



# wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 7, 2016 – 03:13 PM EST

PDB ID : 5EKC  
Title : Thermostable aldehyde dehydrogenase from *Pyrobaculum* sp.1860 complexed with NADP+  
Authors : Petrova, T.; Bezsudnova, E.Y.; Boyko, K.M.; Nikolaeva, A.Y.; Rakitina, T.V.; Shabalin, I.G.; Popov, V.O.  
Deposited on : 2015-11-03  
Resolution : 1.90 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20028320
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)	:	rb-20028320

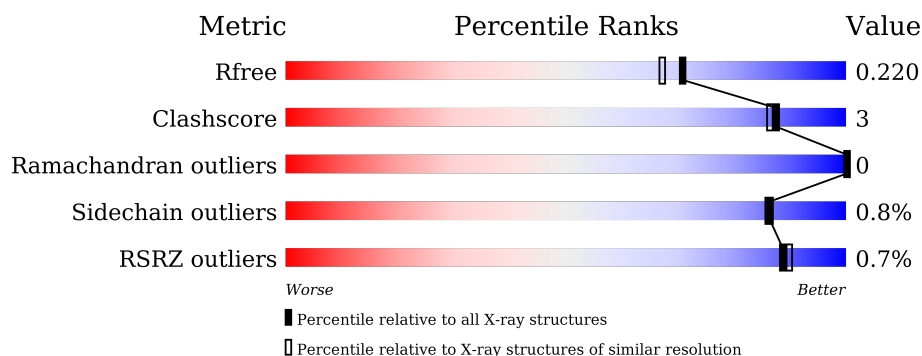
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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.



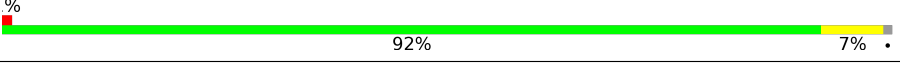
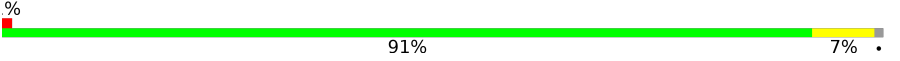
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	491	<div> <div></div> <div>93%</div> <div>6%</div> <div>.</div> </div>
1	B	491	<div> <div></div> <div>94%</div> <div></div> <div>.</div> </div>
1	C	491	<div> <div></div> <div>92%</div> <div>6%</div> <div>.</div> </div>
1	D	491	<div> <div></div> <div>92%</div> <div>7%</div> <div>.</div> </div>
1	E	491	<div> <div></div> <div>92%</div> <div>8%</div> <div></div> </div>
1	F	491	<div> <div></div> <div>92%</div> <div>6%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
1	G	491	 % 92% 7%
1	H	491	 % 91% 7%

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
2	NAP	F	500	-	-	-	X

## 2 Entry composition

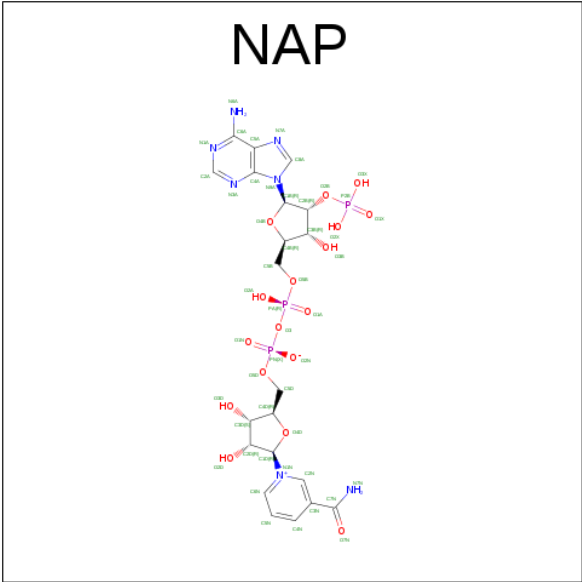
There are 3 unique types of molecules in this entry. The entry contains 33123 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Aldehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	485	Total	C	N	O	S	0	5	1
			3761	2398	651	700	12			
1	B	485	Total	C	N	O	S	0	4	0
			3775	2411	654	697	13			
1	C	483	Total	C	N	O	S	0	5	0
			3761	2399	658	693	11			
1	D	484	Total	C	N	O	S	0	6	0
			3755	2398	650	695	12			
1	E	490	Total	C	N	O	S	0	13	0
			3812	2433	662	704	13			
1	F	482	Total	C	N	O	S	0	5	0
			3760	2400	653	696	11			
1	G	484	Total	C	N	O	S	0	8	0
			3767	2407	653	694	13			
1	H	485	Total	C	N	O	S	0	6	0
			3781	2414	658	696	13			

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C<sub>21</sub>H<sub>28</sub>N<sub>7</sub>O<sub>17</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			39	15	5	16	3		
2	B	1	Total	C	N	O	P	0	0
			39	15	5	16	3		
2	C	1	Total	C	N	O	P	0	0
			39	15	5	16	3		
2	D	1	Total	C	N	O	P	0	0
			39	15	5	16	3		
2	E	1	Total	C	N	O	P	0	0
			39	15	5	16	3		
2	F	1	Total	C	N	O	P	0	0
			39	15	5	16	3		
2	G	1	Total	C	N	O	P	0	0
			39	15	5	16	3		
2	H	1	Total	C	N	O	P	0	0
			39	15	5	16	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	355	Total	O	0	0
			355	355		
3	B	356	Total	O	0	0
			356	356		
3	C	302	Total	O	0	0
			302	302		
3	D	341	Total	O	0	0
			341	341		

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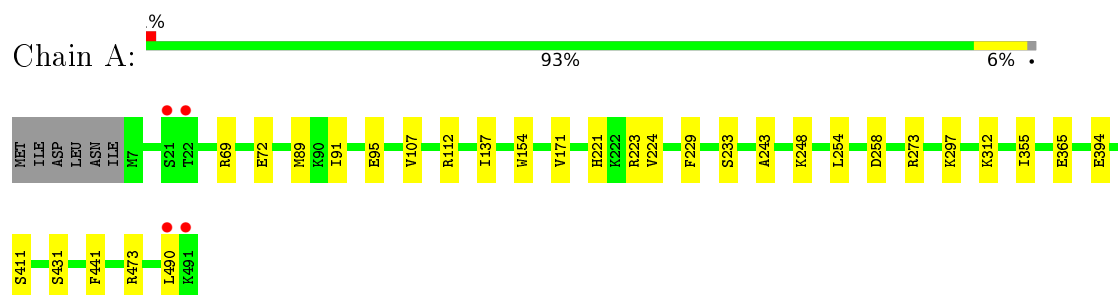
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	324	Total 324	O 324	0	0
3	F	294	Total 294	O 294	0	0
3	G	342	Total 342	O 342	0	0
3	H	325	Total 325	O 325	0	0

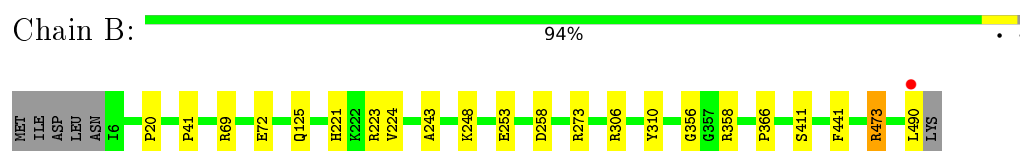
### 3 Residue-property plots [i](#)

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.

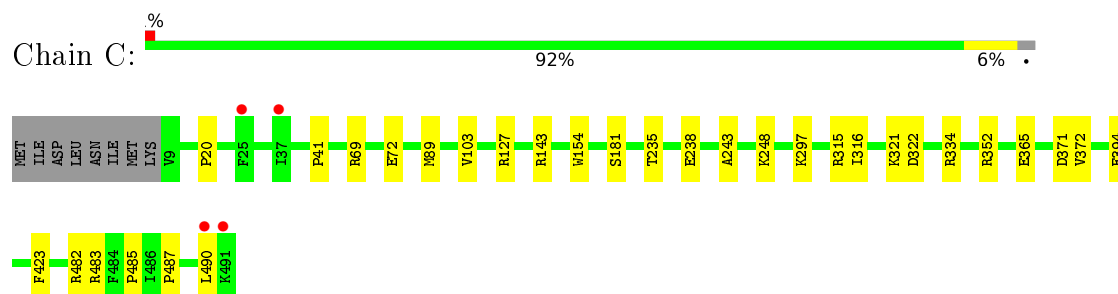
- Molecule 1: Aldehyde dehydrogenase



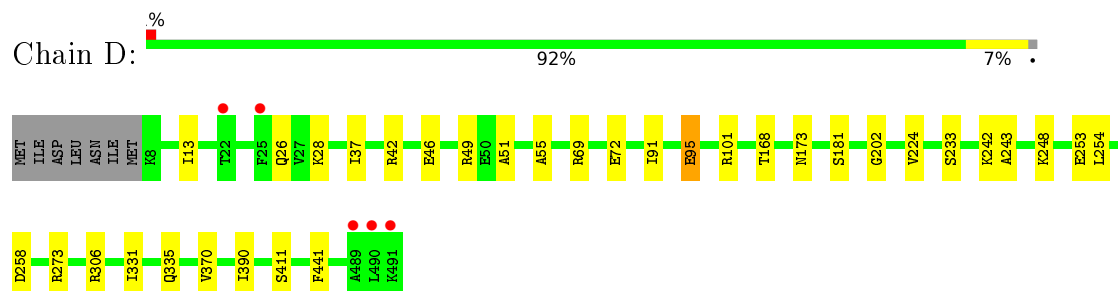
- Molecule 1: Aldehyde dehydrogenase



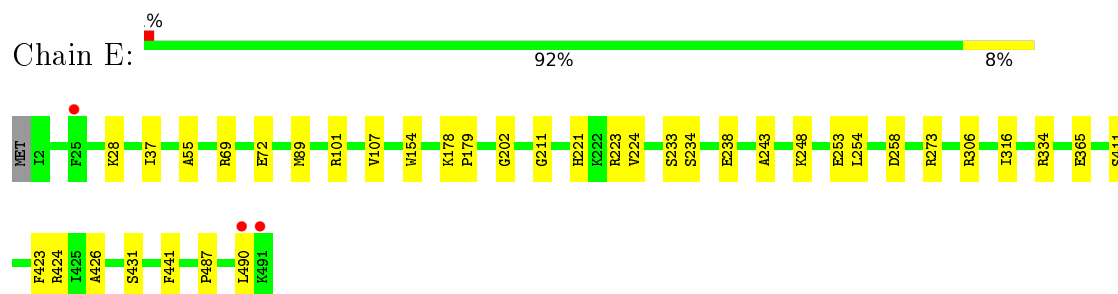
- Molecule 1: Aldehyde dehydrogenase



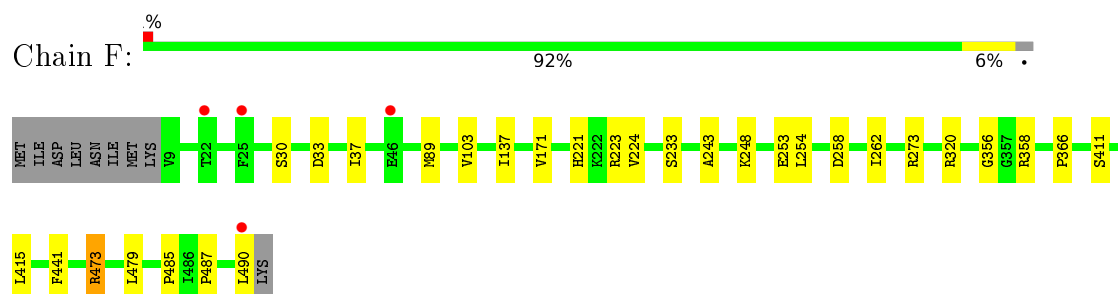
- Molecule 1: Aldehyde dehydrogenase



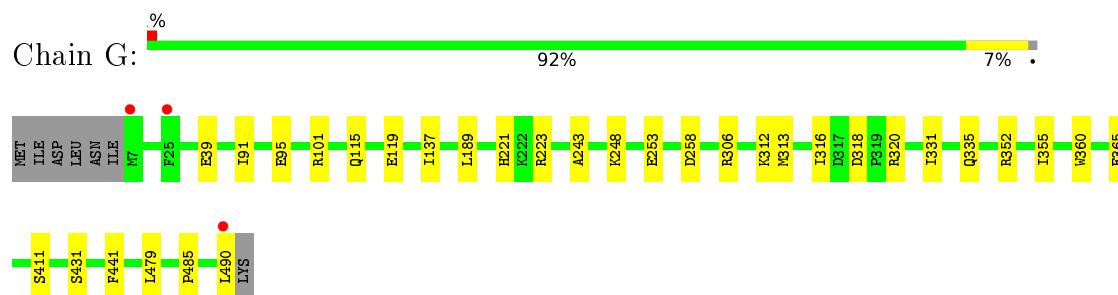
- Molecule 1: Aldehyde dehydrogenase



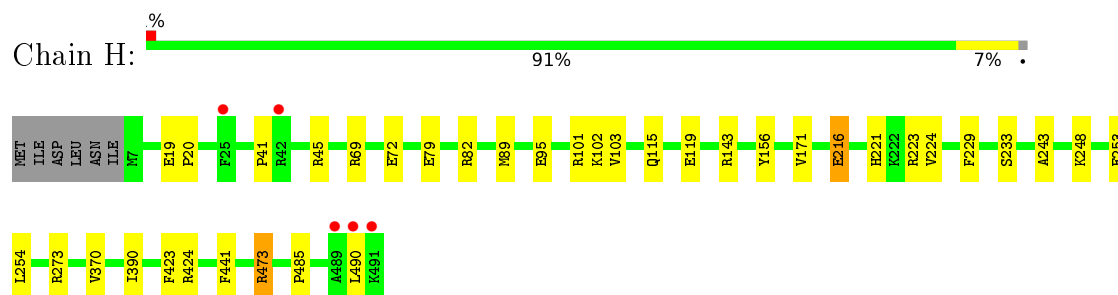
- Molecule 1: Aldehyde dehydrogenase



- Molecule 1: Aldehyde dehydrogenase



- Molecule 1: Aldehyde dehydrogenase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	164.55Å 185.21Å 208.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.65 – 1.90 49.65 – 1.89	Depositor EDS
% Data completeness (in resolution range)	98.5 (49.65-1.90) 98.6 (49.65-1.89)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.90 (at 1.90Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, $R_{free}$	0.193 , 0.221 0.193 , 0.220	Depositor DCC
$R_{free}$ test set	24720 reflections (5.02%)	DCC
Wilson B-factor (Å <sup>2</sup> )	19.7	Xtriage
Anisotropy	0.007	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 45.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	33123	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	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 44.92 % 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 1.4240e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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: NAP

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	0.39	0/3850	0.51	1/5225 (0.0%)
1	B	0.40	0/3865	0.53	0/5240
1	C	0.36	0/3851	0.48	0/5224
1	D	0.39	0/3851	0.51	0/5228
1	E	0.38	0/3905	0.51	0/5295
1	F	0.37	0/3851	0.52	1/5222 (0.0%)
1	G	0.40	0/3871	0.51	0/5252
1	H	0.37	0/3877	0.50	0/5256
All	All	0.38	0/30921	0.51	2/41942 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	473	ARG	NE-CZ-NH2	-5.75	117.43	120.30
1	A	473	ARG	NE-CZ-NH1	-5.19	117.71	120.30

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	3761	0	3732	22	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3775	0	3770	16	0
1	C	3761	0	3745	24	0
1	D	3755	0	3713	20	0
1	E	3812	0	3804	25	0
1	F	3760	0	3755	21	0
1	G	3767	0	3752	21	0
1	H	3781	0	3779	27	0
2	A	39	0	17	3	0
2	B	39	0	17	1	0
2	C	39	0	17	5	0
2	D	39	0	17	2	0
2	E	39	0	17	3	0
2	F	39	0	17	1	0
2	G	39	0	17	1	0
2	H	39	0	17	1	0
3	A	355	0	0	1	0
3	B	356	0	0	1	0
3	C	302	0	0	2	0
3	D	341	0	0	1	0
3	E	324	0	0	0	0
3	F	294	0	0	0	0
3	G	342	0	0	2	0
3	H	325	0	0	2	0
All	All	33123	0	30186	155	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

The worst 5 of 155 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:473:ARG:HH11	1:H:473:ARG:HG3	1.45	0.81
1:B:473:ARG:HG3	1:B:473:ARG:HH11	1.49	0.77
1:C:334:ARG:HE	2:C:500:NAP:H2D	1.56	0.71
1:C:315:ARG:NH2	1:C:322:ASP:OD2	2.23	0.70
1:E:426:ALA:O	1:F:473:ARG:NH2	2.25	0.70

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	487/491 (99%)	474 (97%)	13 (3%)	0	100	100
1	B	486/491 (99%)	475 (98%)	11 (2%)	0	100	100
1	C	485/491 (99%)	474 (98%)	11 (2%)	0	100	100
1	D	487/491 (99%)	474 (97%)	13 (3%)	0	100	100
1	E	492/491 (100%)	479 (97%)	13 (3%)	0	100	100
1	F	484/491 (99%)	471 (97%)	13 (3%)	0	100	100
1	G	489/491 (100%)	477 (98%)	12 (2%)	0	100	100
1	H	488/491 (99%)	475 (97%)	13 (3%)	0	100	100
All	All	3898/3928 (99%)	3799 (98%)	99 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	394/408 (97%)	392 (100%)	2 (0%)	92	92
1	B	398/408 (98%)	395 (99%)	3 (1%)	86	86
1	C	394/408 (97%)	393 (100%)	1 (0%)	94	95
1	D	390/408 (96%)	387 (99%)	3 (1%)	86	86
1	E	402/408 (98%)	399 (99%)	3 (1%)	88	88
1	F	397/408 (97%)	393 (99%)	4 (1%)	82	81

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	395/408 (97%)	392 (99%)	3 (1%)	86	86
1	H	398/408 (98%)	392 (98%)	6 (2%)	72	69
All	All	3168/3264 (97%)	3143 (99%)	25 (1%)	86	86

5 of 25 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	441	PHE
1	F	253[B]	GLU
1	H	441	PHE
1	F	171	VAL
1	F	358	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

8 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
2	NAP	A	500	-	36,42,52	1.89	10 (27%)	45,65,80	2.04	6 (13%)
2	NAP	B	500	-	36,42,52	1.95	11 (30%)	45,65,80	2.01	6 (13%)
2	NAP	C	500	-	36,42,52	1.91	11 (30%)	45,65,80	2.15	5 (11%)
2	NAP	D	500	-	36,42,52	1.92	11 (30%)	45,65,80	2.12	4 (8%)
2	NAP	E	500	-	36,42,52	1.94	12 (33%)	45,65,80	2.31	4 (8%)
2	NAP	F	500	-	36,42,52	1.91	10 (27%)	45,65,80	2.06	5 (11%)
2	NAP	G	500	-	36,42,52	1.93	11 (30%)	45,65,80	2.14	5 (11%)
2	NAP	H	500	-	36,42,52	1.90	12 (33%)	45,65,80	2.04	5 (11%)

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	NAP	A	500	-	-	0/23/56/67	0/4/4/5
2	NAP	B	500	-	-	0/23/56/67	0/4/4/5
2	NAP	C	500	-	-	0/23/56/67	0/4/4/5
2	NAP	D	500	-	-	0/23/56/67	0/4/4/5
2	NAP	E	500	-	-	0/23/56/67	0/4/4/5
2	NAP	F	500	-	-	0/23/56/67	0/4/4/5
2	NAP	G	500	-	-	0/23/56/67	0/4/4/5
2	NAP	H	500	-	-	0/23/56/67	0/4/4/5

The worst 5 of 88 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	NAP	C3B-C2B	-5.12	1.41	1.53
2	C	500	NAP	C3B-C2B	-5.11	1.41	1.53
2	G	500	NAP	C3B-C2B	-5.09	1.41	1.53
2	D	500	NAP	C3B-C2B	-5.07	1.41	1.53
2	E	500	NAP	C3B-C2B	-5.05	1.41	1.53

The worst 5 of 40 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	500	NAP	C1B-N9A-C4A	-10.88	114.66	126.81
2	C	500	NAP	C1B-N9A-C4A	-10.28	115.33	126.81
2	D	500	NAP	C1B-N9A-C4A	-9.67	116.02	126.81
2	G	500	NAP	C1B-N9A-C4A	-9.66	116.02	126.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	500	NAP	C1B-N9A-C4A	-9.64	116.05	126.81

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	NAP	3	0
2	B	500	NAP	1	0
2	C	500	NAP	5	0
2	D	500	NAP	2	0
2	E	500	NAP	3	0
2	F	500	NAP	1	0
2	G	500	NAP	1	0
2	H	500	NAP	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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	485/491 (98%)	-0.29	4 (0%) 87 88	9, 19, 35, 53	1 (0%)
1	B	485/491 (98%)	-0.44	1 (0%) 95 95	10, 17, 32, 44	1 (0%)
1	C	483/491 (98%)	-0.12	4 (0%) 87 88	10, 22, 40, 63	1 (0%)
1	D	484/491 (98%)	-0.15	5 (1%) 84 86	10, 19, 33, 65	1 (0%)
1	E	490/491 (99%)	-0.28	3 (0%) 90 91	9, 19, 35, 51	9 (1%)
1	F	482/491 (98%)	-0.27	4 (0%) 87 88	10, 21, 38, 55	1 (0%)
1	G	484/491 (98%)	-0.25	3 (0%) 90 91	9, 18, 35, 51	1 (0%)
1	H	485/491 (98%)	-0.25	5 (1%) 84 86	9, 20, 33, 63	1 (0%)
All	All	3878/3928 (98%)	-0.26	29 (0%) 89 90	9, 19, 36, 65	16 (0%)

The worst 5 of 29 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	491	LYS	4.8
1	H	491	LYS	4.5
1	D	490	LEU	4.5
1	D	491	LYS	4.2
1	E	491	LYS	4.0

### 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	NAP	F	500	39/48	0.92	0.14	2.14	21,28,42,43	39
2	NAP	A	500	39/48	0.92	0.16	1.88	18,24,39,41	39
2	NAP	H	500	39/48	0.93	0.14	1.83	17,24,42,45	39
2	NAP	B	500	39/48	0.96	0.11	1.74	15,20,38,43	39
2	NAP	D	500	39/48	0.93	0.13	1.70	19,26,41,45	39
2	NAP	C	500	39/48	0.92	0.15	1.64	23,29,43,44	39
2	NAP	E	500	39/48	0.94	0.13	1.31	16,22,39,42	39
2	NAP	G	500	39/48	0.94	0.12	1.20	18,22,36,42	39

## 6.5 Other polymers [i](#)

There are no such residues in this entry.