



wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 16, 2024 – 07:31 AM EDT

PDB ID : 4WKG
Title : The crystal structure of apo ArnA features an unexpected central binding pocket and provides an explanation for enzymatic coop-erativity
Authors : Grimm, C.
Deposited on : 2014-10-02
Resolution : 2.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.37.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.37.1

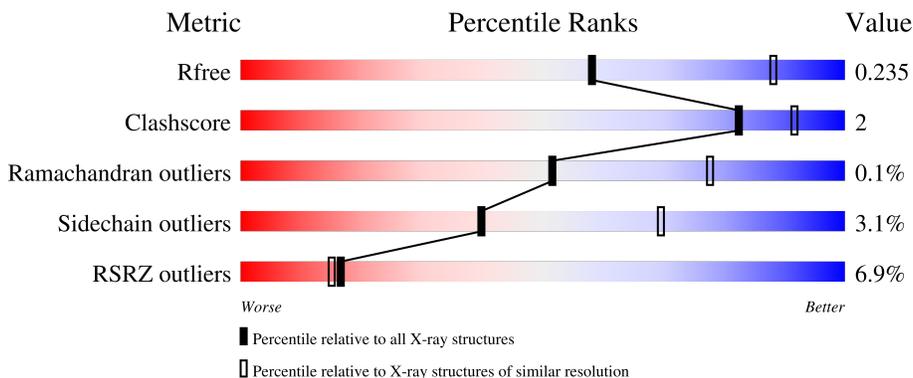
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	A	660	 4% 90% 6% . .
1	B	660	 6% 90% 7% .
1	C	660	 3% 87% 9% .
1	D	660	 18% 87% 10% .
1	E	660	 5% 88% 8% . .

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	660	 <p>5% 87% 9% ..</p>

2 Entry composition [i](#)

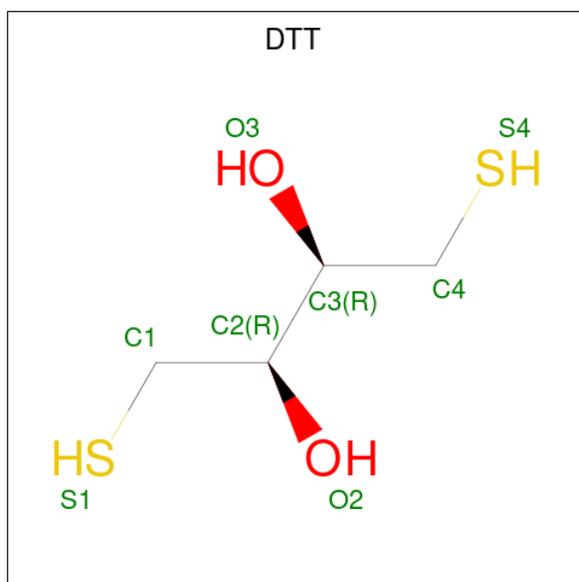
There are 4 unique types of molecules in this entry. The entry contains 60000 atoms, of which 29761 are hydrogens and 0 are deuteriums.

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 Bifunctional polymyxin resistance protein ArnA.

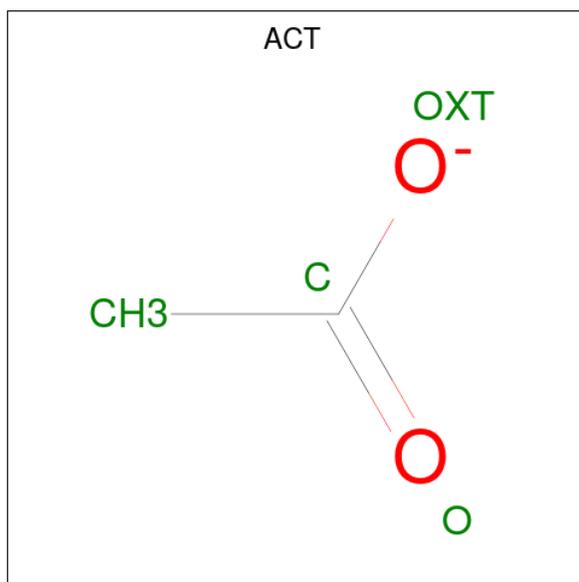
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
1	A	636	Total 9951	C 3195	H 4945	N 889	O 903	S 19	0	0	0
1	B	643	Total 10036	C 3234	H 4972	N 897	O 914	S 19	0	0	0
1	C	637	Total 9963	C 3203	H 4947	N 888	O 906	S 19	0	0	0
1	D	641	Total 9973	C 3218	H 4936	N 889	O 911	S 19	0	0	0
1	E	643	Total 10032	C 3231	H 4974	N 894	O 914	S 19	0	0	0
1	F	637	Total 9950	C 3197	H 4941	N 888	O 905	S 19	0	0	0

- Molecule 2 is 2,3-DIHYDROXY-1,4-DITHIOBUTANE (three-letter code: DTT) (formula: $C_4H_{10}O_2S_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	A	1	Total	C	H	O	S	0	1
			36	8	20	4	4		
2	D	1	Total	C	H	O	S	0	1
			36	8	20	4	4		

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			7	2	3	2		
3	F	1	Total	C	H	O	0	0
			7	2	3	2		

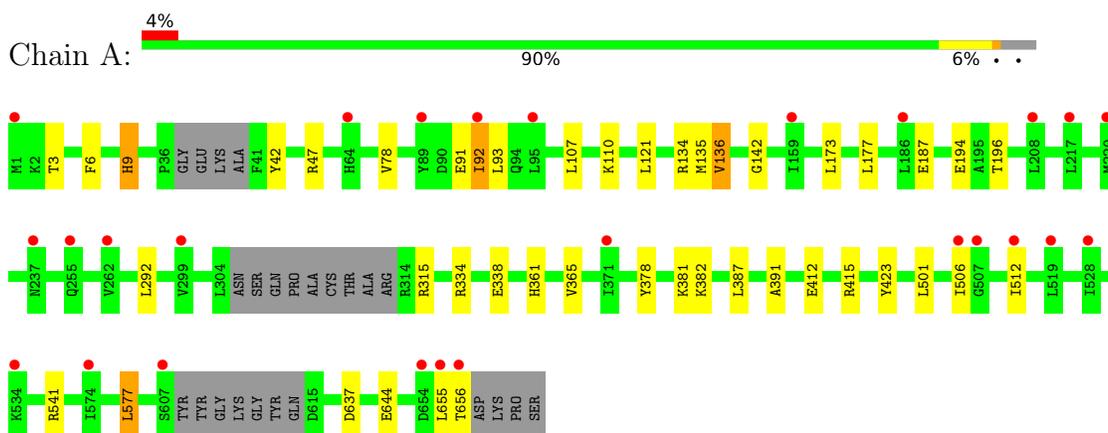
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	O	0	0
			1	1		
4	B	2	Total	O	0	0
			2	2		
4	C	2	Total	O	0	0
			2	2		
4	D	1	Total	O	0	0
			1	1		
4	E	2	Total	O	0	0
			2	2		
4	F	1	Total	O	0	0
			1	1		

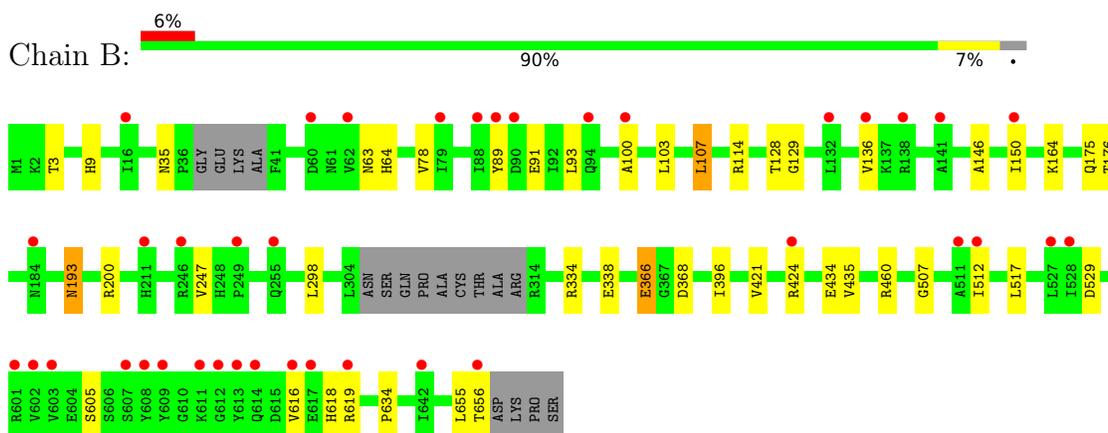
3 Residue-property plots [i](#)

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.

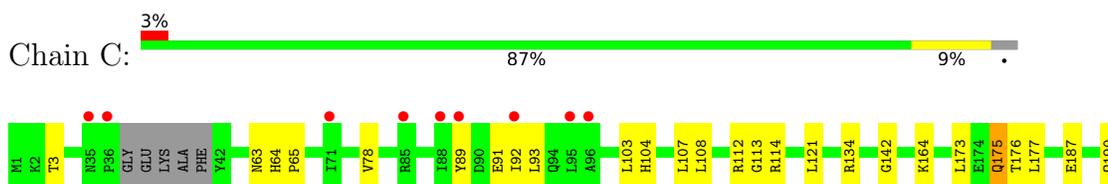
- Molecule 1: Bifunctional polymyxin resistance protein ArnA

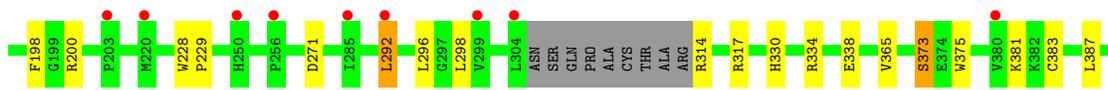


- Molecule 1: Bifunctional polymyxin resistance protein ArnA

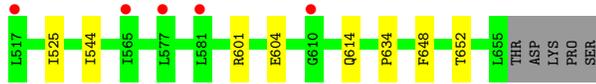
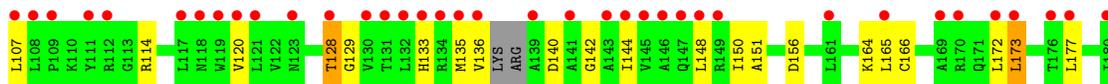
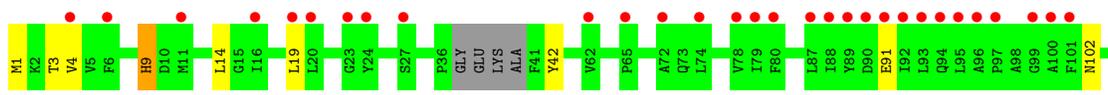
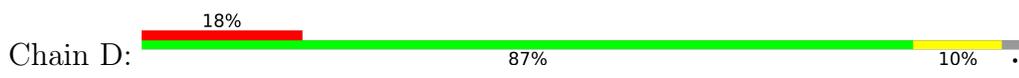


- Molecule 1: Bifunctional polymyxin resistance protein ArnA

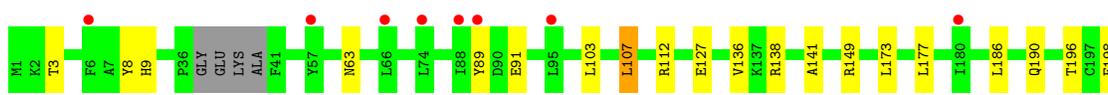
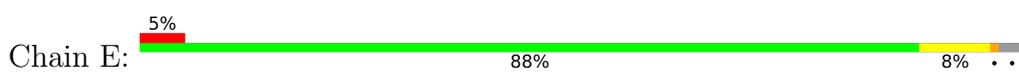




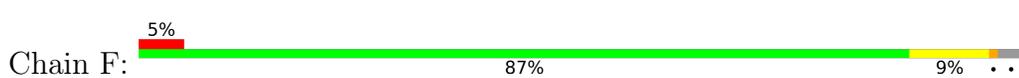
● Molecule 1: Bifunctional polymyxin resistance protein ArnA

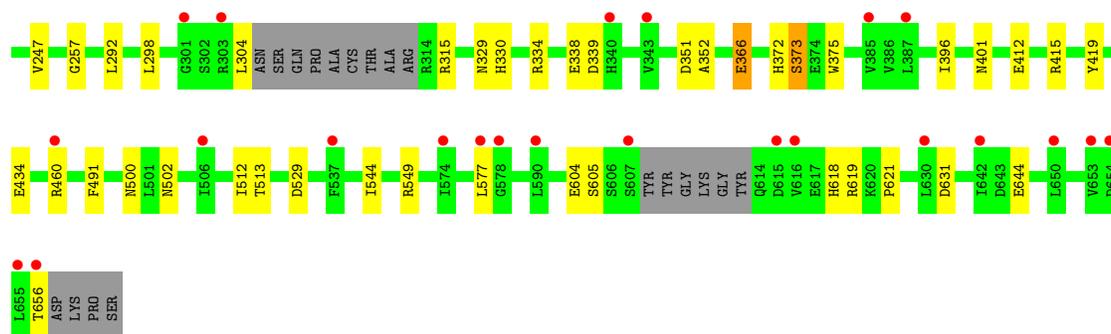


● Molecule 1: Bifunctional polymyxin resistance protein ArnA



● Molecule 1: Bifunctional polymyxin resistance protein ArnA





4 Data and refinement statistics i

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	106.72Å 112.82Å 113.61Å 81.92° 82.96° 83.80°	Depositor
Resolution (Å)	48.16 – 2.70 48.16 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.4 (48.16-2.70) 90.0 (48.16-2.70)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.08 (at 2.69Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.205 , 0.232 0.208 , 0.235	Depositor DCC
R_{free} test set	6967 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	69.0	Xtrriage
Anisotropy	0.218	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 37.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.014 for -h,-l,-k	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	60000	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.50% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.29	0/5126	0.52	1/6965 (0.0%)
1	B	0.30	0/5188	0.54	1/7051 (0.0%)
1	C	0.27	0/5137	0.51	1/6981 (0.0%)
1	D	0.29	0/5160	0.53	0/7013
1	E	0.29	0/5182	0.53	2/7044 (0.0%)
1	F	0.29	0/5129	0.52	0/6970
All	All	0.29	0/30922	0.53	5/42024 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	301	GLY	C-N-CA	6.51	137.98	121.70
1	B	507	GLY	N-CA-C	5.87	127.77	113.10
1	E	507	GLY	N-CA-C	5.35	126.47	113.10
1	A	577	LEU	CB-CG-CD2	-5.18	102.19	111.00
1	C	507	GLY	N-CA-C	5.10	125.85	113.10

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	A	5006	4945	4933	20	0
1	B	5064	4972	4977	23	0
1	C	5016	4947	4937	26	0
1	D	5037	4936	4936	33	0
1	E	5058	4974	4966	25	0
1	F	5009	4941	4930	29	0
2	A	16	20	20	0	0
2	D	16	20	20	1	0
3	A	4	3	3	0	0
3	F	4	3	3	1	0
4	A	1	0	0	0	0
4	B	2	0	0	0	0
4	C	2	0	0	0	0
4	D	1	0	0	0	0
4	E	2	0	0	0	0
4	F	1	0	0	0	0
All	All	30239	29761	29725	150	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 2.

The worst 5 of 150 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:D:434:GLU:OE1	1:D:460:ARG:NH2	2.16	0.77
1:A:382:LYS:NZ	1:B:366:GLU:OE2	2.19	0.76
1:B:434:GLU:HG2	1:B:619:ARG:HH22	1.50	0.75
1:F:500:ASN:ND2	1:F:502:ASN:OD1	2.21	0.74
1:F:415:ARG:NH2	1:F:419:TYR:OH	2.21	0.73

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	628/660 (95%)	613 (98%)	14 (2%)	1 (0%)	47	73
1	B	637/660 (96%)	623 (98%)	13 (2%)	1 (0%)	47	73
1	C	629/660 (95%)	617 (98%)	12 (2%)	0	100	100
1	D	633/660 (96%)	619 (98%)	14 (2%)	0	100	100
1	E	637/660 (96%)	619 (97%)	17 (3%)	1 (0%)	47	73
1	F	629/660 (95%)	618 (98%)	11 (2%)	0	100	100
All	All	3793/3960 (96%)	3709 (98%)	81 (2%)	3 (0%)	51	78

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	136	VAL
1	E	136	VAL
1	B	136	VAL

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	528/559 (94%)	515 (98%)	13 (2%)	47	76
1	B	532/559 (95%)	517 (97%)	15 (3%)	43	73
1	C	529/559 (95%)	513 (97%)	16 (3%)	41	70
1	D	528/559 (94%)	514 (97%)	14 (3%)	44	74
1	E	531/559 (95%)	513 (97%)	18 (3%)	37	66
1	F	528/559 (94%)	507 (96%)	21 (4%)	31	60
All	All	3176/3354 (95%)	3079 (97%)	97 (3%)	40	69

5 of 97 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	634	PRO
1	E	415	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	E	91	GLU
1	E	209	GLU
1	E	656	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	372	HIS
1	F	9	HIS
1	F	363	HIS
1	F	64	HIS
1	C	64	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

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	ACT	F	701	-	3,3,3	0.71	0	3,3,3	1.39	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	DTT	D	701[B]	-	7,7,7	0.57	0	4,8,8	0.29	0
2	DTT	D	701[A]	-	7,7,7	0.60	0	4,8,8	0.32	0
3	ACT	A	702	-	3,3,3	0.75	0	3,3,3	1.41	0
2	DTT	A	701[B]	-	7,7,7	0.61	0	4,8,8	0.35	0
2	DTT	A	701[A]	-	7,7,7	0.72	0	4,8,8	0.53	0

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
2	DTT	A	701[B]	-	-	4/8/8/8	-
2	DTT	D	701[B]	-	-	1/8/8/8	-
2	DTT	A	701[A]	-	-	2/8/8/8	-
2	DTT	D	701[A]	-	-	1/8/8/8	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701[A]	DTT	S1-C1-C2-O2
2	A	701[A]	DTT	S1-C1-C2-C3
2	A	701[B]	DTT	S1-C1-C2-O2
2	A	701[B]	DTT	S1-C1-C2-C3
2	A	701[B]	DTT	O3-C3-C4-S4

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	701	ACT	1	0
2	D	701[B]	DTT	1	0

5.7 Other polymers [\(i\)](#)

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 [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	A	636/660 (96%)	0.43	26 (4%) 37 36	19, 51, 96, 152	0
1	B	643/660 (97%)	0.50	39 (6%) 21 20	20, 56, 105, 170	0
1	C	637/660 (96%)	0.43	23 (3%) 42 42	27, 55, 103, 158	0
1	D	641/660 (97%)	0.98	116 (18%) 1 1	20, 71, 157, 204	0
1	E	643/660 (97%)	0.41	30 (4%) 31 30	16, 49, 98, 147	0
1	F	637/660 (96%)	0.47	32 (5%) 28 27	30, 60, 100, 146	0
All	All	3837/3960 (96%)	0.54	266 (6%) 16 15	16, 55, 121, 204	0

The worst 5 of 266 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	79	ILE	12.9
1	D	100	ALA	12.2
1	D	101	PHE	8.7
1	D	136	VAL	8.0
1	E	608	TYR	7.8

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

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
2	DTT	A	701[A]	8/8	0.80	0.35	32,55,55,55	18
2	DTT	A	701[B]	8/8	0.80	0.35	55,55,55,55	18
2	DTT	D	701[A]	8/8	0.81	0.28	35,64,64,64	18
2	DTT	D	701[B]	8/8	0.81	0.28	46,64,64,64	18
3	ACT	F	701	4/4	0.86	0.28	20,20,58,58	0
3	ACT	A	702	4/4	0.95	0.42	20,20,44,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.