



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 07:23 PM UTC

PDB ID : 4G3I / pdb_00004g3i
Title : Crystal structure of Dpo4 in complex with DNA duplex
Authors : Yan, Y.Y.; Gan, J.H.; Huang, Z.
Deposited on : 2012-07-14
Resolution : 2.50 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

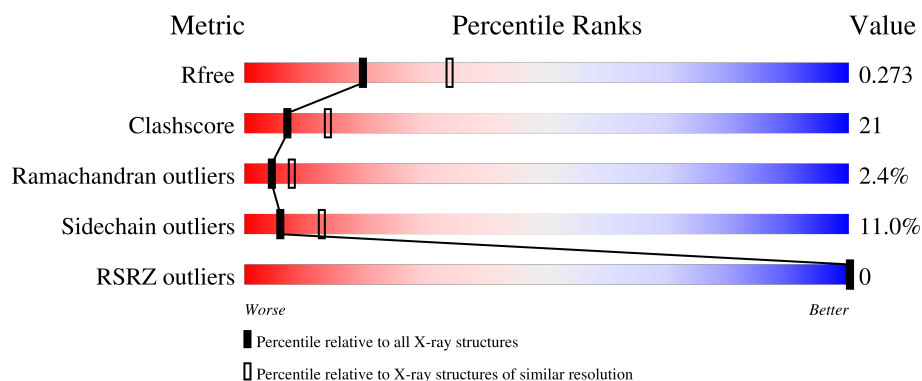
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (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 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	16	
1	E	16	
2	D	13	
2	F	13	
3	A	342	

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Mol	Chain	Length	Quality of chain
3	B	342	 A horizontal bar chart showing the quality of chain B. The bar is divided into three segments: green (61%), yellow (31%), and red (7%). The segments are labeled with their respective percentages: 61%, 31%, and 7%. A small black dot is visible at the end of the red segment.

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6238 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 DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	10	Total	C	N	O	P	0	0	0
			198	96	33	60	9			
1	E	10	Total	C	N	O	P	0	0	0
			198	96	33	60	9			

- Molecule 2 is a DNA chain called DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	8	Total	C	N	O	P	0	0	0
			163	78	33	45	7			
2	F	8	Total	C	N	O	P	0	0	0
			163	78	33	45	7			

- Molecule 3 is a protein called DNA polymerase IV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	342	Total	C	N	O	S	0	0	0
			2750	1763	473	507	7			
3	B	341	Total	C	N	O	S	0	0	0
			2744	1760	472	505	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q97W02
B	0	SER	-	expression tag	UNP Q97W02

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Ca	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total 1	Ca 1	0	0

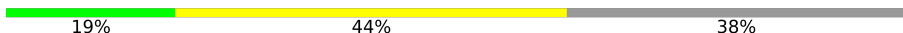
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	1	Total 1	O 1	0	0
5	A	16	Total 16	O 16	0	0
5	B	3	Total 3	O 3	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.

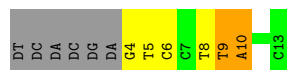
- Molecule 1: DNA

Chain C: 

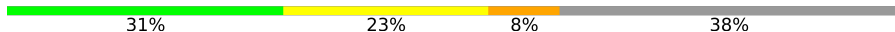


- Molecule 1: DNA

Chain E: 

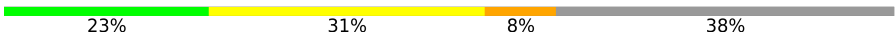


- Molecule 2: DNA

Chain D: 



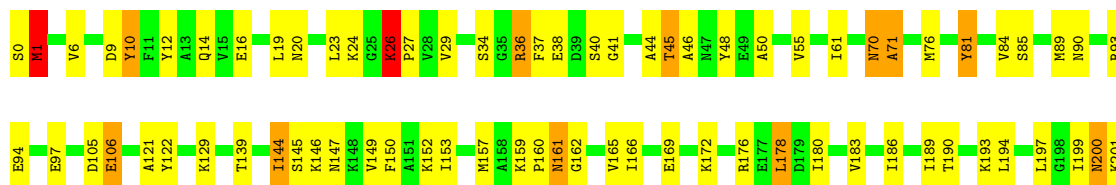
- Molecule 2: DNA

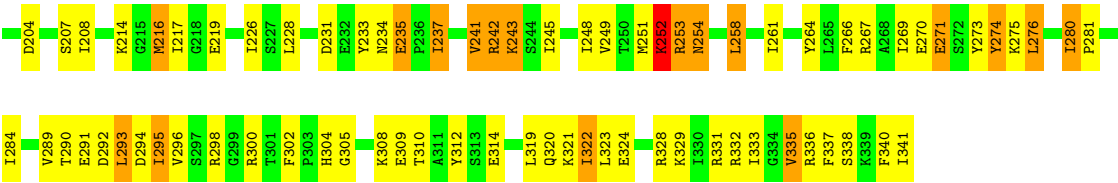
Chain F: 



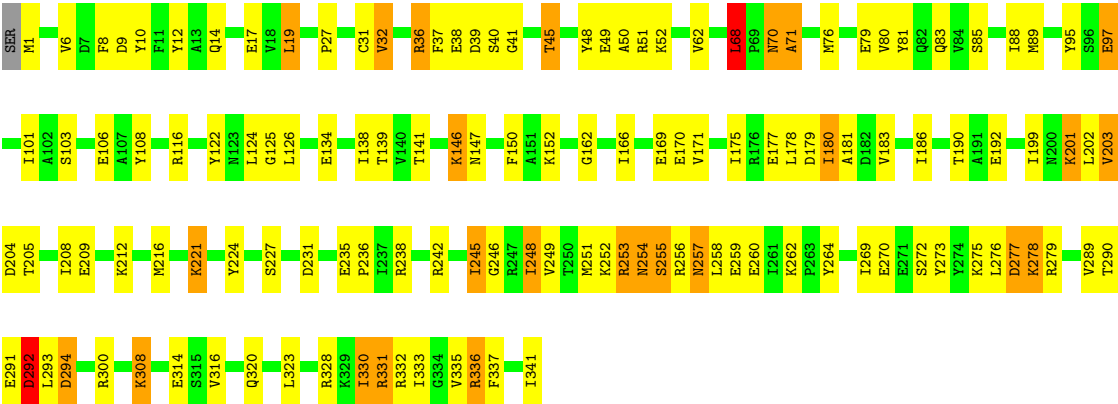
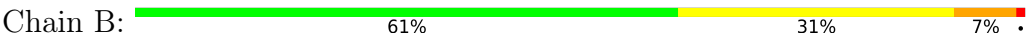
- Molecule 3: DNA polymerase IV

Chain A: 





• Molecule 3: DNA polymerase IV



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	52.47Å 52.63Å 99.44Å 75.40° 80.35° 70.97°	Depositor
Resolution (Å)	95.81 – 2.50 95.81 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.1 (95.81-2.50) 97.1 (95.81-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.78 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.241 , 0.294 0.249 , 0.273	Depositor DCC
R_{free} test set	1650 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	53.7	Xtriage
Anisotropy	0.168	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 39.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6238	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.57% 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

Bond lengths and bond angles in the following residue types are not validated in this section: 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	C	0.61	0/220	1.39	3/337 (0.9%)
1	E	0.60	0/220	1.54	5/337 (1.5%)
2	D	0.59	0/183	1.50	2/281 (0.7%)
2	F	0.57	0/183	1.42	3/281 (1.1%)
3	A	1.02	1/2789 (0.0%)	1.21	13/3744 (0.3%)
3	B	1.05	1/2783 (0.0%)	1.18	12/3736 (0.3%)
All	All	0.99	2/6378 (0.0%)	1.24	38/8716 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	280	ILE	CA-CB	6.16	1.61	1.54
3	B	330	ILE	CA-CB	5.17	1.60	1.54

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	84	VAL	N-CA-C	-7.65	103.34	110.53
3	B	273	TYR	N-CA-C	6.95	118.50	111.07
3	A	105	ASP	N-CA-C	6.57	120.52	111.71
3	B	80	VAL	N-CA-C	-6.56	105.35	111.45
3	B	70	ASN	N-CA-C	6.38	118.32	111.36
2	D	5	DG	N9-C1'-C2'	6.37	123.06	113.50
2	F	10	DC	O3'-P-O5'	-6.33	94.51	104.00
3	A	296	VAL	N-CA-C	6.23	116.85	107.75
1	E	6	DC	P-O3'-C3'	6.08	129.33	120.20
3	A	90	ASN	N-CA-C	6.03	117.52	111.07
1	E	8	DT	C3'-C2'-C1'	-6.03	92.56	101.60
3	B	235	GLU	CA-C-N	-6.00	114.43	120.31
3	B	235	GLU	C-N-CA	-6.00	114.43	120.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	68	LEU	CA-C-N	-5.97	113.47	119.56
3	B	68	LEU	C-N-CA	-5.97	113.47	119.56
3	B	17	GLU	N-CA-C	5.93	118.53	111.71
3	A	61	ILE	CB-CA-C	-5.93	104.04	112.22
3	A	295	ILE	N-CA-C	5.82	121.45	109.34
3	B	95	TYR	N-CA-C	5.79	118.37	111.71
1	C	4	DG	C3'-C2'-C1'	-5.79	92.92	101.60
3	B	41	GLY	N-CA-C	5.77	117.76	110.38
1	C	4	DG	N9-C1'-C2'	5.72	122.07	113.50
3	A	34	SER	N-CA-C	5.61	118.12	111.33
3	A	41	GLY	N-CA-C	5.59	117.53	110.38
2	D	11	DA	P-O3'-C3'	5.35	128.22	120.20
2	F	7	DC	O5'-C5'-C4'	-5.27	102.89	110.80
3	A	338	SER	N-CA-C	5.26	116.00	108.74
1	E	9	DT	O5'-C5'-C4'	5.22	118.63	110.80
2	F	10	DC	C1'-O4'-C4'	-5.22	101.87	109.70
3	A	235	GLU	CA-C-N	5.22	125.50	119.92
3	A	235	GLU	C-N-CA	5.22	125.50	119.92
1	E	10	DA	P-O5'-C5'	-5.20	112.19	120.00
3	B	31	CYS	N-CA-C	5.20	118.20	110.14
3	A	145	SER	N-CA-C	5.20	115.39	108.38
3	B	235	GLU	N-CA-C	5.19	116.42	109.83
3	A	70	ASN	N-CA-C	5.08	118.51	111.71
1	C	8	DT	O5'-C5'-C4'	5.05	118.38	110.80
1	E	8	DT	C1'-O4'-C4'	-5.01	102.18	109.70

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	198	0	115	6	0
1	E	198	0	115	2	0
2	D	163	0	91	6	0
2	F	163	0	91	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	2750	0	2894	130	0
3	B	2744	0	2889	121	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	16	0	0	0	0
5	B	3	0	0	1	0
5	C	1	0	0	0	0
All	All	6238	0	6195	259	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 21.

All (259) 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:B:254:ASN:HB3	3:B:330:ILE:O	1.43	1.18
3:B:256:ARG:O	3:B:323:LEU:HD21	1.46	1.15
3:B:12:TYR:HB2	3:B:45:THR:HG21	1.18	1.13
3:B:116:ARG:HB3	3:B:116:ARG:NH1	1.77	0.99
3:A:294:ASP:CB	3:A:328:ARG:HH12	1.75	0.98
3:A:242:ARG:HG3	3:A:242:ARG:HH11	1.30	0.95
3:B:12:TYR:HB2	3:B:45:THR:CG2	1.98	0.91
3:B:116:ARG:HB3	3:B:116:ARG:HH11	1.36	0.91
3:A:199:ILE:HG23	3:A:204:ASP:HB2	1.55	0.89
3:A:14:GLN:NE2	3:A:139:THR:H	1.71	0.89
3:A:294:ASP:HB3	3:A:328:ARG:NH1	1.87	0.88
3:A:14:GLN:HE22	3:A:139:THR:H	0.93	0.88
3:A:149:VAL:O	3:A:153:ILE:HG13	1.74	0.88
3:B:254:ASN:CB	3:B:330:ILE:O	2.22	0.87
3:A:294:ASP:HB3	3:A:328:ARG:HH12	1.37	0.85
3:B:48:TYR:HA	3:B:51:ARG:HG3	1.59	0.85
3:B:248:ILE:HD11	3:B:332:ARG:HB3	1.58	0.84
3:B:12:TYR:CB	3:B:45:THR:HG21	2.06	0.84
3:B:38:GLU:O	3:B:39:ASP:HB2	1.79	0.83
3:B:251:MET:CE	3:B:330:ILE:HG22	2.09	0.83
3:A:37:PHE:CE1	3:A:40:SER:HB3	2.13	0.81
3:B:278:LYS:HA	3:B:278:LYS:HE2	1.63	0.79
3:A:0:SER:O	3:A:1:MET:HB3	1.83	0.79
3:B:166:ILE:HG23	3:B:170:GLU:HG3	1.63	0.79
3:A:14:GLN:HE22	3:A:139:THR:N	1.77	0.78
3:A:200:ASN:H	3:A:200:ASN:HD22	1.32	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:194:LEU:HD22	3:A:199:ILE:HD12	1.67	0.76
3:B:254:ASN:ND2	3:B:331:ARG:HD3	2.01	0.76
3:A:12:TYR:HE2	3:A:76:MET:HE2	1.50	0.76
1:C:9:DT:H2''	1:C:10:DA:OP2	1.85	0.75
2:F:10:DC:H2'	2:F:11:DA:C8	2.22	0.74
3:B:254:ASN:HB3	3:B:330:ILE:C	2.13	0.74
3:B:14:GLN:NE2	3:B:139:THR:H	1.87	0.73
2:D:8:DT:O3'	3:A:242:ARG:NH1	2.21	0.72
3:A:46:ALA:HB1	3:A:50:ALA:HB3	1.70	0.72
2:D:9:DA:P	3:A:242:ARG:HH12	2.13	0.71
3:B:277:ASP:O	3:B:278:LYS:HB2	1.89	0.71
3:B:146:LYS:HD3	3:B:171:VAL:HG21	1.72	0.71
3:A:48:TYR:OH	3:A:159:LYS:HD3	1.91	0.71
3:B:254:ASN:OD1	3:B:254:ASN:N	2.22	0.71
3:A:304:HIS:HD2	3:A:305:GLY:O	1.75	0.70
3:A:242:ARG:HG3	3:A:242:ARG:NH1	2.04	0.69
3:B:256:ARG:O	3:B:323:LEU:CD2	2.36	0.68
3:A:12:TYR:CE2	3:A:76:MET:HE2	2.27	0.68
2:D:8:DT:OP2	3:A:336:ARG:NH2	2.25	0.67
3:A:37:PHE:CZ	3:A:40:SER:HB3	2.30	0.66
3:B:166:ILE:HG23	3:B:170:GLU:CG	2.25	0.66
3:B:14:GLN:HE22	3:B:139:THR:H	1.41	0.65
3:A:37:PHE:CD1	3:A:40:SER:HB3	2.31	0.65
3:B:147:ASN:HD21	3:B:150:PHE:HD2	1.44	0.64
3:A:271:GLU:O	3:A:274:TYR:HB3	1.98	0.64
3:B:116:ARG:NH1	3:B:116:ARG:CB	2.60	0.64
3:B:14:GLN:HG2	3:B:138:ILE:HD13	1.80	0.63
3:B:175:ILE:HG22	3:B:203:VAL:HG23	1.80	0.63
3:B:36:ARG:NH1	3:B:331:ARG:HD2	2.13	0.63
3:A:44:ALA:CB	3:A:76:MET:HE3	2.29	0.63
3:A:294:ASP:HB3	3:A:328:ARG:CZ	2.29	0.62
2:F:8:DT:OP2	3:B:336:ARG:NH2	2.33	0.62
3:A:190:THR:HG23	3:A:217:ILE:HG21	1.80	0.62
3:A:251:MET:HG2	3:A:264:TYR:CD2	2.35	0.62
3:A:275:LYS:O	3:A:276:LEU:C	2.43	0.61
3:B:253:ARG:O	3:B:253:ARG:HG2	2.00	0.61
3:B:8:PHE:CZ	3:B:88:ILE:HG21	2.36	0.61
3:B:122:TYR:CE2	3:B:126:LEU:HD11	2.35	0.61
2:D:9:DA:OP1	3:A:242:ARG:NH1	2.33	0.61
3:A:6:VAL:O	3:A:106:GLU:HA	2.00	0.61
3:A:294:ASP:HB2	3:A:328:ARG:HH12	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:200:ASN:H	3:A:200:ASN:ND2	1.96	0.61
3:A:292:ASP:O	3:A:293:LEU:HB2	2.00	0.61
3:A:200:ASN:HD22	3:A:200:ASN:N	1.99	0.60
3:A:253:ARG:O	3:A:253:ARG:HG3	1.99	0.60
3:A:309:GLU:O	3:A:312:TYR:HB2	2.00	0.60
3:A:149:VAL:HG11	3:A:228:LEU:HD11	1.84	0.60
3:A:200:ASN:ND2	3:A:200:ASN:N	2.47	0.60
3:B:333:ILE:O	3:B:333:ILE:HG23	2.00	0.60
3:A:121:ALA:HB1	3:A:144:ILE:HD13	1.82	0.60
3:B:101:ILE:HD12	3:B:101:ILE:H	1.67	0.60
3:B:236:PRO:HG2	3:B:238:ARG:NH1	2.16	0.60
3:A:26:LYS:HB3	3:A:27:PRO:HD2	1.84	0.59
3:A:251:MET:HG2	3:A:264:TYR:CG	2.38	0.59
1:E:9:DT:H2''	1:E:10:DA:OP2	2.01	0.59
3:B:27:PRO:HB2	3:B:71:ALA:HB2	1.85	0.58
3:A:9:ASP:O	3:A:10:TYR:C	2.47	0.58
3:A:258:LEU:HG	3:A:323:LEU:HD22	1.86	0.58
3:B:258:LEU:HG	3:B:262:LYS:HD2	1.85	0.57
3:A:237:ILE:HG22	3:A:237:ILE:O	2.02	0.57
3:A:273:TYR:HA	3:A:276:LEU:HD12	1.85	0.57
3:B:27:PRO:HB2	3:B:71:ALA:CB	2.35	0.56
3:A:228:LEU:HD23	3:A:233:TYR:HB2	1.87	0.56
3:B:1:MET:HE2	3:B:146:LYS:O	2.05	0.56
3:A:269:ILE:HG13	3:A:335:VAL:HG11	1.87	0.56
3:B:255:SER:CB	3:B:260:GLU:OE1	2.54	0.56
3:B:9:ASP:O	3:B:10:TYR:C	2.48	0.55
3:A:258:LEU:HD11	3:A:320:GLN:HG2	1.88	0.55
3:B:180:ILE:CG2	3:B:181:ALA:N	2.68	0.55
3:A:1:MET:HG2	3:A:146:LYS:O	2.07	0.55
3:B:116:ARG:CB	3:B:116:ARG:CZ	2.85	0.55
2:F:10:DC:C2'	2:F:11:DA:C8	2.89	0.55
3:A:309:GLU:OE1	3:A:309:GLU:N	2.35	0.55
3:B:36:ARG:HH12	3:B:331:ARG:NE	2.04	0.55
3:A:233:TYR:CZ	3:A:235:GLU:HB2	2.43	0.54
3:A:294:ASP:OD2	3:A:328:ARG:NH2	2.41	0.54
3:B:289:VAL:HA	3:B:294:ASP:O	2.07	0.54
3:B:146:LYS:HE2	3:B:231:ASP:OD1	2.07	0.54
3:A:300:ARG:HD3	3:A:314:GLU:OE1	2.08	0.54
3:B:251:MET:HE2	3:B:330:ILE:HG22	1.89	0.54
1:C:10:DA:H3'	3:A:189:ILE:HB	1.90	0.53
3:B:291:GLU:C	3:B:293:LEU:H	2.16	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:36:ARG:HH12	3:B:331:ARG:HD2	1.72	0.53
3:B:180:ILE:O	3:B:183:VAL:HG23	2.09	0.53
2:D:9:DA:P	3:A:242:ARG:NH1	2.81	0.53
3:A:291:GLU:HB3	3:A:329:LYS:HB2	1.91	0.52
3:A:309:GLU:CD	3:A:309:GLU:H	2.16	0.52
3:B:257:ASN:ND2	3:B:259:GLU:H	2.07	0.52
3:A:322:ILE:HG22	3:A:323:LEU:N	2.24	0.52
3:B:316:VAL:O	3:B:320:GLN:HG3	2.10	0.52
3:A:160:PRO:O	3:A:161:ASN:O	2.28	0.52
3:A:199:ILE:HD11	3:A:208:ILE:HG21	1.91	0.51
3:A:294:ASP:HB3	3:A:328:ARG:NH2	2.26	0.51
3:B:254:ASN:HD21	3:B:331:ARG:HD3	1.73	0.51
3:B:36:ARG:HH12	3:B:331:ARG:CD	2.24	0.51
3:B:146:LYS:HE2	3:B:171:VAL:HG11	1.93	0.51
3:B:246:GLY:HA3	3:B:336:ARG:NH1	2.25	0.51
3:B:186:ILE:HD12	3:B:190:THR:HG21	1.92	0.51
3:A:147:ASN:HD21	3:A:150:PHE:HD2	1.60	0.50
3:A:290:THR:OG1	3:A:292:ASP:OD2	2.29	0.50
3:A:280:ILE:HB	3:A:341:ILE:HG13	1.91	0.50
3:B:180:ILE:HG23	3:B:181:ALA:N	2.25	0.50
3:B:256:ARG:HB2	3:B:256:ARG:CZ	2.42	0.50
3:B:278:LYS:HE2	3:B:278:LYS:CA	2.36	0.50
3:A:178:LEU:O	3:A:201:LYS:HD2	2.12	0.50
3:A:304:HIS:C	3:A:304:HIS:CD2	2.90	0.50
3:A:310:THR:O	3:A:314:GLU:HB2	2.12	0.50
3:A:16:GLU:OE2	3:A:81:TYR:OH	2.29	0.50
3:B:278:LYS:HA	3:B:278:LYS:CE	2.29	0.49
3:A:10:TYR:O	3:A:10:TYR:CD1	2.65	0.49
3:A:12:TYR:HB2	3:A:45:THR:CG2	2.41	0.49
2:D:5:DG:H2'	3:A:332:ARG:NH2	2.28	0.49
3:B:253:ARG:O	3:B:254:ASN:C	2.55	0.49
3:A:289:VAL:HB	3:A:332:ARG:HB2	1.95	0.49
3:B:248:ILE:CD1	3:B:332:ARG:HE	2.25	0.49
3:A:294:ASP:CB	3:A:328:ARG:NH1	2.50	0.49
3:B:32:VAL:HG22	3:B:76:MET:HE2	1.95	0.49
3:A:200:ASN:O	3:A:201:LYS:HD3	2.13	0.49
3:B:254:ASN:CA	3:B:330:ILE:O	2.61	0.49
3:A:169:GLU:O	3:A:172:LYS:HB2	2.13	0.48
3:B:108:TYR:OH	3:B:152:LYS:HD2	2.14	0.48
3:B:300:ARG:HD3	3:B:314:GLU:OE1	2.14	0.48
3:A:157:MET:HE2	3:A:166:ILE:HD11	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:264:TYR:HB2	3:A:333:ILE:HG21	1.96	0.48
1:C:7:DC:OP1	3:A:298:ARG:HD3	2.13	0.48
3:B:256:ARG:HB2	3:B:256:ARG:NH1	2.28	0.48
3:A:245:ILE:HG21	3:A:275:LYS:HB3	1.96	0.48
3:A:10:TYR:CE1	3:A:14:GLN:HG3	2.48	0.48
3:B:179:ASP:O	3:B:180:ILE:C	2.57	0.48
3:A:266:PHE:HA	3:A:269:ILE:HD12	1.95	0.47
3:B:68:LEU:HD13	3:B:71:ALA:HB2	1.96	0.47
3:A:190:THR:HA	3:A:193:LYS:HB2	1.96	0.47
3:B:122:TYR:CD2	3:B:126:LEU:HD11	2.49	0.47
3:A:160:PRO:O	3:A:161:ASN:C	2.55	0.47
3:B:36:ARG:HA	3:B:252:LYS:HE3	1.97	0.47
3:B:101:ILE:HD12	3:B:101:ILE:N	2.29	0.47
1:C:10:DA:H5''	3:A:189:ILE:HG21	1.96	0.47
3:A:26:LYS:CB	3:A:27:PRO:HD2	2.45	0.47
2:F:11:DA:H2'	2:F:11:DA:OP2	2.15	0.47
3:B:251:MET:HA	3:B:264:TYR:CE1	2.49	0.47
3:A:193:LYS:HZ2	3:A:216:MET:HE3	1.80	0.47
3:A:146:LYS:NZ	3:A:231:ASP:OD2	2.44	0.46
3:A:294:ASP:CG	3:A:328:ARG:HH22	2.23	0.46
3:B:289:VAL:HB	3:B:332:ARG:HB2	1.98	0.46
3:B:221:LYS:O	3:B:224:TYR:HB3	2.15	0.46
3:B:116:ARG:HB3	3:B:116:ARG:CZ	2.40	0.46
3:B:122:TYR:O	3:B:125:GLY:N	2.49	0.46
3:A:319:LEU:HD12	3:A:319:LEU:O	2.16	0.46
3:B:85:SER:O	3:B:89:MET:HG2	2.15	0.46
3:A:129:LYS:HZ3	3:A:162:GLY:H	1.64	0.46
3:A:254:ASN:HB3	3:A:331:ARG:HA	1.98	0.46
3:B:49:GLU:HA	3:B:52:LYS:HE2	1.98	0.46
3:A:234:ASN:O	3:A:234:ASN:CG	2.58	0.46
3:A:97:GLU:H	3:A:97:GLU:CD	2.24	0.46
1:E:4:DG:H1'	1:E:5:DT:C2	2.51	0.45
3:A:292:ASP:CG	3:A:328:ARG:HH11	2.24	0.45
3:B:209:GLU:OE1	3:B:212:LYS:HG2	2.16	0.45
3:A:85:SER:O	3:A:89:MET:HG2	2.17	0.45
3:B:79:GLU:HB2	5:B:501:HOH:O	2.15	0.45
3:B:289:VAL:HG13	3:B:294:ASP:O	2.17	0.45
3:A:70:ASN:C	3:A:71:ALA:O	2.60	0.45
3:B:1:MET:CE	3:B:146:LYS:O	2.65	0.45
3:B:97:GLU:CD	3:B:97:GLU:H	2.24	0.45
3:B:276:LEU:O	3:B:277:ASP:C	2.60	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:172:LYS:HB3	3:A:176:ARG:HH21	1.81	0.45
3:A:252:LYS:HD2	3:A:252:LYS:HA	1.64	0.45
3:A:254:ASN:N	3:A:254:ASN:HD22	2.14	0.45
3:A:258:LEU:CD1	3:A:320:GLN:HG2	2.47	0.44
3:B:1:MET:HE3	3:B:147:ASN:HA	1.98	0.44
3:B:8:PHE:CZ	3:B:88:ILE:CG2	3.00	0.44
3:B:292:ASP:OD1	3:B:292:ASP:N	2.51	0.44
3:B:290:THR:HG21	3:B:328:ARG:HE	1.82	0.44
3:A:10:TYR:O	3:A:10:TYR:CG	2.69	0.44
3:A:194:LEU:HA	3:A:197:LEU:HB2	2.00	0.44
3:B:19:LEU:HD12	3:B:19:LEU:HA	1.80	0.44
3:B:178:LEU:O	3:B:201:LYS:HD3	2.18	0.44
3:A:321:LYS:HB2	3:A:321:LYS:HE3	1.64	0.44
3:B:9:ASP:HB3	3:B:10:TYR:CD2	2.53	0.44
3:B:124:LEU:HA	3:B:124:LEU:HD12	1.58	0.43
3:B:269:ILE:HG13	3:B:335:VAL:HG11	2.01	0.43
3:B:278:LYS:O	3:B:279:ARG:C	2.60	0.43
3:A:129:LYS:NZ	3:A:162:GLY:H	2.16	0.43
3:A:281:PRO:HA	3:A:340:PHE:HA	2.00	0.43
3:A:300:ARG:HG2	3:A:302:PHE:CE1	2.53	0.43
3:B:236:PRO:HG2	3:B:238:ARG:HH11	1.83	0.43
3:B:278:LYS:CA	3:B:278:LYS:CE	2.96	0.43
3:A:273:TYR:CD1	3:A:276:LEU:HD12	2.53	0.43
1:C:9:DT:C2'	1:C:10:DA:OP2	2.61	0.43
2:F:8:DT:H2''	2:F:9:DA:C8	2.54	0.43
3:A:29:VAL:HG22	3:A:55:VAL:HG11	2.00	0.43
3:A:10:TYR:CZ	3:A:14:GLN:HG3	2.54	0.42
3:A:261:ILE:HG21	3:A:319:LEU:HD21	2.01	0.42
3:A:300:ARG:NH1	3:A:314:GLU:OE1	2.48	0.42
3:A:242:ARG:NH1	3:A:242:ARG:CG	2.71	0.42
3:A:194:LEU:HD23	3:A:197:LEU:HD12	2.00	0.42
3:A:214:LYS:HD2	3:A:219:GLU:OE2	2.18	0.42
3:B:12:TYR:N	3:B:12:TYR:CD1	2.88	0.42
3:B:275:LYS:O	3:B:279:ARG:NH2	2.51	0.42
3:A:44:ALA:C	3:A:45:THR:HG22	2.44	0.42
3:B:36:ARG:HB2	3:B:37:PHE:H	1.60	0.42
3:B:291:GLU:HG3	3:B:292:ASP:OD1	2.19	0.42
3:A:241:VAL:HB	3:A:243:LYS:NZ	2.35	0.42
3:A:292:ASP:O	3:A:293:LEU:CB	2.68	0.42
3:B:81:TYR:CD1	3:B:81:TYR:N	2.86	0.42
3:A:252:LYS:HE3	3:B:83:GLN:NE2	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:199:ILE:HG21	3:B:205:THR:HG22	2.01	0.42
3:B:272:SER:HB2	3:B:337:PHE:CE2	2.55	0.42
3:A:190:THR:HG23	3:A:217:ILE:CG2	2.49	0.42
3:B:254:ASN:HA	3:B:330:ILE:O	2.20	0.42
3:A:197:LEU:HD23	3:A:197:LEU:HA	1.92	0.41
3:B:103:SER:OG	3:B:106:GLU:HG2	2.20	0.41
3:B:70:ASN:O	3:B:70:ASN:ND2	2.53	0.41
3:B:270:GLU:OE1	3:B:308:LYS:NZ	2.50	0.41
3:B:245:ILE:CD1	3:B:279:ARG:HH21	2.34	0.41
3:A:122:TYR:HD1	3:A:165:VAL:HG23	1.85	0.41
3:A:36:ARG:HH22	3:A:254:ASN:HB3	1.86	0.41
3:A:81:TYR:CD1	3:A:81:TYR:N	2.85	0.41
3:A:270:GLU:OE2	3:A:308:LYS:HD2	2.21	0.41
3:B:183:VAL:O	3:B:186:ILE:HB	2.21	0.41
3:B:199:ILE:HG23	3:B:204:ASP:HB2	2.03	0.41
3:A:292:ASP:OD1	3:A:328:ARG:HD2	2.21	0.41
3:B:180:ILE:HD13	3:B:202:LEU:HA	2.02	0.41
3:A:183:VAL:O	3:A:186:ILE:HB	2.20	0.40
3:A:281:PRO:HG3	3:A:337:PHE:HB3	2.03	0.40
3:B:50:ALA:O	3:B:51:ARG:C	2.64	0.40
3:B:251:MET:CE	3:B:330:ILE:CG2	2.91	0.40
3:B:293:LEU:HD23	3:B:293:LEU:HA	1.85	0.40
3:B:38:GLU:O	3:B:39:ASP:CB	2.56	0.40
3:B:141:THR:HA	3:B:162:GLY:O	2.22	0.40
3:A:20:ASN:HB3	3:A:23:LEU:HG	2.04	0.40
3:B:36:ARG:HG3	3:B:40:SER:OG	2.20	0.40
1:C:12:DC:H2''	1:C:13:DC:H5'	2.03	0.40
3:B:6:VAL:O	3:B:106:GLU:HA	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	
3	A	340/342 (99%)	302 (89%)	26 (8%)	12 (4%)	3	4
3	B	339/342 (99%)	304 (90%)	31 (9%)	4 (1%)	10	20
All	All	679/684 (99%)	606 (89%)	57 (8%)	16 (2%)	4	8

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	161	ASN
3	A	295	ILE
3	A	26	LYS
3	A	216	MET
3	B	71	ALA
3	B	277	ASP
3	A	36	ARG
3	A	1	MET
3	A	10	TYR
3	A	293	LEU
3	B	292	ASP
3	A	71	ALA
3	A	252	LYS
3	A	274	TYR
3	A	276	LEU
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	301/301 (100%)	268 (89%)	33 (11%)	6	13
3	B	300/301 (100%)	267 (89%)	33 (11%)	6	13
All	All	601/602 (100%)	535 (89%)	66 (11%)	6	13

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	1	MET
3	A	19	LEU
3	A	24	LYS
3	A	26	LYS
3	A	38	GLU
3	A	45	THR
3	A	81	TYR
3	A	93	ARG
3	A	94	GLU
3	A	106	GLU
3	A	144	ILE
3	A	152	LYS
3	A	178	LEU
3	A	180	ILE
3	A	200	ASN
3	A	207	SER
3	A	226	ILE
3	A	237	ILE
3	A	241	VAL
3	A	242	ARG
3	A	243	LYS
3	A	248	ILE
3	A	249	VAL
3	A	252	LYS
3	A	253	ARG
3	A	254	ASN
3	A	258	LEU
3	A	267	ARG
3	A	271	GLU
3	A	284	ILE
3	A	322	ILE
3	A	324	GLU
3	A	335	VAL
3	B	19	LEU
3	B	32	VAL
3	B	36	ARG
3	B	45	THR
3	B	62	VAL
3	B	68	LEU
3	B	97	GLU
3	B	134	GLU
3	B	146	LYS
3	B	169	GLU

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Mol	Chain	Res	Type
3	B	177	GLU
3	B	192	GLU
3	B	201	LYS
3	B	203	VAL
3	B	208	ILE
3	B	216	MET
3	B	221	LYS
3	B	227	SER
3	B	242	ARG
3	B	245	ILE
3	B	248	ILE
3	B	249	VAL
3	B	253	ARG
3	B	254	ASN
3	B	255	SER
3	B	257	ASN
3	B	278	LYS
3	B	292	ASP
3	B	294	ASP
3	B	308	LYS
3	B	331	ARG
3	B	336	ARG
3	B	341	ILE

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

Mol	Chain	Res	Type
3	A	14	GLN
3	A	161	ASN
3	A	200	ASN
3	A	254	ASN
3	A	304	HIS
3	B	14	GLN
3	B	70	ASN
3	B	83	GLN
3	B	234	ASN
3	B	257	ASN
3	B	304	HIS

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](#)

Of 2 ligands modelled in this entry, 2 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 [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	10/16 (62%)	-1.66	0 100 100	22, 46, 60, 62	0
1	E	10/16 (62%)	-1.51	0 100 100	21, 46, 61, 62	0
2	D	8/13 (61%)	-1.62	0 100 100	26, 29, 50, 51	0
2	F	8/13 (61%)	-1.64	0 100 100	25, 28, 50, 51	0
3	A	342/342 (100%)	-1.38	0 100 100	10, 29, 52, 72	1 (0%)
3	B	341/342 (99%)	-1.41	0 100 100	11, 30, 51, 62	0
All	All	719/742 (96%)	-1.41	0 100 100	10, 30, 54, 72	1 (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 oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
4	CA	A	401	1/1	0.97	0.07	126,126,126,126	0
4	CA	B	401	1/1	1.00	0.04	99,99,99,99	0

6.5 Other polymers [i](#)

There are no such residues in this entry.