



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:48 AM GMT

PDB ID : 3Q12
Title : Pantoate-beta-alanine ligase from Yersinia pestis in complex with pantoate.
Authors : Osipiuk, J.; Maltseva, N.; Kwon, K.; Anderson, W.F.; Joachimiak, A.; Center
for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2010-12-16
Resolution : 1.58 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 (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

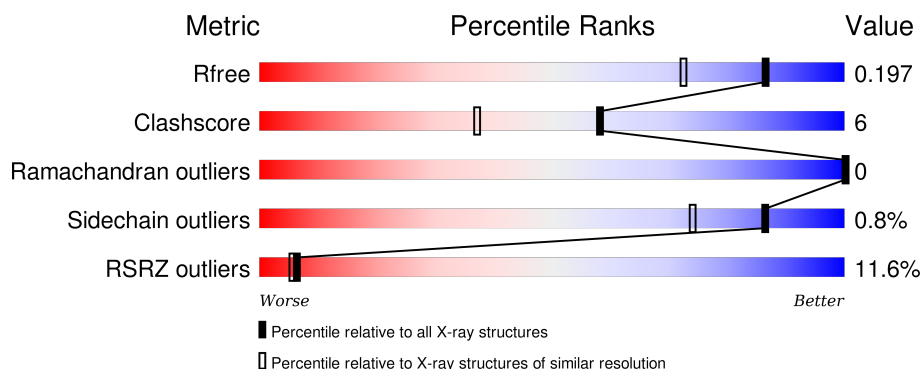
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.58 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	3815 (1.60-1.56)
Clashscore	102246	4131 (1.60-1.56)
Ramachandran outliers	100387	4021 (1.60-1.56)
Sidechain outliers	100360	4018 (1.60-1.56)
RSRZ outliers	91569	3823 (1.60-1.56)

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	287	<div> <div>10%</div> <div>91%</div> <div>8%</div> <div>.</div> </div>
1	B	287	<div> <div>10%</div> <div>92%</div> <div>7%</div> <div>.</div> </div>
1	C	287	<div> <div>2%</div> <div>93%</div> <div>6%</div> <div>.</div> </div>
1	D	287	<div> <div>33%</div> <div>82%</div> <div>16%</div> <div>..</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10600 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.

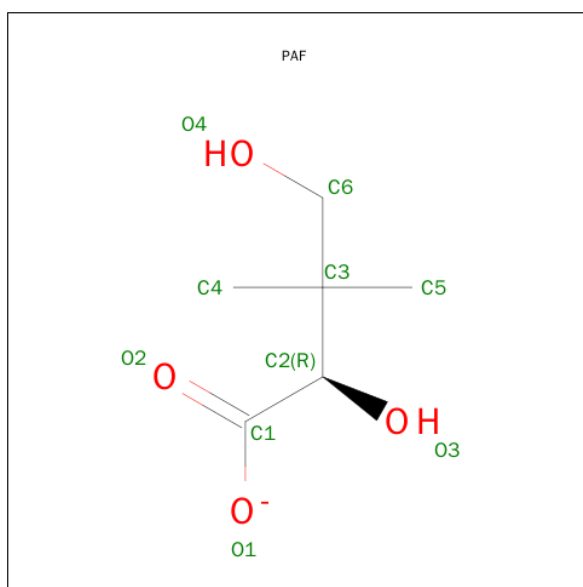
- Molecule 1 is a protein called Pantoate--beta-alanine ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	0	19	0
			2351	1501	405	435	10			
1	B	284	Total	C	N	O	S	0	25	0
			2374	1521	402	441	10			
1	C	285	Total	C	N	O	S	0	22	0
			2370	1512	412	436	10			
1	D	283	Total	C	N	O	S	0	15	0
			2305	1475	395	425	10			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP Q8ZBK7
A	-1	ASN	-	EXPRESSION TAG	UNP Q8ZBK7
A	0	ALA	-	EXPRESSION TAG	UNP Q8ZBK7
B	-2	SER	-	EXPRESSION TAG	UNP Q8ZBK7
B	-1	ASN	-	EXPRESSION TAG	UNP Q8ZBK7
B	0	ALA	-	EXPRESSION TAG	UNP Q8ZBK7
C	-2	SER	-	EXPRESSION TAG	UNP Q8ZBK7
C	-1	ASN	-	EXPRESSION TAG	UNP Q8ZBK7
C	0	ALA	-	EXPRESSION TAG	UNP Q8ZBK7
D	-2	SER	-	EXPRESSION TAG	UNP Q8ZBK7
D	-1	ASN	-	EXPRESSION TAG	UNP Q8ZBK7
D	0	ALA	-	EXPRESSION TAG	UNP Q8ZBK7

- Molecule 2 is PANTOATE (three-letter code: PAF) (formula: C₆H₁₁O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			10	6	4		
2	B	1	Total	C	O	0	0
			10	6	4		
2	C	1	Total	C	O	0	0
			10	6	4		
2	D	1	Total	C	O	0	0
			10	6	4		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	C	1	Total	Cl	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	376	Total	O	0	4
			378	378		
4	B	299	Total	O	0	1
			299	299		
4	C	303	Total	O	0	5
			305	305		

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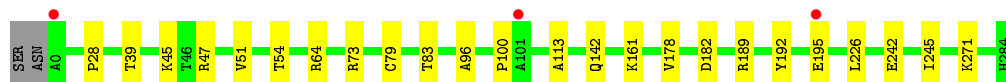
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	176	Total	O	0	2
			176	176		

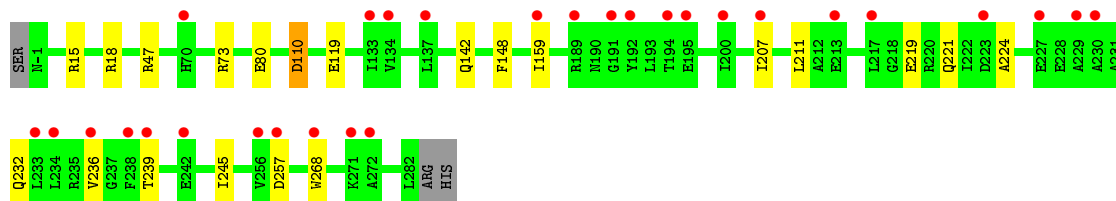
3 Residue-property plots

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: Pantoate--beta-alanine ligase



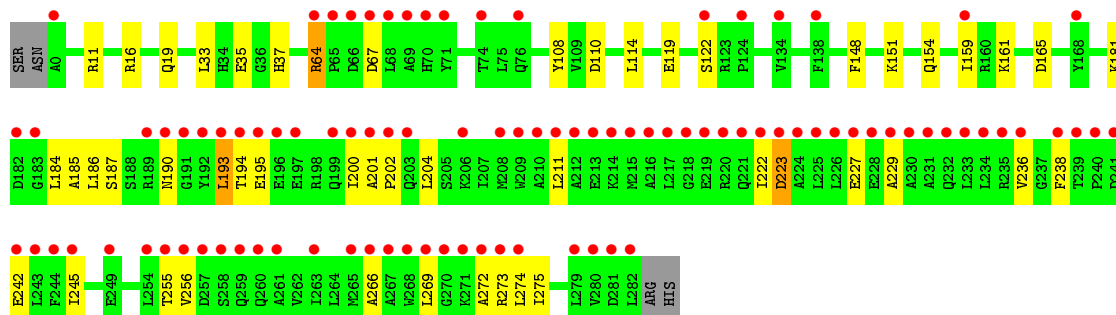
- Molecule 1: Pantoate--beta-alanine ligase



- Molecule 1: Pantoate--beta-alanine ligase



- Molecule 1: Pantoate--beta-alanine ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	44.73 Å 77.05 Å 98.76 Å 100.67° 95.33° 94.43°	Depositor
Resolution (Å)	36.69 – 1.58 36.69 – 1.58	Depositor EDS
% Data completeness (in resolution range)	87.2 (36.69-1.58) 84.1 (36.69-1.58)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 1.58 Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.150 , 0.195 0.153 , 0.197	Depositor DCC
R_{free} test set	7729 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	20.1	Xtriage
Anisotropy	0.700	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 48.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 153765 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10600	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.87% 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, 0.375 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: PAF, 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	0.94	0/2449	0.89	4/3315 (0.1%)
1	B	0.86	0/2490	0.88	6/3378 (0.2%)
1	C	0.83	0/2475	0.84	2/3350 (0.1%)
1	D	0.75	0/2390	0.77	1/3244 (0.0%)
All	All	0.85	0/9804	0.85	13/13287 (0.1%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	110[A]	ASP	CB-CG-OD2	-7.67	111.39	118.30
1	B	110[B]	ASP	CB-CG-OD2	-7.67	111.39	118.30
1	B	47	ARG	NE-CZ-NH2	-6.60	117.00	120.30
1	B	110[A]	ASP	CB-CG-OD1	6.42	124.08	118.30
1	B	110[B]	ASP	CB-CG-OD1	6.42	124.08	118.30
1	A	73	ARG	NE-CZ-NH2	-6.08	117.26	120.30
1	B	47	ARG	NE-CZ-NH1	5.56	123.08	120.30
1	C	66[A]	ASP	CB-CG-OD1	5.48	123.24	118.30
1	C	66[B]	ASP	CB-CG-OD1	5.48	123.24	118.30
1	D	165	ASP	CB-CG-OD2	-5.38	113.46	118.30
1	A	47	ARG	NE-CZ-NH1	5.17	122.89	120.30
1	A	64	ARG	NE-CZ-NH1	5.02	122.81	120.30
1	A	182	ASP	CB-CG-OD1	-5.02	113.78	118.30

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	2351	0	2437	23	0
1	B	2374	0	2469	15	0
1	C	2370	0	2475	21	0
1	D	2305	0	2391	55	0
2	A	10	0	11	1	0
2	B	10	0	11	0	0
2	C	10	0	11	0	0
2	D	10	0	11	0	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
4	A	378	0	0	4	0
4	B	299	0	0	9	0
4	C	305	0	0	3	0
4	D	176	0	0	3	0
All	All	10600	0	9816	113	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 6.

All (113) 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:204:LEU:CD2	1:D:274[B]:LEU:HD13	1.65	1.26
1:A:113:ALA:CB	1:A:161[B]:LYS:HE2	1.73	1.18
1:D:204:LEU:HD22	1:D:274[B]:LEU:CD1	1.74	1.16
1:A:113:ALA:HB2	1:A:161[B]:LYS:HE2	1.31	1.04
1:A:161[B]:LYS:HE3	4:B:417:HOH:O	1.63	0.97
1:C:219[B]:GLU:HG3	1:C:225:LEU:HD11	1.49	0.93
1:D:64[A]:ARG:NE	1:D:67:ASP:HB2	1.84	0.91
2:A:501:PAF:O1	4:A:867:HOH:O	1.89	0.90
1:C:113:ALA:HB2	1:C:161[B]:LYS:HE2	1.53	0.88
1:C:113:ALA:CB	1:C:161[B]:LYS:HE2	2.04	0.87
1:A:113:ALA:HB3	1:A:161[B]:LYS:HE2	1.54	0.87
1:D:64[A]:ARG:CZ	1:D:67:ASP:HB2	2.07	0.85
1:D:238:PHE:HE1	1:D:274[B]:LEU:HD21	1.40	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:76[B]:GLN:NE2	4:C:1059:HOH:O	2.10	0.84
1:C:219[B]:GLU:HG2	1:C:225:LEU:HD21	1.63	0.80
1:B:221[A]:GLN:NE2	4:B:382:HOH:O	2.21	0.72
1:A:113:ALA:HB3	1:A:161[B]:LYS:CE	2.19	0.72
1:D:204:LEU:HD22	1:D:274[B]:LEU:HD13	0.81	0.72
1:A:189[B]:ARG:HD3	4:A:590:HOH:O	1.91	0.70
1:D:242[B]:GLU:OE2	1:D:242[B]:GLU:HA	1.92	0.69
1:A:45[A]:LYS:HD3	1:A:51:VAL:HG23	1.73	0.69
1:D:35[B]:GLU:HG2	1:D:186:LEU:HD11	1.75	0.69
1:B:110[A]:ASP:OD2	4:B:333:HOH:O	2.10	0.67
1:B:221[B]:GLN:HG3	1:B:224:ALA:HB3	1.76	0.67
1:D:184:LEU:HD13	1:D:185:ALA:O	1.96	0.66
1:D:35[B]:GLU:HG2	1:D:186:LEU:CD1	2.25	0.66
1:D:200:ILE:HG23	1:D:269:LEU:HD11	1.78	0.66
1:D:16[B]:ARG:NH2	1:D:19:GLN:OE1	2.29	0.65
1:A:113:ALA:CB	1:A:161[B]:LYS:CE	2.62	0.62
1:D:186:LEU:HD23	1:D:186:LEU:C	2.21	0.61
1:B:80[B]:GLU:OE1	4:B:843:HOH:O	2.16	0.61
1:C:17:TRP:CZ3	1:C:50[B]:VAL:HG11	2.37	0.60
1:D:184:LEU:CD1	1:D:190:ASN:ND2	2.64	0.60
1:D:255:THR:HG22	1:D:256:VAL:N	2.17	0.59
1:B:18:ARG:NH2	4:B:606:HOH:O	2.36	0.59
1:C:113:ALA:HB3	1:C:161[B]:LYS:CD	2.33	0.59
1:A:113:ALA:HB3	1:A:161[B]:LYS:CD	2.34	0.58
1:C:113:ALA:HB3	1:C:161[B]:LYS:HE2	1.85	0.58
1:D:110[B]:ASP:OD2	4:D:538:HOH:O	2.17	0.57
1:A:161[B]:LYS:CE	4:B:417:HOH:O	2.36	0.57
1:A:79:CYS:O	1:A:83[B]:THR:HG23	2.05	0.57
1:D:238:PHE:CE1	1:D:274[B]:LEU:HD21	2.32	0.55
1:D:186:LEU:HD23	1:D:187:SER:N	2.21	0.54
1:D:194:THR:HG22	1:D:195:GLU:N	2.23	0.54
1:D:211:LEU:HD11	1:D:229:ALA:HB2	1.89	0.53
1:D:184:LEU:CD1	1:D:190:ASN:HD21	2.22	0.53
1:B:148:PHE:CE1	1:B:159[B]:ILE:HD13	2.43	0.53
1:B:211:LEU:HD21	1:B:245:ILE:HD11	1.91	0.52
1:D:148:PHE:CE1	1:D:159[B]:ILE:HD13	2.44	0.52
1:A:39[A]:THR:HG21	4:A:375:HOH:O	2.10	0.52
1:C:113:ALA:HB3	1:C:161[B]:LYS:HD3	1.91	0.51
1:C:35:GLU:O	1:C:39:THR:HG23	2.11	0.51
1:D:151:LYS:NZ	1:D:275:ILE:HD11	2.26	0.50
1:C:214:LYS:NZ	4:C:341:HOH:O	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:219[B]:GLU:CG	1:C:225:LEU:HD21	2.38	0.50
1:D:266:ALA:HB1	1:D:273:ARG:NH2	2.26	0.50
1:D:184:LEU:HD11	1:D:190:ASN:ND2	2.26	0.50
1:D:33:LEU:HD22	1:D:37:HIS:CG	2.48	0.49
1:A:192[B]:TYR:CE1	1:A:271:LYS:HD3	2.48	0.48
1:C:76[A]:GLN:NE2	4:C:935:HOH:O	2.37	0.48
1:D:255:THR:HG22	1:D:256:VAL:H	1.78	0.48
1:C:64[A]:ARG:NH1	1:C:66[A]:ASP:OD2	2.47	0.48
1:B:239:THR:O	1:B:268[A]:TRP:HE3	1.97	0.48
1:A:39[B]:THR:HG21	1:A:178:VAL:HG21	1.95	0.48
1:A:113:ALA:HB3	1:A:161[B]:LYS:HD3	1.97	0.47
1:B:142[B]:GLN:NE2	4:B:436:HOH:O	2.46	0.46
1:D:119[B]:GLU:HB2	1:D:154:GLN:OE1	2.15	0.46
1:C:113:ALA:HB3	1:C:161[B]:LYS:CE	2.46	0.46
1:D:64[A]:ARG:CZ	1:D:67:ASP:CB	2.89	0.46
1:D:184:LEU:CD1	1:D:185:ALA:O	2.65	0.45
1:D:200:ILE:CG2	1:D:269:LEU:HD11	2.46	0.45
1:D:238:PHE:CE1	1:D:274[B]:LEU:HD11	2.52	0.45
1:A:142:GLN:OE1	1:B:18:ARG:NH2	2.43	0.45
1:D:35[B]:GLU:CG	1:D:186:LEU:CD1	2.95	0.44
1:D:222:ILE:HD12	1:D:245:ILE:HB	1.98	0.44
1:A:195[B]:GLU:CD	1:A:195[B]:GLU:H	2.21	0.44
1:A:226:LEU:HD21	1:A:245:ILE:HG13	1.99	0.44
1:D:11:ARG:NH2	4:D:881:HOH:O	2.50	0.44
1:D:193:LEU:HG	1:D:269:LEU:CD2	2.48	0.43
1:A:96:ALA:O	1:A:100:PRO:HA	2.18	0.43
1:D:242[B]:GLU:OE2	1:D:242[B]:GLU:CA	2.62	0.43
1:D:222:ILE:HG23	1:D:245:ILE:HD12	2.00	0.43
1:D:114:LEU:HD12	1:D:161[B]:LYS:HD3	1.99	0.43
1:D:64[A]:ARG:CD	1:D:67:ASP:HB2	2.47	0.43
1:D:186:LEU:CD2	1:D:186:LEU:C	2.87	0.43
1:C:17:TRP:CE3	1:C:50[B]:VAL:HG11	2.54	0.43
1:D:255:THR:CG2	1:D:256:VAL:N	2.82	0.43
1:B:73[A]:ARG:NH1	4:B:388:HOH:O	2.52	0.43
1:C:110[A]:ASP:HB2	1:D:108:TYR:CZ	2.54	0.42
1:C:82:LEU:HD13	1:C:90:VAL:HG22	2.01	0.42
1:D:184:LEU:HD11	1:D:190:ASN:CG	2.40	0.42
1:D:184:LEU:HD12	1:D:190:ASN:ND2	2.34	0.42
1:D:200:ILE:HD11	1:D:236:VAL:HB	2.02	0.42
1:D:16[B]:ARG:HA	1:D:16[B]:ARG:CZ	2.50	0.42
1:D:122:SER:N	4:D:999:HOH:O	2.43	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:269:LEU:N	1:D:272:ALA:O	2.49	0.42
1:B:236:VAL:HG12	1:B:236:VAL:O	2.19	0.42
1:C:113:ALA:CB	1:C:161[B]:LYS:CE	2.87	0.41
1:A:45[A]:LYS:HD2	4:A:879:HOH:O	2.20	0.41
1:D:201:ALA:N	1:D:202:PRO:CD	2.84	0.41
1:B:15:ARG:NE	4:B:384:HOH:O	2.52	0.41
1:D:223:ASP:O	1:D:227:GLU:HG2	2.20	0.41
1:D:193:LEU:HD11	1:D:272:ALA:HB3	2.01	0.41
1:A:45[A]:LYS:HG2	1:A:51:VAL:HG21	2.03	0.41
1:D:194:THR:HG22	1:D:195:GLU:H	1.85	0.41
1:D:211:LEU:HD11	1:D:229:ALA:CB	2.51	0.41
1:B:207:ILE:HG12	1:B:232[B]:GLN:HG2	2.02	0.40
1:A:28:PRO:HA	1:A:54:THR:OG1	2.21	0.40
1:C:28:PRO:HA	1:C:54:THR:OG1	2.22	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	301/287 (105%)	297 (99%)	4 (1%)	0	100	100
1	B	307/287 (107%)	302 (98%)	5 (2%)	0	100	100
1	C	305/287 (106%)	300 (98%)	5 (2%)	0	100	100
1	D	296/287 (103%)	289 (98%)	7 (2%)	0	100	100
All	All	1209/1148 (105%)	1188 (98%)	21 (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	257/240 (107%)	255 (99%)	2 (1%)	86	74
1	B	262/240 (109%)	258 (98%)	4 (2%)	72	48
1	C	260/240 (108%)	258 (99%)	2 (1%)	86	74
1	D	251/240 (105%)	246 (98%)	5 (2%)	63	34
All	All	1030/960 (107%)	1017 (99%)	13 (1%)	86	55

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	242[A]	GLU
1	A	242[B]	GLU
1	B	119[A]	GLU
1	B	119[B]	GLU
1	B	219[A]	GLU
1	B	219[B]	GLU
1	C	66[A]	ASP
1	C	66[B]	ASP
1	D	64[A]	ARG
1	D	64[B]	ARG
1	D	181	LYS
1	D	193	LEU
1	D	223	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	HIS
1	C	58	ASN
1	C	61	GLN
1	D	34	HIS

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 ⓘ

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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$
2	PAF	A	501	-	5,9,9	1.38	1 (20%)	7,13,13	5.11	7 (100%)
2	PAF	B	501	-	5,9,9	2.80	2 (40%)	7,13,13	4.98	6 (85%)
2	PAF	C	501	-	5,9,9	2.41	2 (40%)	7,13,13	5.02	5 (71%)
2	PAF	D	501	-	5,9,9	2.80	3 (60%)	7,13,13	6.71	7 (100%)

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	PAF	A	501	-	-	0/9/13/13	0/0/0/0
2	PAF	B	501	-	-	0/9/13/13	0/0/0/0
2	PAF	C	501	-	-	0/9/13/13	0/0/0/0
2	PAF	D	501	-	-	0/9/13/13	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	PAF	C3-C2	-5.66	1.50	1.55
2	D	501	PAF	C3-C2	-5.03	1.50	1.55
2	C	501	PAF	C3-C2	-4.10	1.51	1.55
2	D	501	PAF	C5-C3	-2.82	1.47	1.53
2	D	501	PAF	C4-C3	-2.28	1.49	1.53
2	A	501	PAF	C3-C2	-2.25	1.53	1.55
2	B	501	PAF	C5-C3	-2.14	1.49	1.53
2	C	501	PAF	C4-C3	-2.14	1.49	1.53

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	PAF	C3-C2-C1	-12.37	104.22	115.58
2	B	501	PAF	C3-C2-C1	-9.07	107.25	115.58
2	A	501	PAF	C3-C2-C1	-7.95	108.28	115.58
2	C	501	PAF	C3-C2-C1	-7.57	108.63	115.58
2	D	501	PAF	C5-C3-C6	-6.59	95.70	108.84
2	C	501	PAF	C4-C3-C6	-6.45	95.98	108.84
2	D	501	PAF	C4-C3-C6	-6.24	96.39	108.84
2	A	501	PAF	C5-C3-C6	-6.11	96.66	108.84
2	B	501	PAF	C5-C3-C6	-4.41	100.04	108.84
2	B	501	PAF	C4-C3-C6	-3.81	101.25	108.84
2	A	501	PAF	C4-C3-C6	-3.60	101.66	108.84
2	C	501	PAF	C5-C3-C6	-3.58	101.70	108.84
2	A	501	PAF	O4-C6-C3	2.39	117.44	113.03
2	B	501	PAF	C5-C3-C4	2.73	114.76	109.28
2	D	501	PAF	O4-C6-C3	3.11	118.77	113.03
2	A	501	PAF	C5-C3-C4	3.24	115.79	109.28
2	B	501	PAF	C4-C3-C2	3.60	115.92	109.34
2	D	501	PAF	C4-C3-C2	4.29	117.17	109.34
2	D	501	PAF	C5-C3-C4	4.98	119.29	109.28
2	A	501	PAF	C4-C3-C2	5.00	118.47	109.34
2	C	501	PAF	C4-C3-C2	5.09	118.63	109.34
2	D	501	PAF	C5-C3-C2	5.18	118.80	109.34
2	A	501	PAF	C5-C3-C2	5.29	119.01	109.34
2	C	501	PAF	C5-C3-C4	5.76	120.84	109.28
2	B	501	PAF	C5-C3-C2	5.99	120.28	109.34

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	PAF	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, 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	285/287 (99%)	-0.07	3 (1%) 82 83	10, 18, 30, 38	0
1	B	284/287 (98%)	0.51	29 (10%) 9 8	11, 24, 42, 51	0
1	C	285/287 (99%)	-0.06	5 (1%) 71 72	13, 22, 33, 38	0
1	D	283/287 (98%)	1.71	95 (33%) 0 0	14, 33, 80, 96	0
All	All	1137/1148 (99%)	0.52	132 (11%) 6 5	10, 23, 55, 96	0

All (132) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	217	LEU	11.4
1	D	256	VAL	10.5
1	D	194	THR	9.6
1	D	200	ILE	8.7
1	D	234	LEU	8.6
1	D	225	LEU	8.4
1	D	236	VAL	8.4
1	D	231	ALA	7.9
1	D	269	LEU	7.8
1	D	254	LEU	7.8
1	B	192[A]	TYR	7.6
1	D	226	LEU	7.5
1	D	282	LEU	7.4
1	D	211	LEU	7.3
1	D	224	ALA	7.3
1	D	243	LEU	7.0
1	D	270	GLY	7.0
1	D	267	ALA	6.8
1	D	216	ALA	6.6
1	D	222	ILE	6.6
1	D	268[A]	TRP	6.5

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Mol	Chain	Res	Type	RSRZ
1	D	218	GLY	6.4
1	D	209	TRP	6.2
1	D	213	GLU	6.2
1	D	281	ASP	6.1
1	B	256	VAL	5.9
1	D	199	GLN	5.9
1	D	65	PRO	5.8
1	D	229	ALA	5.5
1	B	271	LYS	5.4
1	D	192	TYR	5.3
1	D	221	GLN	5.3
1	D	232	GLN	5.3
1	D	0	ALA	5.2
1	D	259	GLN	5.1
1	D	239	THR	5.1
1	D	228	GLU	5.1
1	D	238	PHE	5.0
1	D	233	LEU	5.0
1	D	196[A]	GLU	5.0
1	D	255	THR	5.0
1	D	240	PRO	4.9
1	B	268[A]	TRP	4.9
1	D	69	ALA	4.8
1	D	210	ALA	4.8
1	D	257	ASP	4.7
1	D	68	LEU	4.6
1	B	239	THR	4.5
1	D	235	ARG	4.5
1	D	272	ALA	4.5
1	D	189	ARG	4.4
1	D	219	GLU	4.4
1	D	220	ARG	4.3
1	D	245	ILE	4.3
1	D	230	ALA	4.3
1	D	280	VAL	4.2
1	D	223	ASP	4.1
1	B	191	GLY	4.1
1	D	274[A]	LEU	4.1
1	D	215	MET	4.1
1	D	191	GLY	4.1
1	D	260	GLN	4.0
1	D	266	ALA	4.0

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Mol	Chain	Res	Type	RSRZ
1	D	183	GLY	3.9
1	D	195	GLU	3.9
1	D	258	SER	3.8
1	B	257[A]	ASP	3.7
1	C	0	ALA	3.7
1	D	208	MET	3.6
1	B	200	ILE	3.6
1	B	234	LEU	3.6
1	D	271	LYS	3.6
1	D	212	ALA	3.5
1	D	66	ASP	3.5
1	D	70	HIS	3.4
1	A	101	ALA	3.2
1	B	238	PHE	3.1
1	B	133	ILE	3.1
1	D	279	LEU	3.1
1	D	214	LYS	3.1
1	D	202	PRO	3.1
1	D	227	GLU	3.0
1	D	203	GLN	3.0
1	D	193	LEU	3.0
1	B	159[A]	ILE	3.0
1	D	201	ALA	3.0
1	A	0	ALA	2.9
1	D	273	ARG	2.9
1	D	71	TYR	2.9
1	B	236	VAL	2.9
1	D	122	SER	2.8
1	B	134	VAL	2.8
1	D	67	ASP	2.8
1	D	76	GLN	2.8
1	D	265	MET	2.8
1	B	194	THR	2.7
1	D	138	PHE	2.6
1	D	64[A]	ARG	2.6
1	B	207	ILE	2.6
1	D	206	LYS	2.6
1	D	197	GLU	2.6
1	B	195	GLU	2.5
1	D	159[A]	ILE	2.5
1	D	241	ASP	2.5
1	B	230	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	272	ALA	2.4
1	D	261	ALA	2.4
1	D	134	VAL	2.4
1	D	124	PRO	2.4
1	C	109	VAL	2.4
1	B	189	ARG	2.4
1	D	74	THR	2.4
1	B	223[A]	ASP	2.3
1	D	190	ASN	2.3
1	B	229	ALA	2.3
1	C	168	TYR	2.2
1	B	217	LEU	2.2
1	A	195[A]	GLU	2.2
1	B	227	GLU	2.2
1	B	242	GLU	2.2
1	D	244	PHE	2.2
1	D	168	TYR	2.2
1	B	137	LEU	2.1
1	B	213[A]	GLU	2.1
1	D	242[A]	GLU	2.1
1	C	284	HIS	2.1
1	B	70	HIS	2.1
1	C	163	VAL	2.1
1	D	263	ILE	2.0
1	D	182	ASP	2.0
1	B	233	LEU	2.0
1	D	249	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, 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(Å ²)	Q<0.9
2	PAF	C	501	10/10	0.90	0.14	1.41	15,18,26,29	0
2	PAF	A	501	10/10	0.96	0.11	0.66	13,16,23,26	0
3	CL	C	285	1/1	0.99	0.09	0.34	22,22,22,22	0
3	CL	A	285	1/1	1.00	0.10	-0.22	19,19,19,19	0
2	PAF	B	501	10/10	0.95	0.11	-0.57	13,16,27,32	0
2	PAF	D	501	10/10	0.93	0.08	-0.68	19,21,30,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.