Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 01:01 PM GMT

PDB ID	:	3SN7
Title	:	Highly Potent, Selective, and Orally Active Phosphodiestarase 10A Inhibitors
Authors	:	Parris, K.D.
Deposited on	:	2011-06-28
Resolution	:	1.82 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

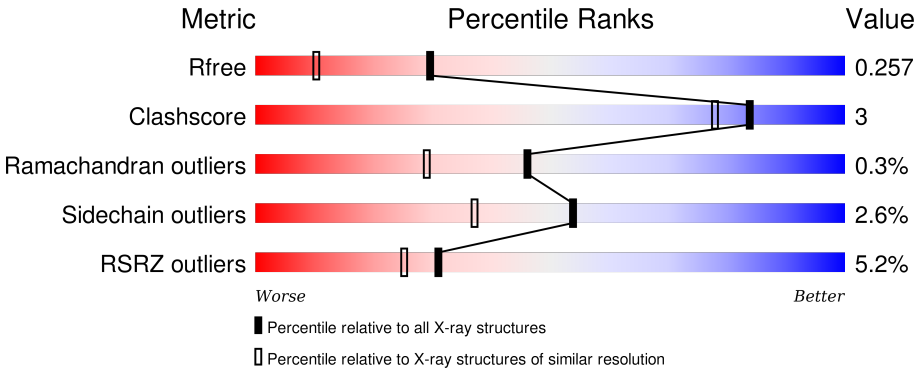
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.82 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	5422 (1.84-1.80)
Clashscore	102246	6347 (1.84-1.80)
Ramachandran outliers	100387	6276 (1.84-1.80)
Sidechain outliers	100360	6276 (1.84-1.80)
RSRZ outliers	91569	5439 (1.84-1.80)

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 <=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	345	3% 84% 8% • 7%
1	B	345	7% 82% 8% • 10%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	MG	A	782	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5469 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 cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	0	3	0
			2627	1678	446	478	25			
1	B	310	Total	C	N	O	S	0	2	0
			2530	1611	431	463	25			

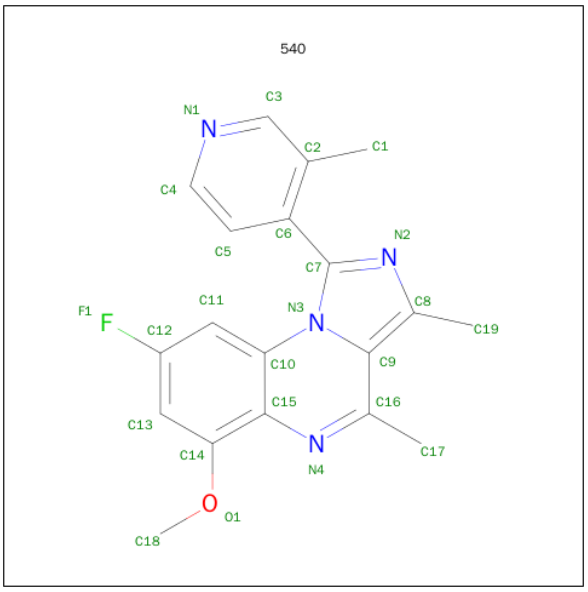
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	435	GLY	-	EXPRESSION TAG	UNP Q9Y233
A	436	SER	-	EXPRESSION TAG	UNP Q9Y233
A	437	HIS	-	EXPRESSION TAG	UNP Q9Y233
A	438	MET	-	EXPRESSION TAG	UNP Q9Y233
B	435	GLY	-	EXPRESSION TAG	UNP Q9Y233
B	436	SER	-	EXPRESSION TAG	UNP Q9Y233
B	437	HIS	-	EXPRESSION TAG	UNP Q9Y233
B	438	MET	-	EXPRESSION TAG	UNP Q9Y233

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cl	0	0
			1	1		
2	A	1	Total	Cl	0	0
			1	1		

- Molecule 3 is 8-FLUORO-6-METHOXY-3,4-DIMETHYL-1-(3-METHYLPYRIDIN-4-YL)IMIDAZO[1,5-A]QUINOXALINE (three-letter code: 540) (formula: $C_{19}H_{17}FN_4O$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C F N O 25 19 1 4 1	0	0
3	B	1	Total C F N O 25 19 1 4 1	0	0

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Zn 1 1	0	0
4	A	1	Total Zn 1 1	0	0

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Mg 1 1	0	0
5	A	1	Total Mg 1 1	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	153	Total O 153 153	0	0

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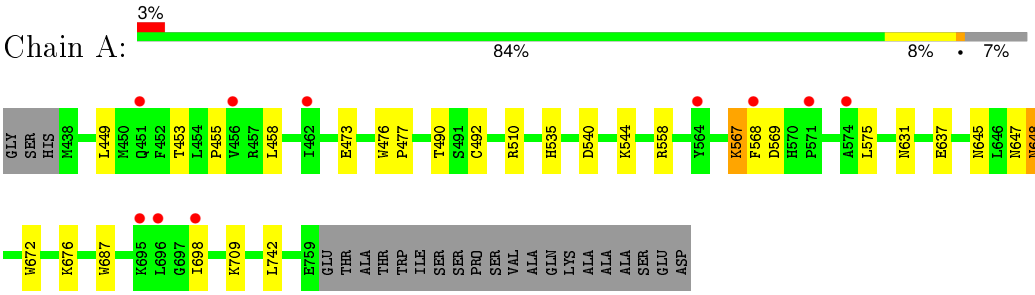
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	103	Total	O	0	0
			103	103		

3 Residue-property plots (i)

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.

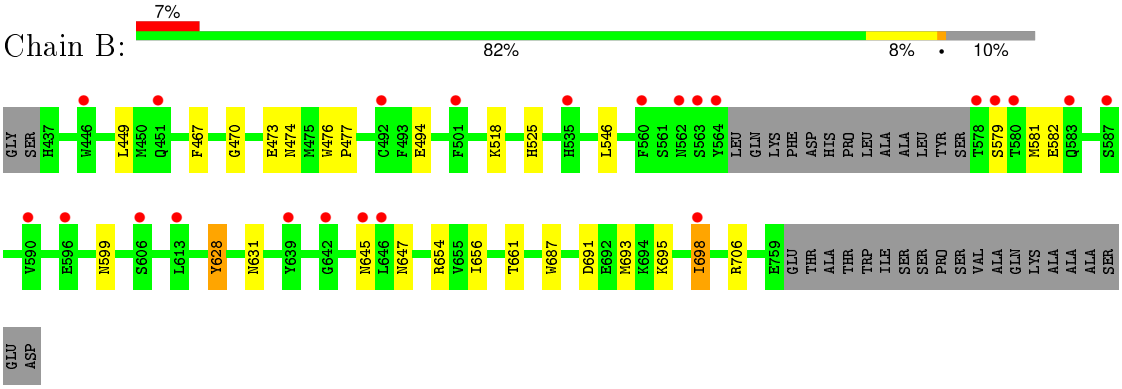
- Molecule 1: cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A

Chain A:



- Molecule 1: cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	50.64Å 81.47Å 158.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.24 – 1.82 43.01 – 1.82	Depositor EDS
% Data completeness (in resolution range)	79.3 (44.24-1.82) 79.3 (43.01-1.82)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.34 (at 1.82Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.247 , 0.276 0.217 , 0.257	Depositor DCC
R_{free} test set	2416 reflections (5.39%)	DCC
Wilson B-factor (Å^2)	27.6	Xtriage
Anisotropy	0.093	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 46.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 47257 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5469	wwPDB-VP
Average B, all atoms (Å^2)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.18% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: MG, ZN, 540, CL

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	0.51	0/2691	0.62	0/3644
1	B	0.49	0/2590	0.61	0/3504
All	All	0.50	0/5281	0.61	0/7148

There are no bond length outliers.

There are no bond angle outliers.

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	A	2627	0	2591	15	0
1	B	2530	0	2487	14	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	25	0	17	1	0
3	B	25	0	17	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	153	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	103	0	0	0	0
All	All	5469	0	5112	31	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 3.

All (31) 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:A:645:ASN:H	1:A:651:HIS:HD2	1.27	0.79
3:B:780:540:H17	3:B:780:540:H19	1.68	0.74
3:A:780:540:H17	3:A:780:540:H19	1.77	0.66
1:B:546:LEU:HD21	1:B:656:ILE:HG23	1.79	0.64
1:A:568:PHE:CZ	1:A:698:ILE:HG13	2.34	0.62
1:A:645:ASN:N	1:A:651:HIS:HD2	1.98	0.59
1:A:645:ASN:H	1:A:651:HIS:CD2	2.16	0.57
1:A:535:HIS:HD2	6:A:76:HOH:O	1.88	0.57
1:A:631:ASN:HD22	1:A:654:ARG:HH11	1.51	0.56
1:B:631:ASN:HD22	1:B:654:ARG:HH11	1.57	0.53
1:B:473:GLU:HA	1:B:476:TRP:CE2	2.43	0.53
1:B:687:TRP:HB3	1:B:706:ARG:HB3	1.89	0.52
1:A:473:GLU:HA	1:A:476:TRP:CE2	2.44	0.52
1:B:645:ASN:HD21	1:B:647:ASN:HB2	1.74	0.52
1:B:691:ASP:O	1:B:695:LYS:HG2	2.12	0.49
1:A:648:ASN:HD22	1:A:648:ASN:C	2.16	0.49
1:A:645:ASN:HD21	1:A:647:ASN:HB2	1.80	0.47
1:B:546:LEU:HD21	1:B:656:ILE:CG2	2.46	0.45
1:A:687:TRP:CZ2	1:A:709:LYS:HG2	2.51	0.45
1:B:467:PHE:HB3	1:B:525:HIS:CE1	2.51	0.45
1:A:455:PRO:HD2	1:A:458:LEU:HD12	1.98	0.44
1:B:470:GLY:HA2	1:B:518:LYS:HE3	1.99	0.44
1:B:449:LEU:HD13	1:B:477:PRO:HB2	1.98	0.44
1:A:672:TRP:CZ2	1:A:676:LYS:HD3	2.54	0.43
1:B:693:MET:HB3	1:B:698:ILE:HB	2.01	0.42
1:A:540:ASP:O	1:A:544:LYS:HG3	2.20	0.42
1:A:449:LEU:HD13	1:A:477:PRO:HB2	2.01	0.42
1:B:631:ASN:HD21	1:B:654:ARG:HA	1.85	0.42
1:A:510:ARG:HE	1:A:558:ARG:HG3	1.85	0.42
1:B:579:SER:HB2	1:B:582:GLU:HB2	2.03	0.41
1:B:628:TYR:CD1	1:B:661:THR:HG21	2.57	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	323/345 (94%)	314 (97%)	7 (2%)	2 (1%)	30 14
1	B	308/345 (89%)	296 (96%)	12 (4%)	0	100 100
All	All	631/690 (91%)	610 (97%)	19 (3%)	2 (0%)	46 29

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	569	ASP
1	A	567	LYS

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	A	293/308 (95%)	284 (97%)	9 (3%)	47 29
1	B	283/308 (92%)	277 (98%)	6 (2%)	61 47
All	All	576/616 (94%)	561 (97%)	15 (3%)	54 37

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

Mol	Chain	Res	Type
1	A	453	THR

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Mol	Chain	Res	Type
1	A	490	THR
1	A	492	CYS
1	A	567	LYS
1	A	575	LEU
1	A	637	GLU
1	A	648	ASN
1	A	651	HIS
1	A	742	LEU
1	B	474	ASN
1	B	494	GLU
1	B	581	MET
1	B	599	ASN
1	B	628	TYR
1	B	698	ILE

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

Mol	Chain	Res	Type
1	A	535	HIS
1	A	631	ASN
1	A	645	ASN
1	A	648	ASN
1	A	651	HIS
1	B	599	ASN
1	B	631	ASN
1	B	640	GLN
1	B	645	ASN
1	B	680	ASN
1	B	714	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 (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	540	A	780	-	20,28,28	3.45	14 (70%)	25,42,42	1.87	12 (48%)
3	540	B	780	-	20,28,28	3.43	12 (60%)	25,42,42	2.12	10 (40%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	540	A	780	-	-	0/6/6/6	0/4/4/4
3	540	B	780	-	-	0/6/6/6	0/4/4/4

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	780	540	C17-C16	-7.96	1.45	1.50
3	B	780	540	C17-C16	-6.86	1.46	1.50
3	A	780	540	C7-N2	-3.19	1.29	1.35
3	B	780	540	C7-N2	-2.39	1.30	1.35
3	B	780	540	C3-N1	2.10	1.38	1.34
3	B	780	540	C5-C4	2.33	1.43	1.38
3	B	780	540	C13-C12	2.61	1.42	1.37
3	A	780	540	C5-C4	2.63	1.44	1.38
3	A	780	540	C3-N1	2.63	1.40	1.34
3	A	780	540	C13-C14	2.75	1.47	1.37
3	B	780	540	C15-N4	2.79	1.43	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	780	540	C15-N4	2.81	1.43	1.37
3	A	780	540	C11-C12	2.91	1.41	1.35
3	A	780	540	C4-N1	3.02	1.42	1.33
3	A	780	540	C14-C15	3.21	1.47	1.42
3	B	780	540	C4-N1	3.34	1.43	1.33
3	B	780	540	C3-C2	3.36	1.45	1.39
3	A	780	540	C3-C2	3.57	1.45	1.39
3	A	780	540	C13-C12	3.79	1.44	1.37
3	B	780	540	C5-C6	3.92	1.46	1.39
3	A	780	540	C5-C6	4.01	1.46	1.39
3	A	780	540	C16-C9	5.33	1.50	1.42
3	A	780	540	C6-C2	5.47	1.49	1.40
3	B	780	540	C16-C9	5.64	1.50	1.42
3	B	780	540	C14-C15	5.85	1.51	1.42
3	B	780	540	C6-C2	6.65	1.51	1.40

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	780	540	C2-C3-N1	-3.66	120.51	124.77
3	B	780	540	C1-C2-C3	-3.64	115.05	120.65
3	A	780	540	C2-C3-N1	-2.93	121.36	124.77
3	A	780	540	C1-C2-C3	-2.78	116.38	120.65
3	B	780	540	F1-C12-C11	-2.76	114.79	119.07
3	A	780	540	C4-C5-C6	-2.74	115.14	118.72
3	B	780	540	C4-C5-C6	-2.63	115.28	118.72
3	A	780	540	F1-C12-C11	-2.09	115.82	119.07
3	A	780	540	C2-C6-C7	-2.05	116.09	119.75
3	B	780	540	C13-C14-C15	-2.04	117.48	120.15
3	A	780	540	F1-C12-C13	2.05	120.95	118.22
3	A	780	540	C4-N1-C3	2.16	120.81	116.84
3	A	780	540	C7-N2-C8	2.22	109.71	104.41
3	A	780	540	C14-C15-C10	2.25	119.85	118.25
3	B	780	540	F1-C12-C13	2.36	121.36	118.22
3	A	780	540	C5-C6-C2	2.47	122.70	119.24
3	B	780	540	C4-N1-C3	2.66	121.74	116.84
3	B	780	540	O1-C14-C15	2.68	118.08	115.14
3	A	780	540	C1-C2-C6	2.81	127.10	122.43
3	B	780	540	C1-C2-C6	3.18	127.73	122.43
3	A	780	540	O1-C14-C15	3.62	119.12	115.14
3	B	780	540	C18-O1-C14	3.80	123.33	117.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	780	540	1	0
3	B	780	540	1	0

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

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, 95^{th} 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	322/345 (93%)	0.13	10 (3%)	52	47	15, 31, 57, 72	0
1	B	310/345 (89%)	0.30	23 (7%)	17	14	20, 36, 66, 93	0
All	All	632/690 (91%)	0.22	33 (5%)	31	25	15, 33, 64, 93	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	696	LEU	6.6
1	B	579	SER	6.1
1	B	562	ASN	5.7
1	B	578	THR	5.1
1	B	698	ILE	4.7
1	B	646	LEU	4.2
1	B	564	TYR	4.0
1	A	698	ILE	3.8
1	A	568	PHE	3.7
1	B	563	SER	3.4
1	B	590	VAL	3.2
1	B	501	PHE	3.0
1	B	606	SER	2.9
1	A	574	ALA	2.9
1	B	583	GLN	2.8
1	B	580	THR	2.8
1	A	456	VAL	2.8
1	A	564	TYR	2.7
1	B	492	CYS	2.6
1	A	571	PRO	2.6
1	B	446	TRP	2.5
1	A	451	GLN	2.5
1	B	587	SER	2.4
1	B	639	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	560	PHE	2.3
1	B	645	ASN	2.2
1	B	535	HIS	2.2
1	A	462	ILE	2.2
1	A	695	LYS	2.1
1	B	613	LEU	2.1
1	B	451	GLN	2.1
1	B	642	GLY	2.0
1	B	596	GLU	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, 95^{th} 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
5	MG	A	782	1/1	0.98	0.14	6.08	16,16,16,16	0
2	CL	B	1	1/1	0.99	0.20	1.85	11,11,11,11	0
3	540	B	780	25/25	0.93	0.10	0.96	25,31,39,41	0
4	ZN	A	781	1/1	1.00	0.10	0.13	24,24,24,24	0
5	MG	B	782	1/1	0.97	0.10	-0.28	30,30,30,30	0
3	540	A	780	25/25	0.94	0.10	-0.33	16,24,30,31	0
4	ZN	B	781	1/1	0.99	0.07	-2.09	28,28,28,28	0
2	CL	A	1	1/1	0.95	0.17	-	38,38,38,38	0

6.5 Other polymers (i)

There are no such residues in this entry.