



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 10:02 AM GMT

PDB ID : 4AA6
Title : THE OESTROGEN RECEPTOR RECOGNIZES AN IMPERFECTLY
PALINDROMIC RESPONSE ELEMENT THROUGH AN ALTERNATIVE
SIDE-CHAIN CONFORMATION
Authors : Schwabe, J.W.; Chapman, L.; Rhodes, D.
Deposited on : 2011-11-30
Resolution : 2.60 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

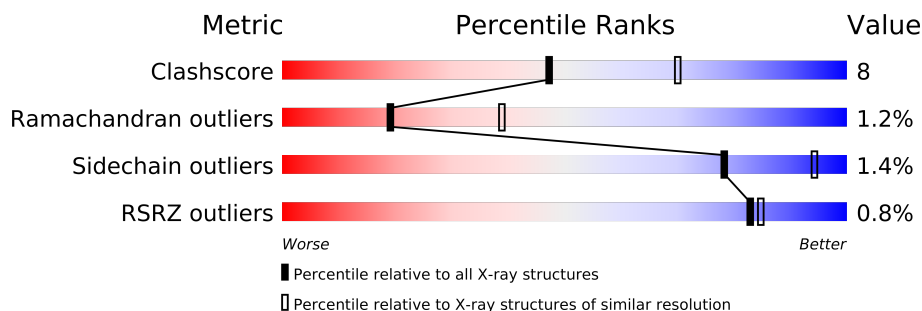
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.60 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))
Clashscore	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)
RSRZ outliers	66119	1718 (2.60-2.60)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	71	
1	B	71	
1	E	71	
1	F	71	
2	C	18	
2	G	18	
3	D	18	
3	H	18	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4615 atoms, of which 886 are hydrogens and 0 are deuterium.

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 ESTROGEN RECEPTOR.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	69	Total	C	H	N	O	S	0	0	0
			675	335	129	103	97	11			
1	B	70	Total	C	H	N	O	S	0	0	0
			678	338	127	103	98	12			
1	E	68	Total	C	H	N	O	S	0	0	0
			609	306	108	92	92	11			
1	F	69	Total	C	H	N	O	S	0	0	0
			649	326	120	97	94	12			

- Molecule 2 is a DNA chain called 5'-D(*CP*TP*AP*AP*GP*TP*CP*AP*CP*AP*GP*TP*GP*AP*CP*CP*TP*G)-3'.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	C	18	Total	C	H	N	O	P	0	0	0
			403	175	38	68	105	17			
2	G	18	Total	C	H	N	O	P	0	0	0
			403	175	38	68	105	17			

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

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	D	18	Total	C	H	N	O	P	0	0	0
			401	176	36	64	108	17			
3	H	18	Total	C	H	N	O	P	0	0	0
			401	176	36	64	108	17			

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	Zn	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Zn	0	0
			2	2		
4	F	2	Total	Zn	0	0
			2	2		
4	E	2	Total	Zn	0	0
			2	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	35	Total	H	O	0	0
			103	68	35		
5	B	18	Total	H	O	0	0
			52	34	18		
5	C	9	Total	H	O	0	0
			25	16	9		
5	D	10	Total	H	O	0	0
			28	18	10		
5	E	25	Total	H	O	0	0
			69	44	25		
5	F	19	Total	H	O	0	0
			57	38	19		
5	G	12	Total	H	O	0	0
			36	24	12		
5	H	6	Total	H	O	0	0
			18	12	6		

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: ESTROGEN RECEPTOR

Chain A: 



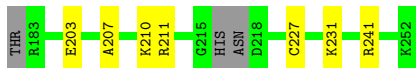
- Molecule 1: ESTROGEN RECEPTOR

Chain B: 



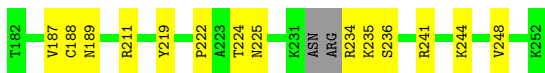
- Molecule 1: ESTROGEN RECEPTOR

Chain E: 



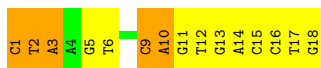
- Molecule 1: ESTROGEN RECEPTOR

Chain F: 



- Molecule 2: 5'-D(*CP*TP*AP*AP*GP*TP*CP*AP*CP*AP*GP*TP*GP*AP*CP*CP*TP*G)-3'

Chain C: 



- Molecule 2: 5'-D(*CP*TP*AP*AP*GP*TP*CP*AP*CP*AP*GP*TP*GP*AP*CP*CP*TP*G)-3'

Chain G: 



- Molecule 3: 5'-D(*TP*CP*AP*GP*GP*TP*CP*AP*CP*TP*GP*TP*GP*AP*CP*TP*TP*A)-3'

Chain D: 

T1	C2	A3	G4	G5	T6	C7	A8	C9	T10	G11	T12	G13	A14	C15	T16	T17	A18
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- Molecule 3: 5'-D(*TP*CP*AP*GP*GP*TP*CP*AP*CP*TP*GP*TP*GP*AP*CP*TP*TP*A)-3'

Chain H: 

T1	C2	A3	G4	G5	T6	C7	A8	C9	T10	G11	T12	G13	A14	C15	T16	T17	A18
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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	121.68Å 113.09Å 62.36Å 90.00° 117.45° 90.00°	Depositor
Resolution (Å)	7.00 – 2.60 15.85 – 2.51	Depositor EDS
% Data completeness (in resolution range)	99.5 (7.00-2.60) 99.1 (15.85-2.51)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.79 (at 2.52Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.222 , (Not available) 0.210 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	34.0	Xtriage
Anisotropy	0.691	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 43.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	3 of 25463 reflections (0.012%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4615	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 41.27 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.4214e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.37	0/557	0.54	0/744
1	B	0.36	0/562	0.56	0/751
1	E	0.34	0/509	0.55	0/676
1	F	0.38	0/539	0.57	0/718
2	C	1.80	5/409 (1.2%)	2.53	34/629 (5.4%)
2	G	1.80	7/409 (1.7%)	2.40	36/629 (5.7%)
3	D	1.82	4/408 (1.0%)	2.74	44/628 (7.0%)
3	H	1.72	5/408 (1.2%)	2.67	43/628 (6.8%)
All	All	1.20	21/3801 (0.6%)	1.81	157/5403 (2.9%)

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	17	DT	C5-C7	7.28	1.54	1.50
3	D	12	DT	O3'-P	-7.27	1.52	1.61
3	D	5	DG	C5'-C4'	-7.24	1.43	1.51
2	C	5	DG	C5'-C4'	-6.98	1.43	1.51
2	C	17	DT	C5-C7	6.87	1.54	1.50
3	H	7	DC	N1-C6	-6.78	1.33	1.37
3	H	6	DT	C5-C7	6.73	1.54	1.50
2	G	2	DT	C5-C7	6.73	1.54	1.50
2	G	7	DC	N1-C6	-6.46	1.33	1.37
3	H	9	DC	N1-C6	-6.23	1.33	1.37
2	C	3	DA	P-O5'	-5.96	1.53	1.59
3	H	12	DT	C5-C7	5.64	1.53	1.50
2	C	13	DG	C5-C4	-5.50	1.34	1.38
2	G	5	DG	N9-C4	5.43	1.42	1.38
2	G	10	DA	N9-C4	-5.41	1.34	1.37
3	D	12	DT	C5-C7	5.40	1.53	1.50
2	C	10	DA	C5-C4	-5.34	1.35	1.38
3	H	1	DT	C5-C7	5.31	1.53	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	15	DC	P-O5'	5.30	1.65	1.59
3	D	15	DC	N1-C2	-5.19	1.34	1.40
2	G	13	DG	C6-N1	-5.11	1.35	1.39

All (157) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	16	DT	C6-C5-C7	-12.42	115.45	122.90
3	D	6	DT	C6-C5-C7	-11.70	115.88	122.90
3	H	12	DT	O4'-C1'-C2'	-10.18	97.76	105.90
3	D	1	DT	P-O3'-C3'	10.12	131.84	119.70
3	H	10	DT	O4'-C1'-C2'	-10.11	97.81	105.90
3	D	1	DT	C6-C5-C7	-10.04	116.88	122.90
3	D	12	DT	O4'-C1'-C2'	-9.80	98.06	105.90
3	D	4	DG	O4'-C1'-C2'	-9.78	98.08	105.90
3	H	12	DT	C6-C5-C7	-9.77	117.04	122.90
3	H	10	DT	N3-C2-O2	-9.70	116.48	122.30
3	H	4	DG	P-O3'-C3'	9.44	131.03	119.70
2	C	9	DC	O4'-C1'-C2'	-9.30	98.46	105.90
3	H	16	DT	C6-C5-C7	-9.12	117.43	122.90
3	H	1	DT	C6-C5-C7	-9.01	117.50	122.90
2	C	12	DT	O4'-C1'-C2'	-8.97	98.73	105.90
3	H	5	DG	P-O3'-C3'	8.97	130.46	119.70
3	D	14	DA	P-O3'-C3'	8.86	130.33	119.70
3	H	15	DC	C1'-O4'-C4'	-8.85	101.25	110.10
3	D	15	DC	O4'-C1'-C2'	-8.78	98.88	105.90
3	D	7	DC	O4'-C1'-C2'	-8.75	98.90	105.90
2	C	2	DT	O4'-C1'-C2'	-8.56	99.05	105.90
2	G	6	DT	C6-C5-C7	-8.54	117.78	122.90
2	G	2	DT	O4'-C1'-C2'	-8.51	99.09	105.90
3	D	5	DG	P-O5'-C5'	-8.50	107.30	120.90
3	H	1	DT	P-O3'-C3'	8.45	129.84	119.70
3	D	12	DT	C6-C5-C7	-8.44	117.83	122.90
2	G	2	DT	O4'-C1'-N1	8.29	113.80	108.00
3	D	17	DT	C6-C5-C7	-8.25	117.95	122.90
2	G	15	DC	P-O3'-C3'	8.21	129.55	119.70
2	G	5	DG	P-O3'-C3'	8.19	129.52	119.70
2	C	6	DT	C6-C5-C7	-8.17	118.00	122.90
2	C	9	DC	C1'-O4'-C4'	-7.99	102.11	110.10
3	D	10	DT	C6-C5-C7	-7.89	118.17	122.90
3	H	6	DT	C6-C5-C7	-7.86	118.19	122.90
2	C	1	DC	P-O3'-C3'	7.80	129.07	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	17	DT	O4'-C1'-C2'	-7.78	99.68	105.90
2	G	15	DC	O4'-C1'-C2'	-7.75	99.70	105.90
2	G	9	DC	C1'-O4'-C4'	-7.75	102.35	110.10
3	D	9	DC	C1'-O4'-C4'	-7.72	102.38	110.10
3	H	9	DC	C1'-O4'-C4'	-7.67	102.43	110.10
3	D	9	DC	O4'-C1'-C2'	-7.59	99.83	105.90
3	D	10	DT	O4'-C1'-C2'	-7.55	99.86	105.90
3	H	15	DC	O4'-C1'-C2'	-7.49	99.91	105.90
2	C	18	DG	N9-C4-C5	7.48	108.39	105.40
2	G	9	DC	P-O3'-C3'	7.35	128.52	119.70
3	D	15	DC	P-O3'-C3'	7.28	128.44	119.70
3	H	16	DT	O4'-C1'-C2'	-7.24	100.11	105.90
2	G	4	DA	O4'-C1'-C2'	-7.20	100.14	105.90
3	H	2	DC	O4'-C1'-N1	7.19	113.03	108.00
3	H	10	DT	C6-C5-C7	-7.17	118.60	122.90
2	G	12	DT	O4'-C1'-C2'	-7.14	100.19	105.90
2	C	9	DC	P-O3'-C3'	7.10	128.22	119.70
2	G	15	DC	N1-C2-O2	7.08	123.15	118.90
2	C	5	DG	C4'-C3'-C2'	-7.05	96.75	103.10
2	C	12	DT	C6-C5-C7	-7.05	118.67	122.90
2	C	15	DC	O4'-C1'-C2'	-7.04	100.27	105.90
2	C	15	DC	P-O3'-C3'	7.02	128.13	119.70
2	C	16	DC	P-O3'-C3'	6.96	128.05	119.70
2	C	1	DC	C2-N3-C4	6.95	123.38	119.90
3	D	16	DT	C4-C5-C7	6.94	123.16	119.00
2	C	9	DC	N3-C2-O2	-6.92	117.06	121.90
2	G	6	DT	C4-C5-C6	6.91	122.15	118.00
3	D	17	DT	C4-C5-C6	6.84	122.10	118.00
2	G	5	DG	C2-N3-C4	6.77	115.29	111.90
2	C	11	DG	P-O3'-C3'	6.76	127.81	119.70
2	C	10	DA	O4'-C1'-N9	6.76	112.73	108.00
2	G	1	DC	C2-N3-C4	6.76	123.28	119.90
2	G	1	DC	N1-C2-O2	6.70	122.92	118.90
2	G	6	DT	O4'-C1'-C2'	-6.69	100.55	105.90
2	G	17	DT	O4'-C1'-C2'	-6.69	100.55	105.90
2	C	6	DT	O4'-C1'-N1	-6.68	103.33	108.00
2	G	5	DG	N3-C4-C5	-6.66	125.27	128.60
2	G	18	DG	N9-C4-C5	6.65	108.06	105.40
3	H	4	DG	N9-C4-C5	6.64	108.06	105.40
2	C	14	DA	N9-C4-C5	6.62	108.45	105.80
3	D	10	DT	N3-C2-O2	-6.61	118.33	122.30
3	H	16	DT	C4-C5-C6	6.50	121.90	118.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	14	DA	P-O3'-C3'	6.50	127.50	119.70
3	H	9	DC	O4'-C1'-C2'	-6.41	100.78	105.90
2	C	9	DC	N1-C2-O2	6.40	122.74	118.90
3	H	3	DA	N9-C4-C5	6.37	108.35	105.80
2	C	15	DC	C1'-O4'-C4'	-6.35	103.75	110.10
3	D	11	DG	C2-N3-C4	6.29	115.04	111.90
2	C	5	DG	O4'-C1'-C2'	-6.28	100.87	105.90
3	H	17	DT	C6-C5-C7	-6.28	119.13	122.90
3	H	6	DT	O4'-C1'-C2'	-6.22	100.92	105.90
2	G	18	DG	N3-C2-N2	-6.21	115.55	119.90
3	D	1	DT	C4-C5-C6	6.15	121.69	118.00
3	D	14	DA	N9-C4-C5	6.15	108.26	105.80
2	G	12	DT	C6-C5-C7	-6.14	119.22	122.90
2	G	17	DT	N3-C2-O2	-6.13	118.62	122.30
3	H	10	DT	C4-C5-C6	6.09	121.66	118.00
2	G	12	DT	C4-C5-C6	6.09	121.65	118.00
3	D	11	DG	N9-C4-C5	6.05	107.82	105.40
2	C	5	DG	P-O5'-C5'	-6.05	111.23	120.90
2	G	17	DT	C4-C5-C6	6.04	121.62	118.00
3	D	14	DA	O3'-P-O5'	-6.00	92.61	104.00
3	D	6	DT	C4-C5-C7	5.96	122.58	119.00
3	D	6	DT	C4-C5-C6	5.90	121.54	118.00
3	D	4	DG	O4'-C1'-N9	5.85	112.10	108.00
3	D	17	DT	C5-C6-N1	-5.83	120.20	123.70
2	G	10	DA	N1-C6-N6	5.83	122.10	118.60
2	C	13	DG	N3-C2-N2	-5.81	115.83	119.90
3	H	18	DA	O4'-C1'-C2'	-5.79	101.27	105.90
2	C	6	DT	C4-C5-C6	5.77	121.46	118.00
2	C	5	DG	P-O3'-C3'	5.76	126.61	119.70
3	D	5	DG	C4'-C3'-C2'	-5.74	97.93	103.10
3	H	17	DT	C4-C5-C6	5.72	121.43	118.00
3	H	6	DT	N1-C2-N3	5.70	118.02	114.60
2	C	14	DA	P-O3'-C3'	5.68	126.51	119.70
3	D	17	DT	N3-C2-O2	-5.68	118.89	122.30
3	D	4	DG	N9-C4-C5	5.64	107.65	105.40
3	H	6	DT	N3-C2-O2	-5.63	118.92	122.30
3	D	3	DA	O4'-C1'-N9	-5.62	104.06	108.00
3	H	11	DG	O4'-C1'-C2'	-5.62	101.41	105.90
2	G	11	DG	P-O3'-C3'	5.57	126.39	119.70
3	H	11	DG	C2-N3-C4	5.54	114.67	111.90
3	D	6	DT	O4'-C1'-C2'	-5.51	101.49	105.90
3	D	6	DT	O3'-P-O5'	-5.50	93.54	104.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	5	DG	O4'-C1'-C2'	-5.49	101.51	105.90
3	H	12	DT	C4-C5-C6	5.48	121.29	118.00
2	G	5	DG	C5-N7-C8	5.47	107.04	104.30
3	H	9	DC	N1-C2-O2	5.47	122.18	118.90
3	D	16	DT	C4-C5-C6	5.47	121.28	118.00
2	C	6	DT	N3-C2-O2	-5.45	119.03	122.30
3	H	6	DT	C4-C5-C6	5.44	121.27	118.00
3	D	13	DG	C8-N9-C4	-5.40	104.24	106.40
3	H	7	DC	N3-C2-O2	-5.40	118.12	121.90
2	C	1	DC	N1-C2-O2	5.39	122.14	118.90
2	G	2	DT	C6-C5-C7	-5.36	119.68	122.90
2	C	11	DG	C5-C6-N1	5.32	114.16	111.50
3	H	14	DA	N7-C8-N9	-5.32	111.14	113.80
3	D	8	DA	N9-C4-C5	5.30	107.92	105.80
2	C	11	DG	O4'-C4'-C3'	5.30	109.18	106.00
3	H	9	DC	N3-C2-O2	-5.21	118.25	121.90
2	G	1	DC	P-O3'-C3'	5.20	125.94	119.70
2	G	4	DA	P-O3'-C3'	5.19	125.92	119.70
2	C	17	DT	C4-C5-C6	5.18	121.11	118.00
3	H	2	DC	O4'-C1'-C2'	-5.17	101.76	105.90
3	H	1	DT	C4-C5-C7	5.17	122.10	119.00
3	H	15	DC	P-O3'-C3'	5.17	125.90	119.70
2	G	14	DA	P-O3'-C3'	5.16	125.89	119.70
3	D	10	DT	O3'-P-O5'	-5.14	94.23	104.00
2	G	15	DC	O4'-C1'-N1	5.13	111.59	108.00
2	G	15	DC	O4'-C4'-C3'	5.11	109.07	106.00
3	H	5	DG	P-O5'-C5'	-5.11	112.73	120.90
2	C	1	DC	O4'-C1'-N1	5.09	111.56	108.00
3	D	11	DG	N3-C4-C5	-5.09	126.05	128.60
2	G	15	DC	N3-C2-O2	-5.09	118.34	121.90
3	D	13	DG	N9-C4-C5	5.07	107.43	105.40
3	D	16	DT	P-O5'-C5'	-5.06	112.80	120.90
2	C	17	DT	C5-C6-N1	-5.06	120.66	123.70
3	H	16	DT	N3-C2-O2	-5.05	119.27	122.30
3	D	5	DG	N7-C8-N9	-5.04	110.58	113.10
3	H	17	DT	N3-C2-O2	-5.03	119.28	122.30
3	H	4	DG	O4'-C4'-C3'	5.03	109.02	106.00
2	G	5	DG	C4-C5-N7	-5.01	108.80	110.80

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	546	129	376	11	0
1	B	551	127	380	11	0
1	E	501	108	343	6	0
1	F	529	120	369	9	0
2	C	365	38	166	4	0
2	G	365	38	167	1	0
3	D	365	36	171	3	0
3	H	365	36	171	6	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
4	E	2	0	0	0	0
4	F	2	0	0	0	0
5	A	35	68	0	4	0
5	B	18	34	0	0	0
5	C	9	16	0	0	0
5	D	10	18	0	1	0
5	E	25	44	0	5	0
5	F	19	38	0	2	0
5	G	12	24	0	0	0
5	H	6	12	0	2	0
All	All	3729	886	2143	49	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

All (49) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:240:CYS:SG	1:A:243:ARG:NH1	2.62	0.72
1:E:203:GLU:HG3	5:E:2007:HOH:O	1.90	0.70
1:B:192:ALA:HB3	1:B:200:TRP:CD1	2.32	0.65
1:F:234:ARG:HD2	1:F:241:ARG:NH2	2.12	0.63
5:E:2006:HOH:H1	3:H:13:DG:H3'	1.62	0.63
1:B:234:ARG:HD2	1:B:241:ARG:NH2	2.14	0.63
2:C:2:DT:H2''	2:C:3:DA:C8	2.35	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:211:ARG:HH11	5:A:2013:HOH:H2	1.49	0.60
1:A:183:ARG:HG3	1:A:183:ARG:HH11	1.70	0.56
1:A:234:ARG:HD2	1:A:241:ARG:NH2	2.22	0.55
1:B:183:ARG:HH11	1:B:183:ARG:HG2	1.72	0.55
1:A:243:ARG:HH22	5:A:2025:HOH:H2	1.55	0.53
1:F:188:CYS:SG	1:F:234:ARG:NH1	2.82	0.51
3:H:17:DT:H6	3:H:17:DT:H5'	1.76	0.51
1:B:195:TYR:HD1	1:B:200:TRP:HE3	1.60	0.50
1:F:234:ARG:HB2	1:F:241:ARG:NE	2.27	0.50
1:A:211:ARG:HH22	5:A:2011:HOH:H2	1.60	0.49
1:B:209:PHE:O	1:B:213:ILE:HG12	2.13	0.48
2:G:2:DT:H2''	2:G:3:DA:C8	2.50	0.47
3:H:16:DT:H2''	5:H:2006:HOH:H2	1.79	0.47
1:B:196:HIS:HE1	5:D:2002:HOH:O	1.96	0.46
1:B:183:ARG:NH1	1:B:183:ARG:HG2	2.30	0.46
1:E:211:ARG:HH22	5:E:2008:HOH:H2	1.61	0.46
1:A:210:LYS:NZ	5:A:2012:HOH:H1	2.13	0.46
1:E:227:CYS:HB3	1:F:222:PRO:HB2	1.98	0.46
1:F:235:LYS:HA	5:F:2017:HOH:O	2.15	0.46
1:B:240:CYS:SG	1:B:243:ARG:NH1	2.88	0.46
1:F:224:THR:O	1:F:225:ASN:HB2	2.16	0.46
1:B:186:ALA:O	1:B:248:VAL:HG11	2.16	0.46
1:A:183:ARG:HG3	1:A:183:ARG:NH1	2.32	0.45
1:E:241:ARG:HH21	5:E:2002:HOH:H1	1.63	0.45
2:C:1:DC:H2'	2:C:2:DT:C6	2.52	0.45
3:H:2:DC:H2''	3:H:3:DA:C8	2.52	0.45
3:D:10:DT:H2''	3:D:11:DG:C8	2.52	0.45
1:F:244:LYS:O	1:F:248:VAL:HG22	2.18	0.44
2:C:9:DC:H2''	2:C:10:DA:C8	2.53	0.44
1:E:207:ALA:HA	1:E:210:LYS:HE2	2.00	0.44
1:A:211:ARG:NH1	3:D:13:DG:N7	2.65	0.43
1:E:231:LYS:HE3	5:E:2003:HOH:H2	1.82	0.43
3:D:1:DT:H2'	3:D:2:DC:C6	2.54	0.42
1:A:228:THR:HG22	1:A:228:THR:O	2.19	0.42
1:B:188:CYS:SG	1:B:234:ARG:NH1	2.93	0.42
3:H:17:DT:H73	5:H:2006:HOH:O	2.20	0.41
1:B:187:VAL:O	1:B:244:LYS:HD3	2.21	0.41
1:F:187:VAL:HG13	1:F:244:LYS:HB3	2.03	0.41
1:F:211:ARG:HH22	5:F:2003:HOH:H1	1.67	0.41
2:C:2:DT:H2''	2:C:3:DA:H8	1.85	0.40
3:H:17:DT:H5'	3:H:17:DT:C6	2.54	0.40
1:A:244:LYS:O	1:A:248:VAL:HG22	2.22	0.40

There are no symmetry-related clashes.

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	65/71 (92%)	63 (97%)	2 (3%)	0	100	100
1	B	66/71 (93%)	60 (91%)	5 (8%)	1 (2%)	15	30
1	E	64/71 (90%)	61 (95%)	3 (5%)	0	100	100
1	F	65/71 (92%)	56 (86%)	7 (11%)	2 (3%)	7	10
All	All	260/284 (92%)	240 (92%)	17 (6%)	3 (1%)	19	39

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	219	TYR
1	B	183	ARG
1	F	236	SER

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	58/61 (95%)	57 (98%)	1 (2%)	73	92
1	B	58/61 (95%)	57 (98%)	1 (2%)	73	92
1	E	51/61 (84%)	51 (100%)	0	100	100
1	F	55/61 (90%)	54 (98%)	1 (2%)	71	91
All	All	222/244 (91%)	219 (99%)	3 (1%)	78	94

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	183	ARG
1	B	214	GLN
1	F	189	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	214	GLN
1	E	189	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 8 ligands modelled in this entry, 8 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.

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	A	69/71 (97%)	-0.40	1 (1%) 72 72	5, 16, 39, 63	0
1	B	70/71 (98%)	-0.20	2 (2%) 49 46	7, 22, 51, 64	0
1	E	68/71 (95%)	-0.34	0 100 100	8, 25, 47, 59	0
1	F	69/71 (97%)	-0.15	0 100 100	7, 20, 52, 59	0
2	C	18/18 (100%)	-0.87	0 100 100	10, 22, 30, 33	0
2	G	18/18 (100%)	-0.74	0 100 100	11, 27, 36, 47	0
3	D	18/18 (100%)	-0.81	0 100 100	4, 23, 29, 36	0
3	H	18/18 (100%)	-0.83	0 100 100	9, 25, 35, 46	0
All	All	348/356 (97%)	-0.38	3 (0%) 83 82	4, 22, 48, 64	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	182	THR	3.3
1	A	216	HIS	2.3
1	B	200	TRP	2.1

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	ZN	B	254	1/1	0.11	-0.41	16,16,16,16	0
4	ZN	A	253	1/1	0.10	-1.43	15,15,15,15	0
4	ZN	A	254	1/1	0.07	-1.76	15,15,15,15	0
4	ZN	F	253	1/1	0.08	-1.86	21,21,21,21	0
4	ZN	E	253	1/1	0.09	-2.00	22,22,22,22	0
4	ZN	E	254	1/1	0.05	-2.89	15,15,15,15	0
4	ZN	F	254	1/1	0.07	-3.41	16,16,16,16	0
4	ZN	B	253	1/1	0.06	-3.88	22,22,22,22	0

6.5 Other polymers

There are no such residues in this entry.