



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:56 AM GMT

PDB ID : 3QE9
Title : Crystal structure of human exonuclease 1 Exo1 (D173A) in complex with DNA (complex I)
Authors : Orans, J.; McSweeney, E.A.; Iyer, R.R.; Hast, M.A.; Hellinga, H.W.; Modrich, P.; Beese, L.S.
Deposited on : 2011-01-20
Resolution : 2.51 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

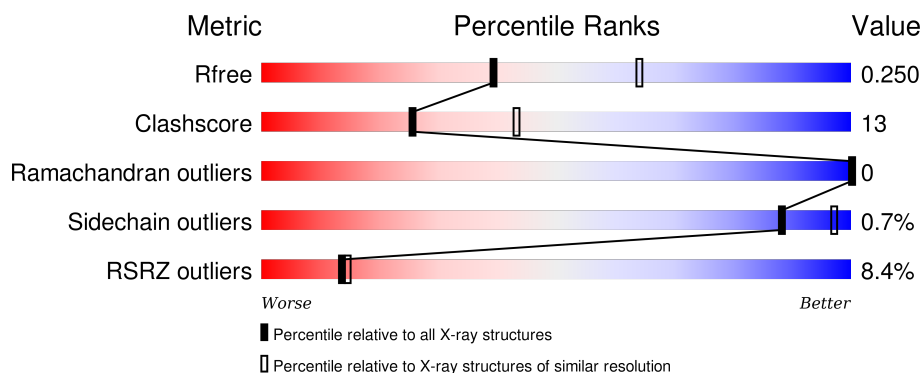
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 Å.

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



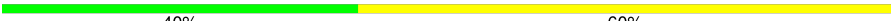
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	Y	352	<div> <div>13%</div> <div> <div></div> <div>72%</div> <div>26%</div> <div>•</div> </div> </div>
1	Z	352	<div> <div>5%</div> <div> <div></div> <div>79%</div> <div>19%</div> <div>•</div> </div> </div>
2	A	13	<div> <div></div> <div> <div>38%</div> <div>54%</div> <div>8%</div> </div> </div>
2	C	13	<div> <div>8%</div> <div> <div></div> <div>54%</div> <div>46%</div> </div> </div>
3	B	10	<div> <div></div> <div> <div>20%</div> <div>70%</div> <div>10%</div> </div> </div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	D	10	 A horizontal bar chart showing the quality of chain 3. The bar is divided into two segments: a green segment on the left and a yellow segment on the right. The green segment is labeled '40%' and the yellow segment is labeled '60%'. The total length of the bar represents 100%.

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6649 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 Exonuclease 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Y	345	Total	C	N	O	S	0	0	0
			2711	1726	471	496	18			
1	Z	344	Total	C	N	O	S	0	0	0
			2706	1723	470	495	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Y	173	ALA	ASP	ENGINEERED MUTATION	UNP Q9UQ84
Z	173	ALA	ASP	ENGINEERED MUTATION	UNP Q9UQ84

- Molecule 2 is a DNA chain called DNA (5'-D(*CP*GP*CP*TP*AP*GP*TP*CP*GP*AP*CP*AP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	13	Total	C	N	O	P	0	0	0
			262	126	48	76	12			
2	A	12	Total	C	N	O	P	0	0	0
			242	116	46	69	11			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	10	Total	C	N	O	P	0	0	0
			206	97	38	61	10			
3	B	10	Total	C	N	O	P	0	0	0
			206	97	38	61	10			

- Molecule 4 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	Z	1	Total K 1 1	0	0
4	Y	1	Total K 1 1	0	0

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	Z	1	Total Ca 1 1	0	0
5	Y	1	Total Ca 1 1	0	0

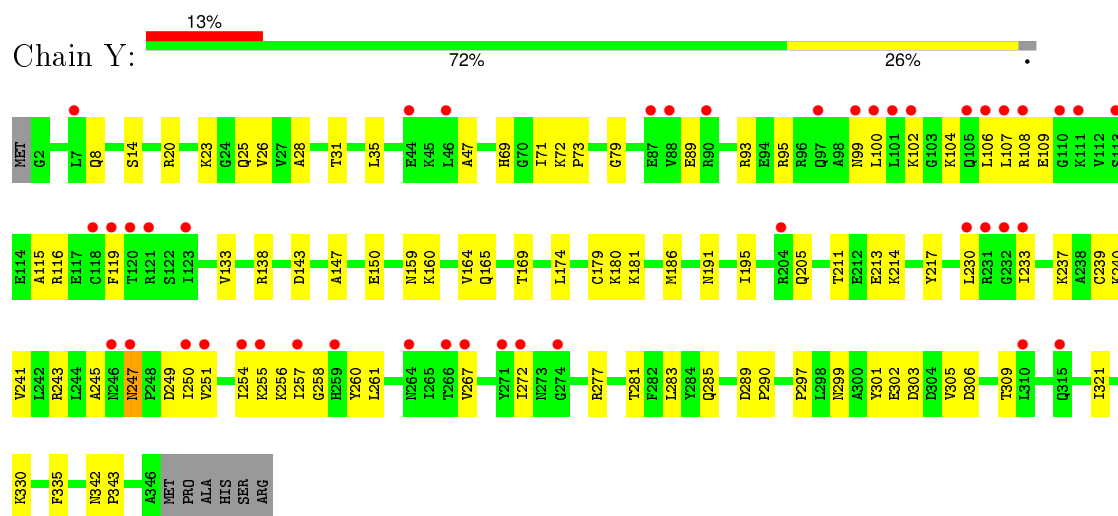
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	Y	102	Total O 102 102	0	0
6	C	4	Total O 4 4	0	0
6	D	7	Total O 7 7	0	0
6	Z	169	Total O 169 169	0	0
6	A	16	Total O 16 16	0	0
6	B	14	Total O 14 14	0	0

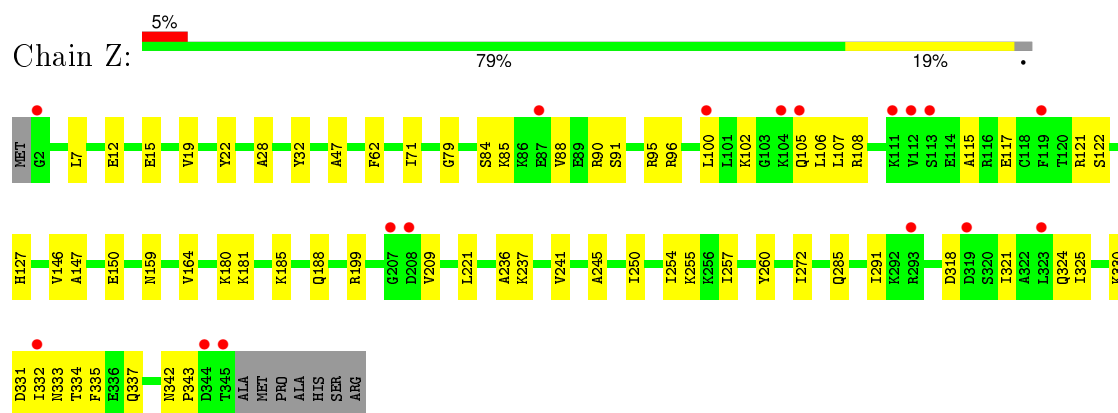
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Exonuclease 1



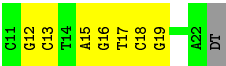
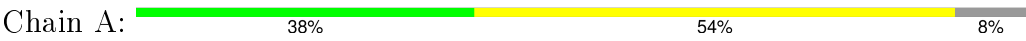
• Molecule 1: Exonuclease 1



• Molecule 2: DNA (5'-D(*CP*GP*CP*TP*AP*GP*TP*CP*GP*AP*CP*AP*T)-3')



• Molecule 2: DNA (5'-D(*CP*GP*CP*TP*AP*GP*TP*CP*GP*AP*CP*AP*T)-3')



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	92.50Å 95.47Å 99.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.51 46.25 – 2.51	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.51) 93.1 (46.25-2.51)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 2.51Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.215 , 0.238 0.215 , 0.250	Depositor DCC
R_{free} test set	1449 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	58.5	Xtriage
Anisotropy	0.331	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 57.4	EDS
Estimated twinning fraction	0.017 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 28797 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6649	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: K, CA

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	Y	0.34	0/2756	0.46	0/3713
1	Z	0.36	0/2751	0.47	0/3706
2	A	0.62	0/271	1.14	0/416
2	C	0.55	0/293	1.06	0/450
3	B	0.87	1/230 (0.4%)	1.13	0/351
3	D	0.64	0/230	1.20	0/351
All	All	0.42	1/6531 (0.0%)	0.63	0/8987

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1	DT	OP3-P	-10.24	1.48	1.61

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	Y	2711	0	2773	74	0
1	Z	2706	0	2768	53	0
2	A	242	0	136	11	0
2	C	262	0	148	4	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	206	0	113	12	0
3	D	206	0	113	8	0
4	Y	1	0	0	0	0
4	Z	1	0	0	0	0
5	Y	1	0	0	0	0
5	Z	1	0	0	0	0
6	A	16	0	0	0	0
6	B	14	0	0	1	0
6	C	4	0	0	0	0
6	D	7	0	0	0	0
6	Y	102	0	0	8	0
6	Z	169	0	0	9	0
All	All	6649	0	6051	155	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 13.

All (155) 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:Y:256:LYS:HB3	1:Y:260:TYR:CE2	1.95	1.02
1:Z:330:LYS:HD2	1:Z:335:PHE:HA	1.43	0.98
1:Y:256:LYS:HB3	1:Y:260:TYR:HE2	1.25	0.97
1:Y:107:LEU:CB	1:Y:115:ALA:HB2	1.96	0.94
3:B:6:DT:H1'	6:B:296:HOH:O	1.69	0.93
1:Y:107:LEU:HB2	1:Y:115:ALA:HB2	1.53	0.90
1:Y:20:ARG:HG3	1:Y:23:LYS:HE3	1.58	0.85
1:Y:8:GLN:NE2	3:D:4:DA:OP2	2.14	0.81
1:Y:159:ASN:HD21	1:Y:181:LYS:H	1.29	0.80
6:Y:454:HOH:O	3:D:2:DC:H5"	1.83	0.79
1:Z:318:ASP:HB3	1:Z:321:ILE:HD13	1.65	0.79
1:Y:102:LYS:O	1:Y:106:LEU:HB2	1.83	0.77
1:Z:85:LYS:HG2	1:Z:88:VAL:CG1	2.15	0.76
1:Z:188:GLN:HB2	6:Z:506:HOH:O	1.85	0.75
1:Y:107:LEU:HB3	1:Y:115:ALA:HB2	1.68	0.75
1:Y:138:ARG:NH1	6:Y:409:HOH:O	2.18	0.75
1:Y:256:LYS:CB	1:Y:260:TYR:HE2	2.01	0.72
1:Y:25:GLN:HB3	1:Y:165:GLN:HE21	1.56	0.71
1:Y:306:ASP:O	1:Y:309:THR:HG22	1.89	0.71
1:Z:330:LYS:CD	1:Z:335:PHE:HA	2.20	0.69
1:Z:330:LYS:HD2	1:Z:335:PHE:CA	2.21	0.69

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Y:35:LEU:HD21	1:Y:133:VAL:HG21	1.74	0.69
3:B:6:DT:H2"	3:B:7:DA:C8	2.27	0.69
3:B:1:DT:H5'	3:B:1:DT:H6	1.57	0.69
1:Y:20:ARG:O	1:Y:23:LYS:HG3	1.95	0.67
1:Y:230:LEU:HB2	1:Y:233:ILE:HD12	1.76	0.67
1:Y:245:ALA:HA	1:Y:260:TYR:CE1	2.29	0.66
1:Z:90:ARG:O	6:Z:512:HOH:O	2.13	0.66
1:Z:221:LEU:HD21	1:Z:254:ILE:HD12	1.77	0.66
1:Z:209:VAL:O	6:Z:481:HOH:O	2.13	0.65
1:Y:104:LYS:O	1:Y:108:ARG:HB2	1.95	0.65
1:Y:277:ARG:HG2	1:Y:277:ARG:HH11	1.62	0.64
1:Z:331:ASP:OD1	1:Z:333:ASN:N	2.28	0.63
3:B:7:DA:H4'	3:B:8:DG:OP1	1.98	0.63
1:Y:159:ASN:ND2	1:Y:180:LYS:H	1.95	0.63
1:Y:89:GLU:O	1:Y:93:ARG:HG2	1.98	0.63
1:Y:160:LYS:HG2	6:Y:393:HOH:O	1.98	0.63
1:Z:32:TYR:HB2	6:Z:370:HOH:O	1.98	0.63
3:B:7:DA:H2'	3:B:8:DG:C8	2.33	0.62
1:Y:277:ARG:HH11	1:Y:277:ARG:CG	2.13	0.62
1:Z:85:LYS:HG2	1:Z:88:VAL:HG11	1.82	0.62
1:Y:26:VAL:H	1:Y:165:GLN:NE2	2.00	0.60
1:Z:91:SER:O	1:Z:95:ARG:HG3	2.01	0.60
2:A:15:DA:H61	3:B:6:DT:H3	1.49	0.60
3:D:7:DA:H2"	3:D:8:DG:H5'	1.83	0.60
1:Y:23:LYS:HG2	1:Y:69:HIS:O	2.02	0.59
1:Y:150:GLU:HB2	1:Y:285:GLN:OE1	2.02	0.59
3:D:4:DA:H1'	3:D:5:DC:H5"	1.85	0.58
1:Y:330:LYS:HE2	6:Y:439:HOH:O	2.03	0.58
1:Z:150:GLU:HB2	1:Z:285:GLN:OE1	2.04	0.58
1:Y:302:GLU:HG2	1:Y:303:ASP:H	1.66	0.58
1:Y:14:SER:HB3	1:Y:195:ILE:HG12	1.85	0.58
1:Z:245:ALA:HB2	1:Z:260:TYR:CZ	2.38	0.57
1:Y:95:ARG:O	1:Y:99:ASN:HB2	2.04	0.57
1:Z:96:ARG:HG2	1:Z:122:SER:HA	1.87	0.56
1:Z:105:GLN:HG3	1:Z:108:ARG:HD2	1.86	0.56
1:Y:249:ASP:OD1	1:Y:251:VAL:HG12	2.05	0.56
1:Y:305:VAL:HG12	1:Y:306:ASP:N	2.20	0.56
1:Z:331:ASP:OD1	1:Z:332:ILE:N	2.40	0.55
1:Z:22:TYR:OH	1:Z:181:LYS:HD3	2.07	0.55
1:Z:291:ILE:HD11	6:Z:505:HOH:O	2.06	0.54
1:Z:334:THR:O	1:Z:335:PHE:HB2	2.07	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Y:180:LYS:HG3	6:Y:393:HOH:O	2.06	0.54
1:Y:213:GLU:HG2	1:Y:217:TYR:CE2	2.43	0.54
1:Y:256:LYS:O	1:Y:260:TYR:HD2	1.90	0.54
2:A:18:DC:H2"	2:A:19:DG:C8	2.43	0.53
1:Z:159:ASN:HD21	1:Z:181:LYS:H	1.56	0.53
1:Y:272:ILE:HG21	6:Y:452:HOH:O	2.07	0.52
1:Y:240:LYS:HG3	1:Y:243:ARG:NH2	2.24	0.52
1:Y:109:GLU:OE1	1:Y:109:GLU:N	2.42	0.52
3:B:2:DC:H2'	3:B:3:DG:C8	2.45	0.52
2:A:15:DA:N1	3:B:6:DT:O2	2.42	0.51
1:Z:159:ASN:ND2	1:Z:180:LYS:H	2.09	0.51
1:Z:255:LYS:HE2	1:Z:272:ILE:HD13	1.91	0.51
1:Z:102:LYS:HA	1:Z:106:LEU:HD12	1.93	0.51
1:Y:247:ASN:N	1:Y:247:ASN:OD1	2.44	0.50
1:Z:127:HIS:CE1	1:Z:332:ILE:HG13	2.46	0.50
1:Z:237:LYS:HG2	2:A:13:DC:OP1	2.12	0.50
1:Y:251:VAL:HG23	1:Y:272:ILE:HG23	1.92	0.50
1:Y:256:LYS:O	1:Y:260:TYR:CD2	2.64	0.50
1:Y:28:ALA:HB2	1:Y:164:VAL:HG11	1.93	0.49
1:Y:169:THR:HG21	1:Y:174:LEU:HD12	1.94	0.49
1:Y:237:LYS:HG2	2:C:14:DT:OP1	2.12	0.49
1:Y:240:LYS:HD3	1:Y:261:LEU:HD23	1.94	0.49
2:C:13:DC:H2"	2:C:14:DT:H5'	1.94	0.49
1:Z:85:LYS:O	1:Z:88:VAL:HG12	2.12	0.49
1:Z:236:ALA:HB3	2:A:12:DG:H3'	1.95	0.49
1:Y:179:CYS:HA	6:Y:393:HOH:O	2.11	0.49
2:A:16:DG:H2'	2:A:17:DT:H72	1.93	0.49
2:A:17:DT:H2'	2:A:17:DT:O5'	2.13	0.49
1:Y:211:THR:HG22	1:Y:213:GLU:H	1.77	0.49
1:Z:102:LYS:O	1:Z:106:LEU:HB2	2.12	0.48
1:Y:79:GLY:HA3	1:Y:147:ALA:O	2.13	0.48
1:Y:72:LYS:HB2	6:Z:509:HOH:O	2.14	0.48
1:Z:19:VAL:HB	1:Z:71:ILE:HD11	1.95	0.48
1:Z:19:VAL:HB	1:Z:71:ILE:CD1	2.43	0.48
3:D:3:DG:H2"	3:D:4:DA:C8	2.49	0.48
1:Y:250:ILE:O	1:Y:254:ILE:HG12	2.13	0.48
1:Z:250:ILE:O	1:Z:254:ILE:HG12	2.14	0.47
1:Y:186:MET:HA	1:Y:191:ASN:O	2.14	0.47
1:Y:239:CYS:O	1:Y:243:ARG:HB2	2.14	0.47
1:Z:47:ALA:HB1	1:Z:100:LEU:HD13	1.96	0.47
3:B:1:DT:H5'	3:B:1:DT:C6	2.44	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Z:7:LEU:HD21	1:Z:185:LYS:HE3	1.96	0.47
1:Z:85:LYS:HG2	1:Z:88:VAL:HG12	1.94	0.46
1:Z:321:ILE:HD12	1:Z:321:ILE:N	2.31	0.46
1:Y:277:ARG:CG	1:Y:277:ARG:NH1	2.75	0.46
1:Y:257:ILE:HG13	1:Y:258:GLY:N	2.30	0.46
1:Y:305:VAL:HG12	1:Y:306:ASP:H	1.81	0.46
1:Z:331:ASP:OD1	1:Z:331:ASP:C	2.54	0.46
1:Y:116:ARG:HA	1:Y:119:PHE:HD2	1.82	0.45
2:A:16:DG:H2'	2:A:17:DT:C7	2.47	0.45
1:Z:337:GLN:HB3	6:Z:477:HOH:O	2.16	0.45
1:Z:28:ALA:HB2	1:Z:164:VAL:HG11	1.99	0.45
1:Y:297:PRO:HG3	1:Y:301:TYR:CE1	2.52	0.45
1:Z:79:GLY:HA3	1:Z:147:ALA:O	2.17	0.45
1:Z:62:PHE:HD2	6:Z:377:HOH:O	1.99	0.45
1:Z:85:LYS:CG	1:Z:88:VAL:HG11	2.46	0.45
1:Z:146:VAL:HB	1:Z:325:ILE:O	2.17	0.45
3:B:9:DC:H2''	3:B:10:DG:C8	2.51	0.45
1:Z:324:GLN:HG3	1:Z:337:GLN:HE22	1.81	0.45
1:Y:281:THR:O	1:Y:285:GLN:HB2	2.17	0.44
1:Y:302:GLU:CG	1:Y:303:ASP:H	2.30	0.44
1:Z:105:GLN:HA	1:Z:108:ARG:HD2	1.98	0.44
1:Y:205:GLN:O	1:Y:239:CYS:HB3	2.18	0.44
1:Z:330:LYS:HE2	1:Z:335:PHE:O	2.18	0.44
1:Y:71:ILE:O	1:Y:73:PRO:HD3	2.17	0.44
3:B:1:DT:H2'	3:B:2:DC:C6	2.53	0.44
1:Y:335:PHE:CD2	1:Y:335:PHE:O	2.71	0.43
1:Z:12:GLU:H	1:Z:12:GLU:CD	2.21	0.43
2:A:15:DA:N6	3:B:6:DT:H3	2.16	0.43
1:Y:321:ILE:HG23	1:Y:335:PHE:CE1	2.53	0.43
1:Z:15:GLU:HB2	1:Z:199:ARG:HH12	1.84	0.43
1:Z:117:GLU:O	1:Z:121:ARG:HG3	2.18	0.43
1:Y:243:ARG:HG2	6:Y:400:HOH:O	2.18	0.43
1:Y:283:LEU:O	1:Y:299:ASN:HB2	2.19	0.43
1:Y:72:LYS:HE2	6:Z:515:HOH:O	2.18	0.42
1:Z:107:LEU:HB3	1:Z:115:ALA:HB2	2.01	0.42
1:Y:26:VAL:HG12	1:Y:164:VAL:HG12	2.01	0.42
2:A:16:DG:H2''	2:A:17:DT:C6	2.54	0.42
3:D:4:DA:H2''	3:D:5:DC:H5'	2.01	0.42
3:D:2:DC:H2''	3:D:3:DG:O5'	2.19	0.42
1:Z:342:ASN:HA	1:Z:343:PRO:HD2	1.90	0.42
2:A:17:DT:C2'	2:A:17:DT:O5'	2.68	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Y:211:THR:HB	1:Y:214:LYS:HG3	2.02	0.41
1:Y:241:VAL:HG11	1:Y:257:ILE:HG22	2.01	0.41
1:Y:47:ALA:HB1	1:Y:100:LEU:CD1	2.50	0.41
1:Y:255:LYS:HG2	1:Y:267:VAL:HG21	2.01	0.41
1:Z:241:VAL:HG21	1:Z:257:ILE:HG13	2.02	0.41
1:Y:289:ASP:HA	1:Y:290:PRO:HD2	1.97	0.41
1:Y:159:ASN:ND2	1:Y:181:LYS:H	2.08	0.41
1:Y:342:ASN:HA	1:Y:343:PRO:HD3	1.81	0.40
3:D:2:DC:H2'	3:D:3:DG:C8	2.55	0.40
2:C:16:DG:H2'	2:C:17:DT:H72	2.04	0.40
2:C:22:DA:H2''	2:C:23:DT:O4'	2.22	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	Y	343/352 (97%)	333 (97%)	10 (3%)	0	100	100
1	Z	342/352 (97%)	338 (99%)	4 (1%)	0	100	100
All	All	685/704 (97%)	671 (98%)	14 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	Y	291/297 (98%)	288 (99%)	3 (1%)	82	95
1	Z	291/297 (98%)	290 (100%)	1 (0%)	94	99
All	All	582/594 (98%)	578 (99%)	4 (1%)	88	97

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	Y	31	THR
1	Y	143	ASP
1	Y	247	ASN
1	Z	84	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	Y	69	HIS
1	Y	159	ASN
1	Y	165	GLN
1	Y	188	GLN
1	Y	197	GLN
1	Y	205	GLN
1	Z	8	GLN
1	Z	99	ASN
1	Z	105	GLN
1	Z	159	ASN
1	Z	337	GLN

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	Y	345/352 (98%)	0.80	44 (12%) 5 5	35, 57, 98, 109	0
1	Z	344/352 (97%)	0.62	17 (4%) 33 38	34, 45, 88, 111	0
2	A	12/13 (92%)	-0.24	0 100 100	50, 72, 83, 89	0
2	C	13/13 (100%)	0.88	1 (7%) 16 18	67, 81, 97, 97	0
3	B	10/10 (100%)	-0.13	0 100 100	54, 76, 84, 85	0
3	D	10/10 (100%)	0.99	0 100 100	70, 86, 92, 95	0
All	All	734/750 (97%)	0.69	62 (8%) 14 14	34, 53, 95, 111	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	Z	112	VAL	10.7
1	Z	111	LYS	8.9
1	Y	110	GLY	6.4
1	Z	345	THR	5.4
1	Y	111	LYS	5.2
1	Y	254	ILE	4.9
1	Z	87	GLU	4.7
2	C	23	DT	4.1
1	Y	101	LEU	4.1
1	Y	247	ASN	4.0
1	Y	106	LEU	4.0
1	Y	251	VAL	4.0
1	Y	259	HIS	4.0
1	Y	264	ASN	3.7
1	Y	120	THR	3.7
1	Z	113	SER	3.6
1	Y	310	LEU	3.4
1	Y	97	GLN	3.3
1	Y	100	LEU	3.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	Z	2	GLY	3.3
1	Y	233	ILE	3.1
1	Y	272	ILE	3.0
1	Y	108	ARG	3.0
1	Y	123	ILE	3.0
1	Y	231	ARG	3.0
1	Y	88	VAL	2.9
1	Y	267	VAL	2.8
1	Y	266	THR	2.8
1	Y	257	ILE	2.8
1	Z	293	ARG	2.8
1	Y	87	GLU	2.8
1	Y	232	GLY	2.8
1	Z	344	ASP	2.6
1	Y	90	ARG	2.6
1	Y	246	ASN	2.6
1	Y	271	TYR	2.6
1	Y	118	CYS	2.6
1	Y	204	ARG	2.5
1	Y	99	ASN	2.5
1	Y	119	PHE	2.5
1	Y	46	LEU	2.5
1	Y	121	ARG	2.5
1	Z	100	LEU	2.5
1	Y	274	GLY	2.5
1	Y	102	LYS	2.5
1	Z	207	GLY	2.4
1	Y	255	LYS	2.4
1	Y	230	LEU	2.4
1	Y	113	SER	2.3
1	Z	119	PHE	2.3
1	Y	250	ILE	2.3
1	Y	315	GLN	2.2
1	Z	104	LYS	2.2
1	Z	105	GLN	2.2
1	Y	105	GLN	2.2
1	Z	319	ASP	2.2
1	Y	44	GLU	2.1
1	Y	7	LEU	2.1
1	Y	107	LEU	2.1
1	Z	332	ILE	2.1
1	Z	208	ASP	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	Z	323	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	K	Z	353	1/1	0.95	0.15	-1.23	61,61,61,61	0
4	K	Y	353	1/1	0.66	0.08	-1.35	98,98,98,98	0
5	CA	Z	354	1/1	0.76	0.17	-	73,73,73,73	0
5	CA	Y	354	1/1	0.87	0.24	-	78,78,78,78	0

6.5 Other polymers [i](#)

There are no such residues in this entry.