



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2017 – 01:42 AM EDT

PDB ID : 5MSR
Title : Structure of the unmodified PCP-R domain of carboxylic acid reductase (CAR) from *Segniliparus rugosus* in complex with NADPH, P43 form
Authors : Gahloth, D.; Leys, D.
Deposited on : unknown
Resolution : 2.37 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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-20029824

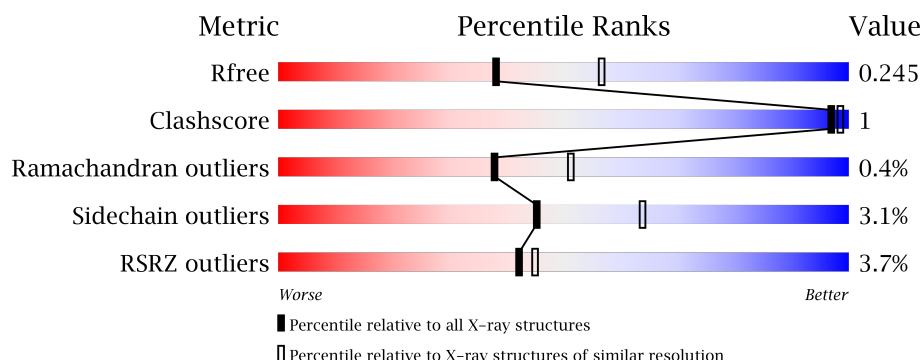
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.37 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}	100719	4388 (2.40-2.36)
Clashscore	112137	4984 (2.40-2.36)
Ramachandran outliers	110173	4907 (2.40-2.36)
Sidechain outliers	110143	4909 (2.40-2.36)
RSRZ outliers	101464	4423 (2.40-2.36)

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 $\leq 5\%$. 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	1188	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -5px; left: 0; width: 100%; height: 10px; background-color: red;"></div> <div style="position: absolute; top: 5px; left: 0; width: 100%; height: 10px; background-color: green;"></div> <div style="position: absolute; top: 5px; left: 41%; width: 10%; height: 10px; background-color: yellow;"></div> <div style="position: absolute; top: 5px; left: 57%; width: 41%; height: 10px; background-color: grey;"></div> </div> <div> % 41% 57% </div> </div>
1	B	1188	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -5px; left: 0; width: 100%; height: 10px; background-color: red;"></div> <div style="position: absolute; top: 5px; left: 0; width: 100%; height: 10px; background-color: green;"></div> <div style="position: absolute; top: 5px; left: 42%; width: 10%; height: 10px; background-color: yellow;"></div> <div style="position: absolute; top: 5px; left: 56%; width: 42%; height: 10px; background-color: grey;"></div> </div> <div> % 42% 56% </div> </div>
1	C	1188	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -5px; left: 0; width: 100%; height: 10px; background-color: red;"></div> <div style="position: absolute; top: 5px; left: 0; width: 100%; height: 10px; background-color: green;"></div> <div style="position: absolute; top: 5px; left: 2%; width: 10%; height: 10px; background-color: yellow;"></div> <div style="position: absolute; top: 5px; left: 57%; width: 41%; height: 10px; background-color: grey;"></div> </div> <div> 2% 41% 57% </div> </div>
1	D	1188	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -5px; left: 0; width: 100%; height: 10px; background-color: red;"></div> <div style="position: absolute; top: 5px; left: 0; width: 100%; height: 10px; background-color: green;"></div> <div style="position: absolute; top: 5px; left: 3%; width: 10%; height: 10px; background-color: yellow;"></div> <div style="position: absolute; top: 5px; left: 58%; width: 40%; height: 10px; background-color: grey;"></div> </div> <div> 3% 40% 58% </div> </div>

2 Entry composition ⓘ

There are 3 unique types of molecules in this entry. The entry contains 16105 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 Thioester reductase domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	514	Total 3926	2492	687	746	1	0	0	0
1	B	520	Total 3973	2519	693	760	1	0	0	0
1	C	516	Total 3944	2504	689	750	1	0	0	0
1	D	504	Total 3869	2461	676	731	1	0	0	0

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O P		
2	A	1	Total 40	15	6	16 3	0	0
2	B	1	Total 40	15	6	16 3	0	0

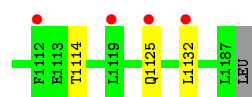
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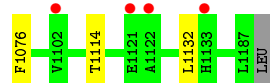
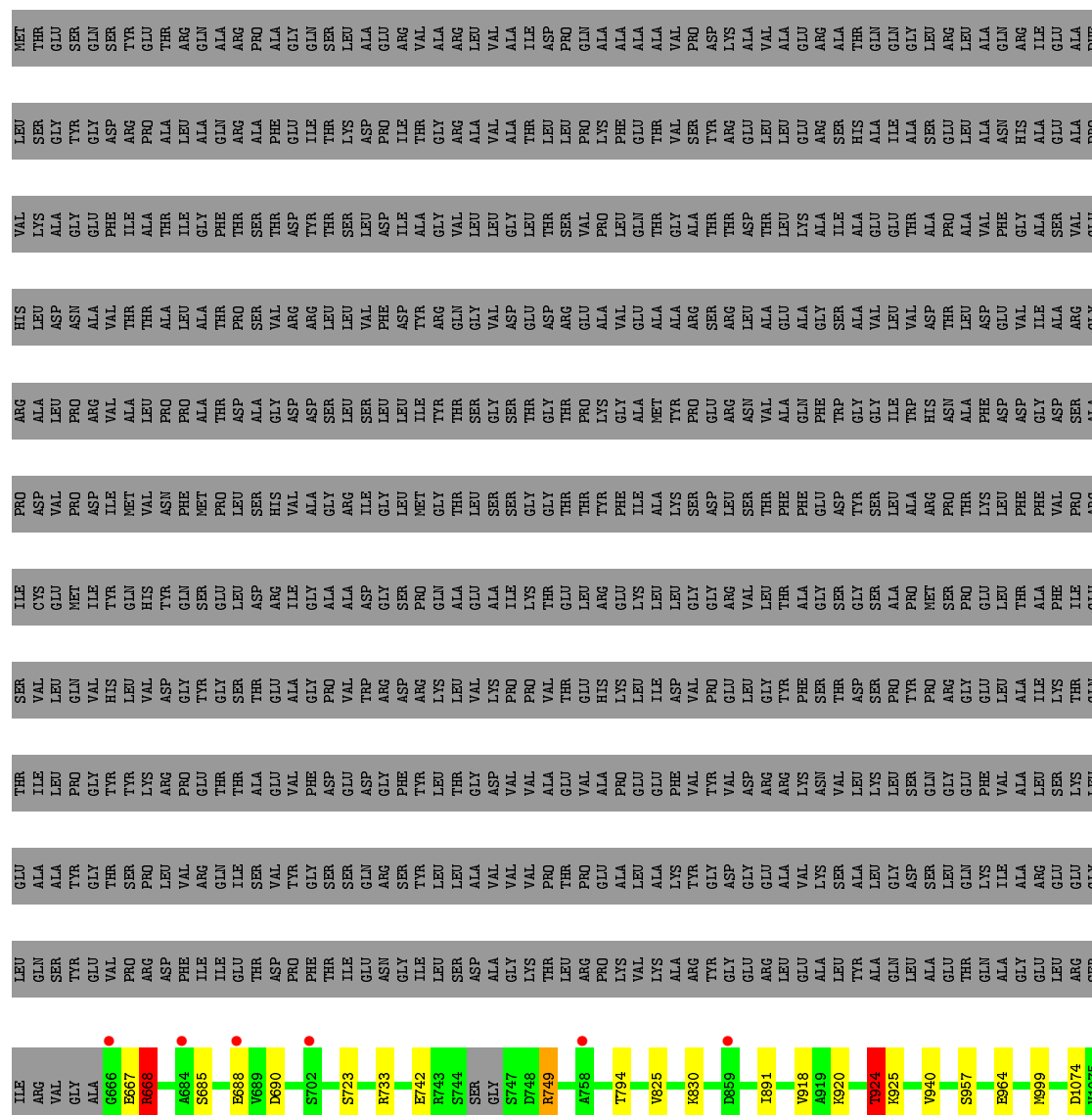
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	0	0
			40	15	6	16	3		
2	D	1	Total	C	N	O	P	0	0
			40	15	6	16	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	63	Total	O	0	0
			63	63		
3	B	75	Total	O	0	0
			75	75		
3	C	58	Total	O	0	0
			58	58		
3	D	37	Total	O	0	0
			37	37		

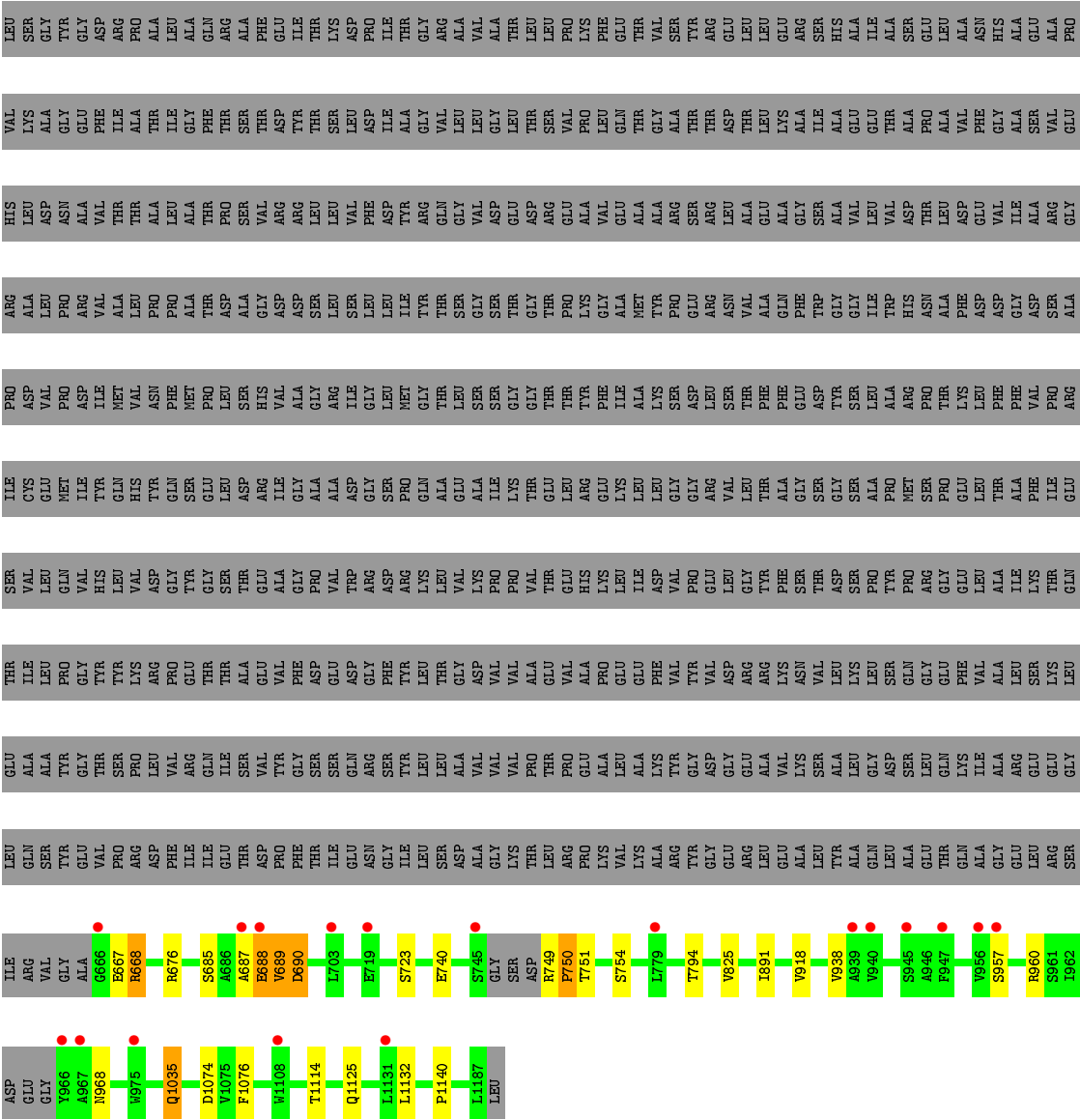


- Molecule 1: Thioester reductase domain-containing protein



- Molecule 1: Thioester reductase domain-containing protein







4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	94.94Å 94.94Å 335.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	200.00 – 2.37 91.36 – 2.37	Depositor EDS
% Data completeness (in resolution range)	98.7 (200.00-2.37) 98.7 (91.36-2.37)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.33 (at 2.37Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.219 , 0.241 0.229 , 0.245	Depositor DCC
R_{free} test set	5667 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	50.2	Xtriage
Anisotropy	0.011	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 40.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.057 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16105	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.57% 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: 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	0.41	0/4011	0.65	2/5457 (0.0%)
1	B	0.41	0/4060	0.69	5/5525 (0.1%)
1	C	0.41	0/4030	0.67	3/5483 (0.1%)
1	D	0.42	0/3953	0.69	10/5376 (0.2%)
All	All	0.41	0/16054	0.67	20/21841 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	733	ARG	NE-CZ-NH1	8.04	124.32	120.30
1	B	733	ARG	NE-CZ-NH2	-7.19	116.70	120.30
1	D	793	ARG	NE-CZ-NH1	7.10	123.85	120.30
1	D	764	ARG	NE-CZ-NH1	6.84	123.72	120.30
1	D	1101	ARG	NE-CZ-NH1	6.78	123.69	120.30
1	D	1103	ASP	CB-CG-OD1	6.60	124.24	118.30
1	A	1103	ASP	CB-CG-OD1	6.54	124.19	118.30
1	D	785	LEU	CA-CB-CG	6.50	130.24	115.30
1	C	688	GLU	C-N-CA	6.18	137.16	121.70
1	D	793	ARG	NE-CZ-NH2	-6.18	117.21	120.30
1	D	774	LEU	CA-CB-CG	6.16	129.48	115.30
1	C	750	PRO	N-CA-C	5.80	127.17	112.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	1149	THR	CB-CA-C	-5.47	96.82	111.60
1	B	940	VAL	N-CA-C	5.36	125.46	111.00
1	C	676	ARG	NE-CZ-NH1	5.35	122.98	120.30
1	B	733	ARG	CG-CD-NE	5.26	122.85	111.80
1	D	982	ARG	NE-CZ-NH2	-5.13	117.73	120.30
1	A	1103	ASP	CB-CG-OD2	-5.12	113.69	118.30
1	B	957	SER	N-CA-C	-5.08	97.27	111.00
1	D	764	ARG	NE-CZ-NH2	-5.05	117.77	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	924	THR	Peptide

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	3926	0	3880	5	0
1	B	3973	0	3912	4	1
1	C	3944	0	3894	9	1
1	D	3869	0	3825	15	0
2	A	40	0	19	0	0
2	B	40	0	19	0	0
2	C	40	0	19	0	0
2	D	40	0	19	0	0
3	A	63	0	0	0	0
3	B	75	0	0	0	0
3	C	58	0	0	0	0
3	D	37	0	0	0	0
All	All	16105	0	15587	33	1

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 1.

All (33) 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:782:ALA:HA	1:D:785:LEU:HD12	1.57	0.86
1:C:750:PRO:O	1:C:754:SER:OG	2.05	0.75
1:D:785:LEU:HD23	1:D:926:ARG:HA	1.67	0.74
1:A:667:GLU:HB2	1:A:668:ARG:HB2	1.75	0.69
1:D:1076:PHE:CE1	1:D:1149:THR:OG1	2.46	0.68
1:C:667:GLU:HB2	1:C:668:ARG:HB2	1.75	0.67
1:B:667:GLU:HB2	1:B:668:ARG:HB2	1.76	0.67
1:D:1149:THR:HG22	1:D:1149:THR:O	2.00	0.60
1:D:785:LEU:HD23	1:D:926:ARG:CA	2.33	0.59
1:C:687:ALA:HA	1:C:688:GLU:HG2	1.86	0.57
1:D:1076:PHE:CZ	1:D:1147:PHE:O	2.59	0.56
1:C:938:VAL:O	1:C:960:ARG:NH2	2.42	0.52
1:D:1076:PHE:CZ	1:D:1149:THR:OG1	2.61	0.52
1:D:782:ALA:HA	1:D:785:LEU:CD1	2.37	0.51
1:A:938:VAL:O	1:A:960:ARG:NH2	2.43	0.51
1:D:1149:THR:O	1:D:1151:ASN:N	2.42	0.51
1:D:938:VAL:O	1:D:960:ARG:NH2	2.43	0.51
1:D:1149:THR:CG2	1:D:1149:THR:O	2.59	0.51
1:C:688:GLU:HB3	1:C:690:ASP:N	2.26	0.50
1:C:750:PRO:O	1:C:751:THR:HG22	2.11	0.50
1:D:1149:THR:HG22	1:D:1153:GLN:H	1.80	0.47
1:A:760:ALA:O	1:A:962:ILE:HD11	2.16	0.45
1:A:1074:ASP:HB3	1:A:1076:PHE:CE2	2.52	0.44
1:C:1074:ASP:HB3	1:C:1076:PHE:CE2	2.53	0.43
1:B:1074:ASP:HB3	1:B:1076:PHE:CE2	2.53	0.43
1:D:1149:THR:HG21	1:D:1153:GLN:HB2	1.99	0.43
1:B:891:ILE:HG21	1:B:918:VAL:HG13	2.01	0.42
1:D:891:ILE:HG21	1:D:918:VAL:HG13	2.00	0.42
1:C:1035:GLN:HG2	1:C:1140:PRO:HA	2.02	0.42
1:C:891:ILE:HG21	1:C:918:VAL:HG13	2.01	0.42
1:A:891:ILE:HG21	1:A:918:VAL:HG13	2.01	0.41
1:B:920:LYS:O	1:B:924:THR:CG2	2.69	0.41
1:D:1076:PHE:HZ	1:D:1147:PHE:O	2.00	0.41

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:925:LYS:NZ	1:C:740:GLU:OE1[3_664]	2.14	0.06

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	508/1188 (43%)	491 (97%)	15 (3%)	2 (0%)	38	51
1	B	516/1188 (43%)	499 (97%)	15 (3%)	2 (0%)	38	51
1	C	510/1188 (43%)	490 (96%)	18 (4%)	2 (0%)	38	51
1	D	494/1188 (42%)	477 (97%)	14 (3%)	3 (1%)	28	39
All	All	2028/4752 (43%)	1957 (96%)	62 (3%)	9 (0%)	38	51

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	668	ARG
1	B	668	ARG
1	B	749	ARG
1	C	668	ARG
1	D	690	ASP
1	D	749	ARG
1	D	1150	LYS
1	A	689	VAL
1	C	689	VAL

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	409/947 (43%)	399 (98%)	10 (2%)	54	72
1	B	414/947 (44%)	399 (96%)	15 (4%)	40	58

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	411/947 (43%)	398 (97%)	13 (3%)	44	62
1	D	405/947 (43%)	392 (97%)	13 (3%)	44	62
All	All	1639/3788 (43%)	1588 (97%)	51 (3%)	45	64

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	685	SER
1	A	688	GLU
1	A	690	ASP
1	A	723	SER
1	A	749	ARG
1	A	825	VAL
1	A	957	SER
1	A	1114	THR
1	A	1125	GLN
1	A	1132	LEU
1	B	668	ARG
1	B	685	SER
1	B	688	GLU
1	B	690	ASP
1	B	723	SER
1	B	742	GLU
1	B	749	ARG
1	B	794	THR
1	B	825	VAL
1	B	830	LYS
1	B	924	THR
1	B	964	GLU
1	B	999	MET
1	B	1114	THR
1	B	1132	LEU
1	C	685	SER
1	C	689	VAL
1	C	690	ASP
1	C	723	SER
1	C	749	ARG
1	C	794	THR
1	C	825	VAL
1	C	957	SER
1	C	968	ASN

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Mol	Chain	Res	Type
1	C	1035	GLN
1	C	1114	THR
1	C	1125	GLN
1	C	1132	LEU
1	D	723	SER
1	D	742	GLU
1	D	749	ARG
1	D	774	LEU
1	D	785	LEU
1	D	793	ARG
1	D	825	VAL
1	D	957	SER
1	D	999	MET
1	D	1057	GLU
1	D	1114	THR
1	D	1125	GLN
1	D	1132	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1126	HIS
1	B	675	GLN
1	B	1126	HIS
1	C	675	GLN
1	C	1011	ASN
1	C	1157	GLN
1	D	675	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

4 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	1201	-	36,43,52	0.91	2 (5%)	42,67,80	1.47	3 (7%)
2	NAP	B	1201	-	36,43,52	0.90	2 (5%)	42,67,80	1.49	2 (4%)
2	NAP	C	1201	-	36,43,52	0.94	2 (5%)	42,67,80	1.50	2 (4%)
2	NAP	D	1201	-	36,43,52	0.89	1 (2%)	42,67,80	1.60	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	1201	-	-	0/23/59/67	0/4/4/5
2	NAP	B	1201	-	-	0/23/59/67	0/4/4/5
2	NAP	C	1201	-	-	0/23/59/67	0/4/4/5
2	NAP	D	1201	-	-	0/23/59/67	0/4/4/5

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1201	NAP	C2A-N3A	2.05	1.35	1.32
2	C	1201	NAP	C2A-N3A	2.14	1.35	1.32
2	B	1201	NAP	C2A-N3A	2.20	1.35	1.32
2	B	1201	NAP	C5A-C4A	2.93	1.47	1.40
2	D	1201	NAP	C5A-C4A	2.97	1.47	1.40
2	A	1201	NAP	C5A-C4A	3.02	1.47	1.40
2	C	1201	NAP	C5A-C4A	3.10	1.47	1.40

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1201	NAP	N3A-C2A-N1A	-7.64	122.21	128.86
2	B	1201	NAP	N3A-C2A-N1A	-7.37	122.44	128.86
2	C	1201	NAP	N3A-C2A-N1A	-7.19	122.59	128.86
2	A	1201	NAP	N3A-C2A-N1A	-6.91	122.84	128.86
2	D	1201	NAP	C4A-C5A-N7A	-2.79	106.71	109.41
2	A	1201	NAP	C4A-C5A-N7A	-2.70	106.81	109.41
2	B	1201	NAP	C4A-C5A-N7A	-2.68	106.82	109.41
2	C	1201	NAP	C4A-C5A-N7A	-2.49	107.00	109.41
2	D	1201	NAP	C1B-N9A-C4A	-2.29	122.67	126.64
2	A	1201	NAP	O4B-C1B-C2B	-2.00	103.09	106.59
2	D	1201	NAP	C2A-N1A-C6A	2.08	122.41	118.77
2	D	1201	NAP	O3X-P2B-O2X	2.30	116.87	107.61

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

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	514/1188 (43%)	0.36	11 (2%) 64 65	44, 79, 126, 182	0
1	B	520/1188 (43%)	0.34	10 (1%) 67 68	42, 75, 126, 187	0
1	C	516/1188 (43%)	0.49	18 (3%) 44 47	41, 84, 141, 203	0
1	D	504/1188 (42%)	0.78	38 (7%) 15 16	46, 93, 153, 185	0
All	All	2054/4752 (43%)	0.49	77 (3%) 42 45	41, 82, 142, 203	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	966	TYR	7.3
1	D	945	SER	6.9
1	D	1146	PRO	6.7
1	C	687	ALA	6.7
1	D	966	TYR	6.5
1	B	684	ALA	6.3
1	D	956	VAL	5.0
1	D	947	PHE	5.0
1	D	1135	PHE	4.7
1	D	1142	ILE	4.5
1	D	1149	THR	4.5
1	D	1133	HIS	4.4
1	B	666	GLY	4.3
1	D	1076	PHE	4.1
1	C	956	VAL	4.0
1	D	946	ALA	4.0
1	D	944	PRO	3.9
1	D	1148	GLN	3.8
1	D	684	ALA	3.6
1	D	967	ALA	3.2
1	A	703	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	943	GLU	3.1
1	D	1121	GLU	3.0
1	C	967	ALA	3.0
1	D	1105	TYR	2.9
1	D	1147	PHE	2.9
1	D	761	THR	2.9
1	A	1125	GLN	2.9
1	A	956	VAL	2.8
1	C	719	GLU	2.8
1	D	940	VAL	2.7
1	A	1112	PHE	2.6
1	D	1131	LEU	2.6
1	C	939	ALA	2.6
1	D	1112	PHE	2.6
1	D	1034	TYR	2.6
1	B	1122	ALA	2.6
1	B	758	ALA	2.5
1	C	666	GLY	2.5
1	B	702	SER	2.4
1	D	670	VAL	2.4
1	D	942	VAL	2.4
1	C	1131	LEU	2.4
1	B	1121	GLU	2.4
1	D	739	ILE	2.4
1	D	774	LEU	2.4
1	C	745	SER	2.3
1	A	764	ARG	2.3
1	C	945	SER	2.3
1	D	939	ALA	2.3
1	A	1038	ALA	2.3
1	B	1102	VAL	2.2
1	A	1132	LEU	2.2
1	D	957	SER	2.2
1	A	942	VAL	2.2
1	D	1144	GLY	2.2
1	A	962	ILE	2.2
1	D	1064	LEU	2.2
1	D	1108	TRP	2.2
1	C	947	PHE	2.1
1	C	703	LEU	2.1
1	C	940	VAL	2.1
1	B	859	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	688	GLU	2.1
1	A	943	GLU	2.1
1	B	688	GLU	2.1
1	B	1133	HIS	2.1
1	D	747	SER	2.1
1	A	1119	LEU	2.1
1	C	779	LEU	2.1
1	D	785	LEU	2.1
1	D	736	ALA	2.0
1	D	886	ASP	2.0
1	C	975	TRP	2.0
1	C	1108	TRP	2.0
1	D	953	ILE	2.0
1	C	957	SER	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, 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	LLDF	B-factors(\AA^2)	Q<0.9
2	NAP	C	1201	40/48	0.98	0.12	-0.85	37,44,65,67	0
2	NAP	D	1201	40/48	0.97	0.12	-1.02	37,42,66,68	0
2	NAP	A	1201	40/48	0.98	0.11	-1.15	37,43,60,60	0
2	NAP	B	1201	40/48	0.98	0.10	-1.20	34,43,61,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.