58
1:C:190:ARG:HH11	1:C:194:ARG:HH22	1.50	0.58
1:A:72:TYR:HE1	1:A:292:THR:HG22	1.69	0.58
1:B:99:ALA:HA	1:B:128:GLN:OE1	2.04	0.58
1:C:22:SER:O	1:C:23:GLN:C	2.45	0.58
2:E:16:DA:H2''	2:E:17:DA:C8	2.39	0.58
1:A:190:ARG:HH11	1:A:194:ARG:HH22	1.52	0.58
1:A:269:ARG:HG2	1:A:294:LYS:O	2.05	0.57
1:B:3:ARG:HD3	1:B:6:GLN:HG3	1.86	0.57
1:C:42:GLN:O	1:C:46:GLU:HB2	2.04	0.57
1:C:43:TYR:CE1	1:C:47:LYS:HD3	2.40	0.57
1:B:99:ALA:N	1:B:123:ILE:HG23	2.20	0.57
1:C:123:ILE:O	1:C:127:MET:HG2	2.04	0.56
1:C:251:ILE:HD12	1:C:254:SER:HA	1.86	0.56
3:G:17:DA:H2'	3:G:18:DT:H71	1.88	0.56
1:A:218:LEU:HD23	1:A:244:PHE:HZ	1.70	0.56
1:D:277:VAL:HG22	1:D:300:ILE:HB	1.86	0.56
1:D:60:ASP:O	1:D:61:LEU:HB2	2.05	0.56
1:D:229:LYS:HG2	1:D:233:LYS:HE3	1.88	0.55
1:A:111:ILE:HG22	1:A:202:ALA:CA	2.34	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:ALA:C	1:A:65:GLY:N	2.61	0.55
1:A:63:ALA:O	1:A:64:LEU:C	2.49	0.55
1:D:47:LYS:CB	1:D:49:TYR:CE2	2.90	0.55
2:H:9:DT:H2'	2:H:10:DT:H72	1.89	0.55
1:A:16:TYR:CA	1:A:26:ILE:HD11	2.36	0.54
1:D:3:ARG:O	1:D:4:GLU:C	2.50	0.54
2:E:4:DG:C2	2:E:5:DA:C4	2.95	0.54
1:B:13:ALA:HA	1:B:41:LEU:CD2	2.37	0.54
1:A:42:GLN:HA	1:A:42:GLN:OE1	2.08	0.54
1:A:243:ARG:NH2	1:A:263:VAL:O	2.40	0.54
1:D:232:LEU:O	1:D:236:ALA:HB3	2.07	0.54
1:D:213:ARG:HH22	1:D:282:ARG:NH1	2.06	0.54
1:A:284:VAL:HG22	1:A:305:THR:HG23	1.89	0.53
1:D:95:SER:CB	1:D:122:GLN:CG	2.86	0.53
1:A:81:PHE:CZ	1:B:15:LEU:HD23	2.40	0.53
1:D:23:GLN:HG2	3:F:11:DA:OP2	2.08	0.53
1:D:99:ALA:HB1	1:D:126:ASN:O	2.09	0.53
1:C:14:ARG:HH21	1:D:56:ASP:CG	2.13	0.53
3:G:13:DT:H2''	3:G:14:DT:H72	1.91	0.53
1:A:35:PRO:CB	1:A:39:ARG:HH12	2.22	0.53
2:E:3:DT:H2''	2:E:4:DG:C8	2.44	0.53
1:D:72:TYR:CD2	1:D:291:LEU:CB	2.85	0.53
1:C:3:ARG:O	1:C:6:GLN:HB3	2.09	0.53
1:C:238:GLY:O	1:C:244:PHE:HA	2.09	0.53
1:C:228:GLU:HA	1:C:231:LEU:HD23	1.91	0.52
1:A:228:GLU:HB3	1:A:256:ILE:HD11	1.91	0.52
1:C:183:LYS:HE3	1:C:258:ASP:O	2.10	0.52
1:A:209:VAL:HG22	1:A:240:ILE:CG1	2.38	0.52
1:C:42:GLN:OE1	1:C:42:GLN:HA	2.08	0.52
1:D:131:GLN:OE1	1:D:131:GLN:HA	2.08	0.52
1:A:72:TYR:HB3	1:A:74:LEU:HD21	1.92	0.52
3:F:3:DT:H2''	3:F:4:DG:C8	2.44	0.52
1:C:280:GLY:O	1:C:284:VAL:HG23	2.09	0.52
1:D:203:ASN:HB2	1:D:272:GLU:CD	2.34	0.52
1:D:229:LYS:HG2	1:D:233:LYS:CE	2.40	0.52
1:D:42:GLN:HA	1:D:42:GLN:OE1	2.09	0.52
1:D:227:GLU:O	1:D:231:LEU:HD13	2.10	0.52
2:H:14:DT:C2'	2:H:15:DC:C6	2.93	0.52
1:A:211:THR:HG22	1:A:214:ASP:CG	2.36	0.51
1:C:23:GLN:HE22	2:E:11:DT:H3'	1.75	0.51
1:A:16:TYR:CD1	1:A:16:TYR:C	2.88	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:ALA:O	1:A:66:SER:N	2.44	0.51
2:H:14:DT:H2''	2:H:15:DC:C6	2.45	0.51
1:C:1:THR:HG21	1:C:4:GLU:HG2	1.92	0.51
1:D:34:ARG:NH2	3:F:12:DT:O4	2.44	0.51
1:B:61:LEU:HD11	1:B:81:PHE:HB2	1.93	0.51
1:C:246:ASP:OD2	1:C:250:ASN:HB2	2.10	0.51
1:A:211:THR:HG23	1:A:214:ASP:H	1.75	0.51
1:D:42:GLN:O	1:D:46:GLU:HB2	2.11	0.51
1:C:231:LEU:HD12	1:C:231:LEU:C	2.35	0.51
3:F:15:DC:C2'	3:F:16:DA:H5''	2.40	0.51
1:C:102:MET:HB2	1:C:127:MET:HE1	1.93	0.51
3:G:8:DA:H2''	3:G:9:DA:C8	2.46	0.51
1:C:69:GLU:HG2	1:C:77:ALA:HB3	1.92	0.51
1:D:34:ARG:N	1:D:35:PRO:CD	2.74	0.51
1:C:6:GLN:HG2	1:D:58:PHE:CZ	2.46	0.50
2:H:3:DT:H2''	2:H:4:DG:H8	1.75	0.50
1:B:3:ARG:HD3	1:B:6:GLN:CG	2.41	0.50
1:C:10:ILE:O	1:C:14:ARG:HG3	2.11	0.50
1:A:209:VAL:HG23	1:A:276:LEU:HD11	1.92	0.50
1:D:243:ARG:HD2	1:D:265:LEU:HG	1.92	0.50
1:C:21:TYR:HA	1:C:25:GLN:OE1	2.11	0.50
1:D:232:LEU:HD11	1:D:244:PHE:CG	2.46	0.50
1:B:42:GLN:OE1	1:B:42:GLN:HA	2.11	0.50
1:B:277:VAL:HG22	1:B:300:ILE:HB	1.93	0.50
1:C:52:ILE:HG22	1:D:52:ILE:CG1	2.42	0.50
1:B:3:ARG:HE	1:B:3:ARG:HA	1.75	0.50
3:F:13:DT:H2''	3:F:14:DT:H72	1.93	0.50
1:D:227:GLU:HG3	1:D:228:GLU:N	2.25	0.50
1:D:179:ASN:OD1	1:D:181:ASP:HB2	2.12	0.49
1:D:264:GLU:O	1:D:267:ASP:HB2	2.12	0.49
2:H:10:DT:C6	2:H:11:DT:H72	2.47	0.49
1:A:226:GLU:HA	1:A:229:LYS:HB2	1.94	0.49
1:A:23:GLN:NE2	1:A:38:SER:OG	2.44	0.49
1:B:34:ARG:N	1:B:35:PRO:CD	2.74	0.49
1:D:43:TYR:O	1:D:46:GLU:HB3	2.12	0.49
1:D:209:VAL:HG22	1:D:278:ALA:CB	2.41	0.49
1:D:16:TYR:HB2	1:D:26:ILE:HD11	1.93	0.49
1:A:68:LEU:HA	1:A:71:LYS:HB3	1.95	0.49
1:C:71:LYS:HE3	1:C:310:VAL:HA	1.94	0.49
1:C:200:LYS:C	1:C:271:LYS:HZ1	2.20	0.49
1:A:34:ARG:N	1:A:35:PRO:CD	2.75	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:LEU:HD23	1:A:244:PHE:CZ	2.47	0.49
1:C:34:ARG:N	1:C:35:PRO:CD	2.76	0.49
1:A:35:PRO:HB2	1:A:39:ARG:HH12	1.77	0.49
1:A:273:ARG:HG3	1:A:273:ARG:NH1	2.27	0.49
1:C:179:ASN:OD1	1:C:181:ASP:HB2	2.13	0.49
2:H:1:DA:H1'	2:H:2:DT:H5'	1.93	0.49
1:B:32:ILE:HD11	1:B:37:VAL:HG22	1.93	0.48
1:B:108:ASP:CG	1:B:133:LYS:HG2	2.38	0.48
1:A:248:LYS:HB3	1:A:250:ASN:HD21	1.79	0.48
1:A:115:SER:H	1:A:120:MET:HE2	1.79	0.48
1:A:103:HIS:ND1	1:A:128:GLN:CB	2.77	0.48
1:A:218:LEU:HD12	1:A:223:TYR:HD2	1.79	0.48
1:A:111:ILE:CB	1:A:202:ALA:HA	2.43	0.48
1:B:16:TYR:HB2	1:B:26:ILE:CD1	2.39	0.48
1:C:200:LYS:HE2	1:C:267:ASP:OD2	2.14	0.48
1:A:277:VAL:HG22	1:A:300:ILE:HB	1.95	0.47
1:C:69:GLU:HA	1:C:74:LEU:HB2	1.96	0.47
3:G:9:DA:H1'	3:G:10:DA:H5'	1.95	0.47
1:A:5:LYS:O	1:A:9:SER:HB3	2.14	0.47
1:A:17:TYR:CE2	1:A:50:VAL:HG11	2.48	0.47
1:C:183:LYS:CE	1:C:258:ASP:O	2.62	0.47
1:D:47:LYS:HG3	1:D:49:TYR:HE2	1.79	0.47
1:C:22:SER:C	1:C:24:GLN:N	2.68	0.47
1:D:95:SER:OG	1:D:122:GLN:CG	2.62	0.47
1:D:115:SER:H	1:D:120:MET:HE2	1.78	0.47
1:A:63:ALA:O	1:A:65:GLY:N	2.48	0.47
1:D:95:SER:OG	1:D:122:GLN:HG2	2.14	0.47
1:B:179:ASN:OD1	1:B:181:ASP:HB2	2.15	0.47
1:A:111:ILE:CG2	1:A:202:ALA:CA	2.88	0.47
1:C:64:LEU:HD23	1:C:79:VAL:HG11	1.97	0.47
1:D:23:GLN:HE22	1:D:38:SER:HB3	1.78	0.47
1:D:124:ALA:HB2	1:D:160:PHE:CE2	2.49	0.47
1:A:131:GLN:OE1	1:A:131:GLN:HA	2.15	0.47
1:D:102:MET:HE2	1:D:206:LEU:HD22	1.96	0.47
1:D:209:VAL:HG13	1:D:276:LEU:CD1	2.34	0.47
1:C:61:LEU:HD22	1:C:79:VAL:HG12	1.96	0.47
1:D:43:TYR:OH	1:D:47:LYS:HE2	2.14	0.47
1:B:128:GLN:N	1:B:129:PRO:HD3	2.31	0.46
1:C:103:HIS:ND1	1:C:128:GLN:HB2	2.30	0.46
1:B:124:ALA:HB2	1:B:160:PHE:CE2	2.51	0.46
1:A:99:ALA:HB1	1:A:126:ASN:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:115:SER:H	1:C:120:MET:HE2	1.81	0.46
1:D:23:GLN:NE2	1:D:38:SER:HB3	2.31	0.46
3:F:18:DT:H5''	2:H:1:DA:H4'	1.97	0.46
1:C:138:VAL:HB	1:C:198:MET:CE	2.41	0.46
2:H:2:DT:C2'	2:H:3:DT:H71	2.46	0.46
1:B:2:ASP:O	1:B:3:ARG:NE	2.48	0.46
1:B:264:GLU:O	1:B:267:ASP:HB2	2.15	0.46
1:C:1:THR:HG22	1:C:3:ARG:H	1.81	0.46
1:A:61:LEU:O	1:A:62:ASP:HB2	2.15	0.46
1:A:207:PHE:CE1	1:A:276:LEU:HD13	2.51	0.46
1:A:17:TYR:HB3	1:B:52:ILE:HD12	1.97	0.45
1:A:179:ASN:OD1	1:A:181:ASP:HB2	2.16	0.45
1:A:244:PHE:HB2	1:A:252:CYS:SG	2.56	0.45
1:B:119:THR:O	1:B:123:ILE:HD12	2.16	0.45
3:G:13:DT:H2''	3:G:14:DT:C7	2.46	0.45
1:A:226:GLU:C	1:A:228:GLU:N	2.72	0.45
1:C:24:GLN:O	1:C:28:GLU:HG2	2.15	0.45
1:B:243:ARG:HD2	1:B:265:LEU:HG	1.98	0.45
1:C:207:PHE:O	1:C:276:LEU:HA	2.17	0.45
1:C:212:VAL:CG1	1:C:233:LYS:HA	2.43	0.45
1:C:102:MET:CB	1:C:127:MET:HE1	2.47	0.45
1:C:246:ASP:OD1	1:C:249:GLY:N	2.50	0.45
2:E:9:DT:C6	2:E:10:DT:H72	2.51	0.45
1:A:209:VAL:HG23	1:A:276:LEU:CD1	2.46	0.45
1:B:209:VAL:HG13	1:B:276:LEU:CD1	2.47	0.45
1:A:111:ILE:HB	1:A:202:ALA:HA	1.97	0.45
1:A:103:HIS:ND1	1:A:128:GLN:HB2	2.32	0.45
1:B:118:THR:O	1:B:122:GLN:HG2	2.17	0.45
1:D:184:ARG:O	1:D:188:LYS:HE3	2.17	0.45
1:A:79:VAL:HA	1:A:301:ILE:O	2.17	0.44
1:D:22:SER:HB3	1:D:25:GLN:CG	2.45	0.44
1:D:64:LEU:HD12	1:D:64:LEU:H	1.81	0.44
1:D:79:VAL:HA	1:D:301:ILE:O	2.16	0.44
2:E:3:DT:H2''	2:E:4:DG:H8	1.80	0.44
1:A:224:PHE:O	1:A:225:ASN:C	2.60	0.44
1:C:227:GLU:HG2	1:C:228:GLU:N	2.33	0.44
2:H:1:DA:H2'	2:H:2:DT:H71	1.99	0.44
1:C:23:GLN:NE2	2:E:11:DT:H3'	2.32	0.44
1:C:217:LEU:HA	1:C:220:ARG:CB	2.47	0.44
1:C:43:TYR:CE2	1:C:47:LYS:HG3	2.53	0.44
1:C:138:VAL:CB	1:C:198:MET:HE2	2.42	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:251:ILE:CD1	1:C:254:SER:HA	2.46	0.44
1:B:23:GLN:NE2	2:H:11:DT:H3'	2.32	0.44
1:C:124:ALA:HB2	1:C:160:PHE:CE2	2.52	0.44
1:D:72:TYR:CE2	1:D:291:LEU:CB	3.01	0.44
1:A:218:LEU:HD12	1:A:223:TYR:CD2	2.53	0.44
1:B:17:TYR:CE2	1:B:50:VAL:HG11	2.53	0.44
3:F:15:DC:H2''	3:F:16:DA:C5'	2.47	0.44
3:F:11:DA:H2'	3:F:12:DT:C7	2.47	0.44
1:A:32:ILE:HD11	1:A:37:VAL:HG22	2.00	0.43
1:D:34:ARG:HG3	1:D:34:ARG:HH11	1.82	0.43
2:E:6:DA:C6	2:E:7:DA:C6	3.05	0.43
1:A:56:ASP:OD2	1:A:60:ASP:HB2	2.18	0.43
1:A:58:PHE:O	1:A:61:LEU:O	2.36	0.43
1:A:243:ARG:HD2	1:A:265:LEU:HG	1.99	0.43
1:D:135:VAL:O	1:D:166:THR:HB	2.18	0.43
1:B:45:LYS:NZ	4:B:401:HOH:O	2.51	0.43
3:G:2:DT:C2'	3:G:3:DT:H72	2.47	0.43
1:B:135:VAL:O	1:B:166:THR:HB	2.19	0.43
1:A:251:ILE:HD12	1:A:251:ILE:H	1.84	0.43
1:C:79:VAL:HA	1:C:301:ILE:O	2.18	0.43
1:C:209:VAL:HG12	1:C:278:ALA:CB	2.49	0.43
2:E:10:DT:C2	2:E:11:DT:C5	3.07	0.43
2:E:12:DG:H1	3:F:7:DC:H42	1.67	0.43
3:G:8:DA:C2	2:H:12:DG:C2	3.07	0.43
1:D:224:PHE:CE1	1:D:256:ILE:HB	2.52	0.43
1:A:72:TYR:OH	1:A:288:HIS:ND1	2.52	0.43
1:B:111:ILE:HG22	1:B:202:ALA:CB	2.49	0.43
1:D:7:GLN:O	1:D:10:ILE:HG13	2.19	0.43
1:D:226:GLU:O	1:D:229:LYS:HB3	2.19	0.43
1:C:209:VAL:CG2	1:C:286:SER:HB3	2.49	0.43
1:D:43:TYR:OH	1:D:47:LYS:CE	2.66	0.43
2:E:4:DG:C2	2:E:5:DA:C5	3.07	0.43
3:F:11:DA:C2'	3:F:12:DT:H72	2.48	0.43
1:A:90:ILE:HG23	1:A:302:ASP:HB3	2.01	0.43
1:D:33:SER:O	1:D:37:VAL:HG23	2.19	0.43
1:D:72:TYR:CE2	1:D:291:LEU:HB2	2.54	0.43
1:A:61:LEU:O	1:A:62:ASP:CB	2.66	0.42
1:B:128:GLN:N	1:B:129:PRO:CD	2.82	0.42
1:C:22:SER:O	1:C:25:GLN:N	2.52	0.42
1:A:124:ALA:HB2	1:A:160:PHE:CE2	2.54	0.42
1:B:79:VAL:HA	1:B:301:ILE:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:13:DT:C2'	2:H:14:DT:H72	2.49	0.42
2:H:2:DT:H2'	2:H:3:DT:H71	2.01	0.42
1:B:16:TYR:O	1:B:16:TYR:CG	2.72	0.42
1:C:16:TYR:O	1:C:20:ASP:CA	2.65	0.42
1:C:50:VAL:CG1	1:D:54:VAL:HG22	2.44	0.42
1:C:210:GLY:N	1:C:239:ASP:HB3	2.34	0.42
1:A:23:GLN:NE2	1:A:38:SER:HG	2.18	0.42
1:C:6:GLN:HG2	1:D:58:PHE:CE1	2.55	0.42
1:D:25:GLN:O	1:D:28:GLU:HB2	2.20	0.42
1:A:23:GLN:NE2	1:A:38:SER:HA	2.35	0.42
1:B:209:VAL:HG13	1:B:276:LEU:HD13	2.01	0.42
1:D:33:SER:HB2	1:D:35:PRO:HD2	2.02	0.42
1:D:52:ILE:HG23	1:D:52:ILE:O	2.20	0.42
1:B:146:HIS:CD2	1:B:148:ARG:NH1	2.88	0.42
1:C:33:SER:O	1:C:37:VAL:HG23	2.20	0.42
1:C:43:TYR:CZ	1:C:47:LYS:HE2	2.48	0.42
1:D:284:VAL:HG22	1:D:305:THR:HG23	2.00	0.42
1:B:113:GLY:HA3	1:B:140:LEU:HD11	2.01	0.41
2:H:12:DG:C8	2:H:13:DT:H72	2.55	0.41
1:C:196:ILE:O	1:C:200:LYS:HG3	2.21	0.41
1:C:277:VAL:HG22	1:C:300:ILE:HB	2.01	0.41
1:D:14:ARG:O	1:D:18:GLN:HB2	2.20	0.41
1:A:224:PHE:O	1:A:226:GLU:N	2.51	0.41
1:C:210:GLY:H	1:C:239:ASP:HB3	1.85	0.41
1:A:199:GLY:HA3	1:A:263:VAL:HG21	2.02	0.41
1:B:114:VAL:HG12	1:B:115:SER:N	2.36	0.41
3:F:13:DT:C2'	3:F:14:DT:H72	2.50	0.41
1:D:23:GLN:HE21	1:D:34:ARG:HD3	1.86	0.41
1:A:92:HIS:ND1	1:A:122:GLN:OE1	2.49	0.41
1:A:287:ILE:H	1:A:287:ILE:HG13	1.73	0.41
1:B:24:GLN:O	1:B:24:GLN:HG3	2.20	0.41
1:D:222:GLY:O	1:D:223:TYR:C	2.62	0.41
1:B:22:SER:CB	1:B:25:GLN:HG3	2.38	0.41
1:C:88:ALA:O	1:C:91:THR:HB	2.21	0.41
1:C:167:MET:HB3	1:C:168:PRO:CD	2.51	0.41
1:D:237:VAL:HG21	1:D:289:GLY:HA3	2.03	0.41
1:A:103:HIS:CE1	1:A:128:GLN:HB3	2.56	0.41
1:C:33:SER:HB2	1:C:35:PRO:HD2	2.03	0.41
1:C:212:VAL:HA	1:C:219:PHE:HE2	1.86	0.41
1:C:232:LEU:HD21	1:C:256:ILE:HD12	2.02	0.41
1:A:33:SER:O	1:A:37:VAL:HG23	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:13:ALA:HA	1:B:41:LEU:HD22	2.02	0.41
1:C:3:ARG:O	1:C:6:GLN:N	2.54	0.41
1:C:218:LEU:HD23	1:C:244:PHE:CZ	2.56	0.41
1:C:264:GLU:O	1:C:267:ASP:HB2	2.20	0.41
1:C:283:LYS:O	1:C:286:SER:HB2	2.21	0.41
1:D:65:GLY:H	1:D:68:LEU:HB3	1.86	0.41
1:D:164:PHE:O	1:D:165:GLN:C	2.64	0.41
1:D:245:PHE:HA	1:D:250:ASN:O	2.21	0.41
2:E:12:DG:C4	2:E:13:DT:C5	3.09	0.41
3:F:18:DT:C5'	2:H:1:DA:H4'	2.51	0.41
1:B:33:SER:HB2	1:B:35:PRO:HD2	2.02	0.40
1:B:127:MET:C	1:B:129:PRO:HD3	2.46	0.40
3:G:1:DA:C2	3:G:2:DT:C2	3.10	0.40
1:A:185:MET:HE3	1:A:185:MET:HB2	1.88	0.40
1:D:59:GLU:CB	1:D:63:ALA:HB2	2.52	0.40
1:C:33:SER:HB3	3:F:4:DG:O5'	2.21	0.40
1:C:141:LYS:HE3	1:C:242:SER:OG	2.22	0.40
1:C:177:PHE:CD2	1:C:261:ILE:HD11	2.56	0.40
1:D:253:SER:CB	1:D:256:ILE:HG12	2.45	0.40
1:A:92:HIS:HA	1:A:122:GLN:OE1	2.22	0.40
1:A:164:PHE:O	1:A:165:GLN:C	2.65	0.40
1:A:223:TYR:C	1:A:225:ASN:H	2.30	0.40
1:B:8:LEU:HA	1:B:8:LEU:HD12	1.79	0.40
1:B:237:VAL:HG23	1:B:238:GLY:N	2.37	0.40
1:A:148:ARG:HD3	1:A:148:ARG:HA	1.32	0.40
2:H:1:DA:C4	2:H:2:DT:C6	3.09	0.40
2:H:7:DA:H2'	2:H:8:DT:H72	2.02	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:ARG:NH2	2:E:16:DA:C5'[12_554]	1.54	0.66
3:G:1:DA:O5'	2:H:17:DA:OP2[11_454]	1.60	0.60
2:E:18:DT:O2	2:E:18:DT:O2[10_445]	2.15	0.05

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	310/318 (98%)	300 (97%)	8 (3%)	2 (1%)	21	52
1	B	309/318 (97%)	300 (97%)	9 (3%)	0	100	100
1	C	310/318 (98%)	299 (96%)	11 (4%)	0	100	100
1	D	310/318 (98%)	299 (96%)	10 (3%)	1 (0%)	36	65
All	All	1239/1272 (97%)	1198 (97%)	38 (3%)	3 (0%)	43	72

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	62	ASP
1	A	64	LEU
1	D	4	GLU

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	263/271 (97%)	262 (100%)	1 (0%)	84	81
1	B	262/271 (97%)	259 (99%)	3 (1%)	65	73
1	C	259/271 (96%)	252 (97%)	7 (3%)	39	59
1	D	262/271 (97%)	257 (98%)	5 (2%)	50	65
All	All	1046/1084 (96%)	1030 (98%)	16 (2%)	57	68

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	263	VAL
1	B	3	ARG
1	B	10	ILE
1	B	217	LEU
1	C	7	GLN
1	C	20	ASP
1	C	242	SER
1	C	260	THR
1	C	271	LYS
1	C	283	LYS
1	C	284	VAL
1	D	3	ARG
1	D	10	ILE
1	D	21	TYR
1	D	34	ARG
1	D	84	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	250	ASN
1	A	257	ASN
1	B	6	GLN
1	B	165	GLN
1	C	51	GLN
1	C	257	ASN
1	D	23	GLN
1	D	125	GLN
1	D	146	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [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	A	312/318 (98%)	-0.48	2 (0%) 85 66	107, 176, 231, 265	0
1	B	311/318 (97%)	-0.51	2 (0%) 85 66	96, 165, 222, 250	0
1	C	312/318 (98%)	-0.43	3 (0%) 79 55	98, 194, 242, 275	0
1	D	312/318 (98%)	-0.48	0 100 100	96, 160, 209, 268	0
2	E	18/18 (100%)	-0.60	0 100 100	131, 158, 183, 192	0
2	H	18/18 (100%)	-0.26	0 100 100	138, 160, 243, 243	0
3	F	18/18 (100%)	-0.58	0 100 100	124, 150, 206, 207	0
3	G	18/18 (100%)	-0.69	0 100 100	111, 150, 192, 206	0
All	All	1319/1344 (98%)	-0.48	7 (0%) 87 69	96, 173, 232, 275	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	34	ARG	2.9
1	C	145	SER	2.4
1	C	160	PHE	2.3
1	A	34	ARG	2.3
1	A	39	ARG	2.3
1	B	34	ARG	2.3
1	B	244	PHE	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 [i](#)

There are no such residues in this entry.