



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 31, 2021 – 09:09 AM EDT

PDB ID : 3FBD  
Title : Crystal structure of the nuclease domain of COLE7(D493Q mutant) in complex with an 18-BP duplex DNA  
Authors : Wang, Y.T.; Doudeva, L.G.; Yuan, H.S.  
Deposited on : 2008-11-19  
Resolution : 2.90 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](mailto: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.

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

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.23.2  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.23.2

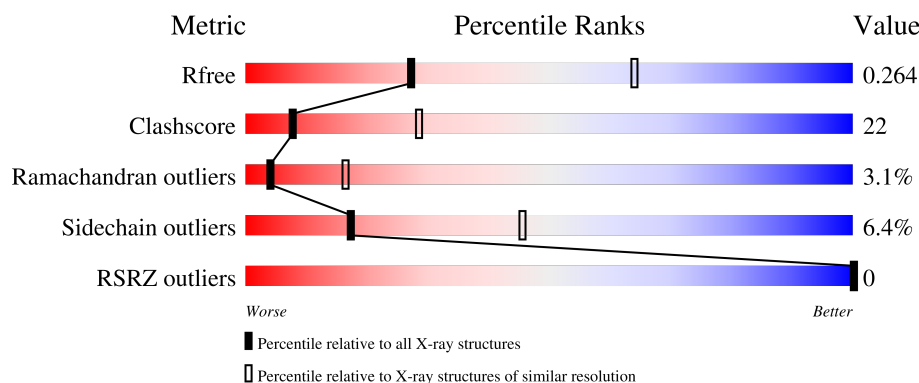
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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))
$R_{free}$	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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  $\geq 3$ , 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	132	<div> <div>55%</div> <div>40%</div> <div>5%</div> </div>
1	D	132	<div> <div>59%</div> <div>37%</div> <div>..</div> </div>
2	B	18	<div> <div>28%</div> <div>61%</div> <div>11%</div> </div>
2	C	18	<div> <div>44%</div> <div>50%</div> <div>6%</div> </div>
2	E	18	<div> <div>56%</div> <div>28%</div> <div>17%</div> </div>

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Mol	Chain	Length	Quality of chain
2	F	18	<div><div></div><div>28%</div><div>67%</div><div>6%</div></div>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3663 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 Colicin-E7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	132	Total	C	N	O	S	0	0	0
			1067	661	208	196	2			
1	D	131	Total	C	N	O	S	0	0	0
			1061	658	207	194	2			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	493	GLN	ASP	engineered mutation	UNP Q47112
D	493	GLN	ASP	engineered mutation	UNP Q47112

- Molecule 2 is a DNA chain called 5'-D(\*DGP\*DGP\*DAP\*DAP\*DTP\*DTP\*DCP\*DGP\*DAP\*DTP\*DCP\*DGP\*DAP\*DAP\*DTP\*DTP\*DCP\*DC)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	18	Total	C	N	O	P	0	0	0
			366	176	67	106	17			
2	C	18	Total	C	N	O	P	0	0	0
			366	176	67	106	17			
2	E	18	Total	C	N	O	P	0	0	0
			366	176	67	106	17			
2	F	18	Total	C	N	O	P	0	0	0
			366	176	67	106	17			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	13	Total	O	0	0
			13	13		
3	B	11	Total	O	0	0
			11	11		

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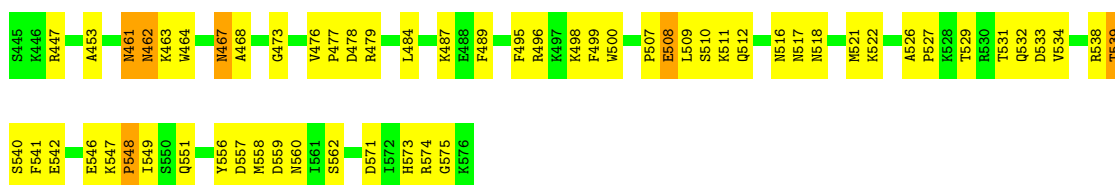
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	8	Total 8	O 8	0	0
3	D	18	Total 18	O 18	0	0
3	E	7	Total 7	O 7	0	0
3	F	14	Total 14	O 14	0	0

### 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.

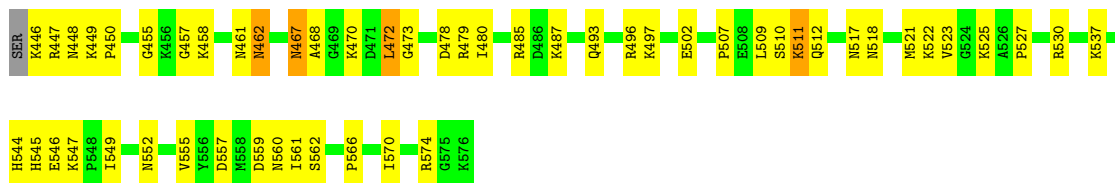
#### • Molecule 1: Colicin-E7

Chain A: 

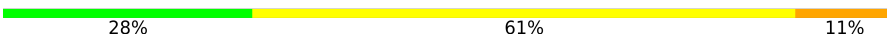


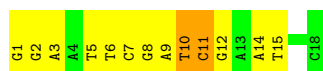
#### • Molecule 1: Colicin-E7

Chain D: 



#### • Molecule 2: 5'-D(\*DGP\*DGP\*DAP\*DAP\*DTP\*DTP\*DCP\*DGP\*DAP\*DTP\*DCP\*DGP\*DAP\*DAP\*DTP\*DTP\*DCP\*DC)-3'

Chain B: 



#### • Molecule 2: 5'-D(\*DGP\*DGP\*DAP\*DAP\*DTP\*DTP\*DCP\*DGP\*DAP\*DTP\*DCP\*DGP\*DAP\*DAP\*DTP\*DTP\*DCP\*DC)-3'

Chain C: 




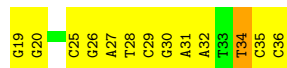
#### • Molecule 2: 5'-D(\*DGP\*DGP\*DAP\*DAP\*DTP\*DTP\*DCP\*DGP\*DAP\*DTP\*DCP\*DGP\*DAP\*DAP\*DTP\*DTP\*DCP\*DC)-3'

Chain E:  56% 28% 17%



● Molecule 2: 5'-D(\*DGP\*DGP\*DAP\*DAP\*DTP\*DTP\*DCP\*DGP\*DAP\*DTP\*DCP\*DGP\*DAP\*DAP\*DTP\*DTP\*DCP\*DC)-3'

Chain F:  28% 67% 6%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	60.59Å 49.42Å 92.40Å 90.00° 102.64° 90.00°	Depositor
Resolution (Å)	27.20 – 2.90 37.92 – 2.89	Depositor EDS
% Data completeness (in resolution range)	93.9 (27.20-2.90) 93.2 (37.92-2.89)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.11 (at 2.90Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.201 , 0.264 0.201 , 0.264	Depositor DCC
$R_{free}$ test set	1166 reflections (10.07%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	41.5	Xtriage
Anisotropy	0.919	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 45.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	3663	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.35% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.58	0/1089	0.72	0/1453
1	D	0.62	0/1083	0.73	0/1445
2	B	0.65	0/410	0.80	0/631
2	C	0.60	0/410	0.73	0/631
2	E	0.74	0/410	0.94	0/631
2	F	0.68	0/410	0.89	0/631
All	All	0.63	0/3812	0.78	0/5422

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
2	B	0	2
2	C	0	1
2	E	0	5
2	F	0	1
All	All	0	9

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	10	DT	Sidechain
2	B	11	DC	Sidechain
2	C	34	DT	Sidechain
2	E	10	DT	Sidechain
2	E	11	DC	Sidechain

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Mol	Chain	Res	Type	Group
2	E	15	DT	Sidechain
2	E	7	DC	Sidechain
2	E	9	DA	Sidechain
2	F	34	DT	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	1067	0	1078	47	0
1	D	1061	0	1073	42	0
2	B	366	0	205	23	0
2	C	366	0	205	12	0
2	E	366	0	205	7	0
2	F	366	0	205	15	0
3	A	13	0	0	0	0
3	B	11	0	0	4	0
3	C	8	0	0	3	0
3	D	18	0	0	3	0
3	E	7	0	0	4	0
3	F	14	0	0	7	0
All	All	3663	0	2971	141	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 22.

All (141) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:31:DA:H3'	3:F:7:HOH:O	1.45	1.12
1:D:549:ILE:HD12	1:D:555:VAL:HG22	1.31	1.08
2:C:26:DG:H2''	2:C:27:DA:H5'	1.42	0.99
2:B:10:DT:H2''	2:B:11:DC:H5''	1.44	0.98
1:A:510:SER:HB2	1:A:518:ASN:HD21	1.31	0.91
1:D:547:LYS:H	1:D:560:ASN:HD21	1.14	0.90
2:B:9:DA:H2'	3:B:161:HOH:O	1.74	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:547:LYS:H	1:A:560:ASN:HD21	1.19	0.86
2:E:7:DC:H3'	3:E:168:HOH:O	1.76	0.85
1:A:507:PRO:O	1:A:511:LYS:HG3	1.81	0.81
2:F:26:DG:H5''	3:F:2:HOH:O	1.81	0.80
2:C:26:DG:H2''	2:C:27:DA:C5'	2.16	0.75
1:A:467:ASN:H	1:A:467:ASN:HD22	1.34	0.73
1:A:521:MET:SD	1:A:527:PRO:HG3	2.29	0.73
1:D:457:GLY:O	1:D:458:LYS:HG3	1.89	0.72
1:A:478:ASP:OD1	1:A:479:ARG:N	2.22	0.72
1:A:547:LYS:N	1:A:560:ASN:HD21	1.90	0.70
1:A:453:ALA:HB2	1:A:489:PHE:HE2	1.55	0.69
1:A:573:HIS:CE1	2:B:10:DT:H4'	2.29	0.68
2:F:25:DC:H5'	3:F:119:HOH:O	1.93	0.68
1:A:479:ARG:NH1	1:A:508:GLU:HB2	2.10	0.67
1:A:479:ARG:HH12	1:A:508:GLU:CB	2.07	0.67
2:B:9:DA:H2''	2:B:10:DT:H5'	1.76	0.66
2:C:25:DC:H2''	2:C:26:DG:H5''	1.77	0.66
1:A:547:LYS:H	1:A:560:ASN:ND2	1.93	0.65
2:E:11:DC:H3'	3:E:151:HOH:O	1.96	0.65
2:B:2:DG:H2'	2:B:2:DG:OP2	1.96	0.65
1:A:509:LEU:O	1:A:512:GLN:HG3	1.98	0.63
1:D:521:MET:SD	1:D:527:PRO:HG3	2.39	0.62
1:D:446:LYS:N	3:D:190:HOH:O	2.32	0.62
1:A:529:THR:HG22	1:A:539:THR:O	2.00	0.62
3:D:116:HOH:O	2:F:29:DC:H4'	2.01	0.61
2:F:28:DT:H2''	2:F:29:DC:H5'	1.82	0.60
1:D:546:GLU:HB3	1:D:560:ASN:ND2	2.17	0.59
2:E:5:DT:H2''	2:E:6:DT:OP2	2.01	0.59
1:A:461:ASN:HB3	1:A:463:LYS:H	1.66	0.59
2:F:29:DC:H2''	2:F:30:DG:C8	2.38	0.59
1:A:529:THR:CG2	1:A:534:VAL:HG12	2.33	0.58
1:A:447:ARG:NH1	2:B:11:DC:OP2	2.36	0.58
2:C:26:DG:C5'	3:C:122:HOH:O	2.50	0.58
1:D:478:ASP:OD2	1:D:479:ARG:N	2.38	0.57
1:D:467:ASN:ND2	1:D:467:ASN:H	2.01	0.57
1:D:517:ASN:O	1:D:521:MET:HE2	2.05	0.57
1:D:509:LEU:O	1:D:512:GLN:HB2	2.04	0.57
1:A:479:ARG:HH12	1:A:508:GLU:HB2	1.68	0.56
1:D:549:ILE:CD1	1:D:555:VAL:HG22	2.21	0.56
1:A:477:PRO:HG2	1:A:509:LEU:HD11	1.87	0.56
2:E:9:DA:C5'	3:E:178:HOH:O	2.54	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:34:DT:H2''	2:F:35:DC:O5'	2.06	0.56
1:D:566:PRO:O	1:D:570:ILE:HG12	2.06	0.56
2:F:19:DG:H2''	2:F:20:DG:O5'	2.06	0.56
1:A:479:ARG:HH12	1:A:508:GLU:HB3	1.70	0.56
1:D:493:GLN:HG3	1:D:496:ARG:NH2	2.22	0.55
1:D:507:PRO:O	1:D:511:LYS:HE2	2.05	0.55
1:D:480:ILE:HD12	1:D:502:GLU:HB3	1.87	0.55
2:B:8:DG:H2''	2:B:9:DA:OP2	2.07	0.54
2:C:25:DC:H5''	3:C:133:HOH:O	2.08	0.54
1:D:446:LYS:O	1:D:446:LYS:HG2	2.07	0.54
2:B:8:DG:H2'	2:B:8:DG:OP2	2.08	0.54
1:A:462:ASN:HA	1:A:512:GLN:NE2	2.22	0.53
2:C:34:DT:H2''	2:C:35:DC:H5'	1.88	0.53
1:D:507:PRO:O	1:D:511:LYS:HG2	2.08	0.53
1:D:537:LYS:HB2	2:F:27:DA:H4'	1.89	0.53
2:B:1:DG:H2''	2:B:2:DG:OP2	2.08	0.53
1:D:455:GLY:HA2	1:D:559:ASP:OD2	2.09	0.53
1:A:538:ARG:HD2	1:A:542:GLU:OE2	2.09	0.53
2:F:26:DG:P	3:F:2:HOH:O	2.66	0.53
1:D:574:ARG:O	1:D:574:ARG:HG3	2.08	0.53
2:B:10:DT:H2''	2:B:11:DC:C5'	2.31	0.53
2:F:25:DC:H1'	2:F:26:DG:H5'	1.91	0.52
1:A:453:ALA:HB2	1:A:489:PHE:CE2	2.41	0.52
1:D:518:ASN:O	1:D:522:LYS:HG3	2.11	0.51
2:B:14:DA:H5''	3:B:153:HOH:O	2.11	0.51
1:D:546:GLU:HB3	1:D:560:ASN:HD22	1.76	0.51
2:C:27:DA:C8	2:C:28:DT:H72	2.46	0.50
2:B:14:DA:C5'	3:B:153:HOH:O	2.59	0.50
2:C:25:DC:C2'	2:C:26:DG:H5''	2.40	0.50
2:B:9:DA:C2'	2:B:10:DT:H5'	2.41	0.49
1:A:531:THR:C	1:A:533:ASP:H	2.15	0.49
1:A:517:ASN:O	1:A:521:MET:HE2	2.13	0.49
2:B:5:DT:H2''	2:B:6:DT:OP2	2.13	0.48
2:E:7:DC:H2''	2:E:8:DG:C8	2.49	0.48
1:D:485:ARG:O	1:D:487:LYS:HG2	2.14	0.48
2:F:32:DA:OP2	3:F:7:HOH:O	2.20	0.47
1:A:484:LEU:O	1:A:487:LYS:HB2	2.14	0.47
2:E:8:DG:H2''	2:E:9:DA:OP2	2.14	0.47
1:A:468:ALA:HA	1:A:473:GLY:HA2	1.97	0.47
1:D:523:VAL:HG23	1:D:525:LYS:HG3	1.97	0.47
2:F:26:DG:C5'	3:F:2:HOH:O	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:35:DC:H2'	2:F:36:DC:C6	2.50	0.46
1:A:542:GLU:HG2	2:B:12:DG:H5'	1.96	0.46
1:D:510:SER:O	1:D:512:GLN:N	2.48	0.46
1:A:557:ASP:O	1:A:559:ASP:N	2.49	0.46
1:A:529:THR:HG23	1:A:534:VAL:HG12	1.98	0.46
1:D:480:ILE:CD1	1:D:502:GLU:HB3	2.46	0.46
1:A:467:ASN:HD22	1:A:467:ASN:N	2.03	0.45
1:A:479:ARG:NH1	1:A:508:GLU:CB	2.72	0.45
1:A:549:ILE:C	1:A:551:GLN:H	2.19	0.45
1:A:546:GLU:HB3	1:A:560:ASN:ND2	2.32	0.45
2:C:26:DG:C2'	2:C:27:DA:C5'	2.94	0.44
1:A:476:VAL:HG13	1:A:499:PHE:HE1	1.81	0.44
1:A:518:ASN:O	1:A:522:LYS:HG3	2.18	0.44
1:D:468:ALA:HA	1:D:473:GLY:HA2	2.00	0.44
1:D:447:ARG:CG	1:D:448:ASN:ND2	2.81	0.43
1:A:496:ARG:O	1:A:499:PHE:HB3	2.18	0.43
1:D:530:ARG:HD2	3:D:1:HOH:O	2.18	0.43
1:A:498:LYS:HA	1:A:498:LYS:HD3	1.84	0.43
2:B:9:DA:H1'	2:B:10:DT:H5'	2.00	0.43
1:A:526:ALA:HB3	1:A:540:SER:HB3	2.00	0.43
2:B:7:DC:H3'	3:B:102:HOH:O	2.17	0.43
1:A:531:THR:C	1:A:533:ASP:N	2.73	0.42
2:C:26:DG:P	3:C:122:HOH:O	2.78	0.42
1:D:552:ASN:O	1:D:552:ASN:OD1	2.37	0.42
2:C:19:DG:H2''	2:C:20:DG:O5'	2.20	0.42
1:D:467:ASN:ND2	1:D:467:ASN:N	2.65	0.42
2:B:9:DA:H2''	2:B:10:DT:OP2	2.20	0.42
2:E:9:DA:H5'	3:E:178:HOH:O	2.20	0.42
1:A:573:HIS:NE2	2:B:10:DT:H4'	2.35	0.42
1:D:544:HIS:O	1:D:561:ILE:HA	2.20	0.42
1:D:561:ILE:HG22	1:D:562:SER:N	2.34	0.42
2:B:7:DC:H2''	2:B:8:DG:C8	2.55	0.41
1:A:538:ARG:CD	1:A:542:GLU:OE2	2.68	0.41
1:D:545:HIS:CD2	1:D:549:ILE:HD11	2.56	0.41
2:F:35:DC:H5''	3:F:159:HOH:O	2.20	0.41
1:A:489:PHE:HB2	1:A:556:TYR:CE1	2.55	0.41
1:A:500:TRP:HB3	1:A:521:MET:HG2	2.02	0.41
2:B:2:DG:H2''	2:B:3:DA:OP2	2.20	0.41
1:D:447:ARG:HG3	1:D:448:ASN:ND2	2.36	0.41
1:A:484:LEU:HD22	1:A:489:PHE:HZ	1.85	0.41
2:B:9:DA:H1'	2:B:10:DT:C5'	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:484:LEU:HD13	1:A:495:PHE:CE2	2.55	0.41
1:D:470:LYS:HB2	1:D:470:LYS:NZ	2.36	0.41
1:D:455:GLY:CA	1:D:559:ASP:OD2	2.69	0.40
1:D:561:ILE:CG2	1:D:562:SER:N	2.84	0.40
1:D:510:SER:C	1:D:512:GLN:N	2.75	0.40
1:D:557:ASP:OD1	1:D:560:ASN:HB2	2.21	0.40
2:C:31:DA:H2"	2:C:32:DA:OP2	2.20	0.40
1:D:462:ASN:HA	1:D:512:GLN:NE2	2.36	0.40
2:B:14:DA:C2	2:B:15:DT:C2	3.08	0.40
1:D:472:LEU:HD12	1:D:472:LEU:HA	1.85	0.40
1:A:527:PRO:HB2	1:A:541:PHE:HD2	1.86	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	130/132 (98%)	115 (88%)	9 (7%)	6 (5%)	2	9
1	D	129/132 (98%)	111 (86%)	16 (12%)	2 (2%)	9	32
All	All	259/264 (98%)	226 (87%)	25 (10%)	8 (3%)	4	16

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	464	TRP
1	A	461	ASN
1	A	548	PRO
1	A	558	MET
1	D	511	LYS
1	D	450	PRO
1	A	574	ARG

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Mol	Chain	Res	Type
1	A	575	GLY

### 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	118/118 (100%)	109 (92%)	9 (8%)	13	36
1	D	117/118 (99%)	111 (95%)	6 (5%)	24	56
All	All	235/236 (100%)	220 (94%)	15 (6%)	17	45

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	462	ASN
1	A	467	ASN
1	A	508	GLU
1	A	516	ASN
1	A	532	GLN
1	A	539	THR
1	A	548	PRO
1	A	562	SER
1	A	571	ASP
1	D	449	LYS
1	D	461	ASN
1	D	462	ASN
1	D	467	ASN
1	D	472	LEU
1	D	497	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	467	ASN
1	A	493	GLN

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Mol	Chain	Res	Type
1	A	518	ASN
1	A	532	GLN
1	A	551	GLN
1	A	560	ASN
1	D	448	ASN
1	D	461	ASN
1	D	466	ASN
1	D	467	ASN
1	D	551	GLN
1	D	552	ASN
1	D	560	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 monosaccharides 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	132/132 (100%)	-0.39	0 100 100	23, 49, 70, 80	0
1	D	131/132 (99%)	-0.45	0 100 100	25, 44, 69, 79	0
2	B	18/18 (100%)	-0.44	0 100 100	34, 57, 76, 85	0
2	C	18/18 (100%)	-0.57	0 100 100	44, 57, 75, 80	0
2	E	18/18 (100%)	-0.40	0 100 100	31, 54, 70, 74	0
2	F	18/18 (100%)	-0.58	0 100 100	31, 50, 68, 72	0
All	All	335/336 (99%)	-0.44	0 100 100	23, 49, 72, 85	0

There are no RSRZ outliers to report.

### 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 monosaccharides 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.