Full wwPDB X-ray Structure Validation Report (i)

Aug 16, 2023 – 01:35 PM EDT

PDB ID	:	1ZY1
Title	:	X-ray structure of peptide deformylase from Arabidopsis thaliana (AtPDF1A) in complex with Met-Ala-Ser
Authors	:	Fieulaine, S.; Juillan-Binard, C.; Serero, A.; Dardel, F.; Giglione, C.; Meinnel, T.; Ferrer, J.-L.
Deposited on	:	2005-06-09
Resolution	:	3.00 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

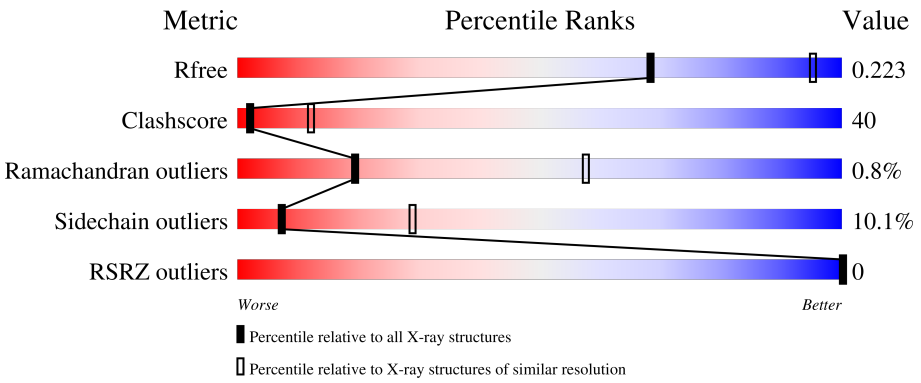
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 3.00 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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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 <=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	197	37% 50% 9% • •
1	B	197	37% 51% 9% •
2	D	3	33% 67%
2	E	3	33% 33% 33%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3056 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 Peptide deformylase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	191	1478	934	259	277	8	0	0	0
1	B	191	1470	929	260	273	8	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP Q9FV53
A	191	SER	-	expression tag	UNP Q9FV53
A	192	HIS	-	expression tag	UNP Q9FV53
A	193	HIS	-	expression tag	UNP Q9FV53
A	194	HIS	-	expression tag	UNP Q9FV53
A	195	HIS	-	expression tag	UNP Q9FV53
A	196	HIS	-	expression tag	UNP Q9FV53
A	197	HIS	-	expression tag	UNP Q9FV53
B	1	MET	-	initiating methionine	UNP Q9FV53
B	191	SER	-	expression tag	UNP Q9FV53
B	192	HIS	-	expression tag	UNP Q9FV53
B	193	HIS	-	expression tag	UNP Q9FV53
B	194	HIS	-	expression tag	UNP Q9FV53
B	195	HIS	-	expression tag	UNP Q9FV53
B	196	HIS	-	expression tag	UNP Q9FV53
B	197	HIS	-	expression tag	UNP Q9FV53

- Molecule 2 is a protein called tripeptide fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	D	3	20	11	3	5	1	0	0	0
2	E	3	20	11	3	5	1	0	0	0

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 4 is water.

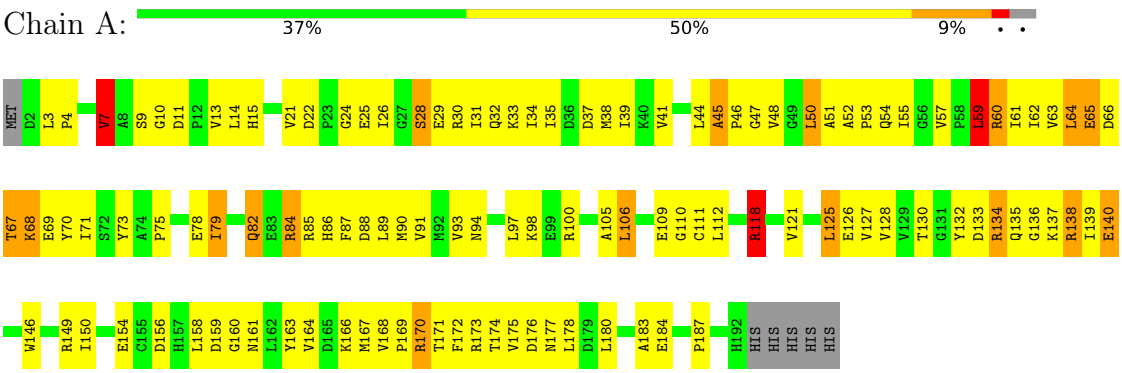
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	30	30	30	0	0
4	B	35	35	35	0	0
4	D	1	1	1	0	0

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.

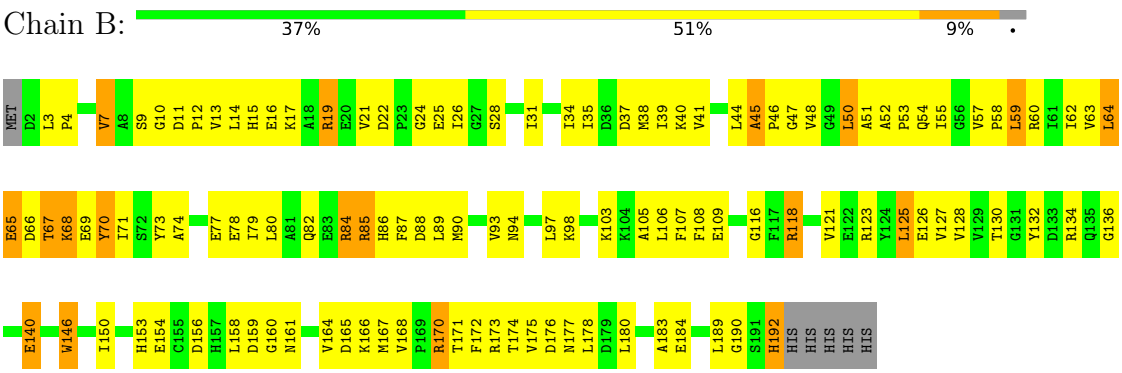
- Molecule 1: Peptide deformylase, mitochondrial

Chain A:



- Molecule 1: Peptide deformylase, mitochondrial

Chain B:

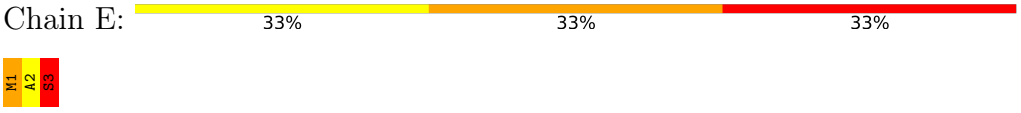


- Molecule 2: tripeptide fragment

Chain D:

- Molecule 2: tripeptide fragment

Chain E:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	51.60Å 76.30Å 109.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 3.00 46.66 – 2.98	Depositor EDS
% Data completeness (in resolution range)	97.2 (30.00-3.00) 99.4 (46.66-2.98)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	1.19 (at 2.96Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.226 , 0.256 0.228 , 0.223	Depositor DCC
R_{free} test set	461 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	57.1	Xtriage
Anisotropy	0.713	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 71.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	3056	wwPDB-VP
Average B, all atoms (Å^2)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.66% 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, 0.333 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: ZN

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	1.18	15/1505 (1.0%)	1.31	24/2040 (1.2%)
1	B	1.13	13/1497 (0.9%)	1.16	25/2032 (1.2%)
2	D	0.85	0/19	0.57	0/22
2	E	1.81	1/19 (5.3%)	2.11	1/22 (4.5%)
All	All	1.16	29/3040 (1.0%)	1.24	50/4116 (1.2%)

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

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	70	TYR	CE1-CZ	-17.20	1.16	1.38
1	A	70	TYR	CE1-CZ	-16.66	1.16	1.38
1	A	70	TYR	CE2-CZ	-16.65	1.17	1.38
1	B	70	TYR	CE2-CZ	-16.27	1.17	1.38
1	A	70	TYR	CG-CD2	-15.09	1.19	1.39
1	B	70	TYR	CG-CD2	-14.95	1.19	1.39
1	A	70	TYR	CG-CD1	-14.35	1.20	1.39
1	B	70	TYR	CG-CD1	-13.96	1.21	1.39
1	A	140	GLU	CD-OE2	-9.36	1.15	1.25
1	B	140	GLU	CD-OE2	-9.02	1.15	1.25
1	A	59	LEU	CG-CD1	-8.01	1.22	1.51
1	B	59	LEU	CG-CD1	-7.78	1.23	1.51
1	A	118	ARG	CZ-NH2	-7.59	1.23	1.33
1	A	73	TYR	CE1-CZ	-7.57	1.28	1.38

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	7	VAL	CB-CG1	-7.47	1.37	1.52
1	A	140	GLU	CD-OE1	-7.39	1.17	1.25
1	B	73	TYR	CE1-CZ	-6.95	1.29	1.38
1	A	84	ARG	CZ-NH1	-6.91	1.24	1.33
1	B	146	TRP	NE1-CE2	6.86	1.46	1.37
1	B	7	VAL	CB-CG2	-6.56	1.39	1.52
1	A	7	VAL	CB-CG2	-6.54	1.39	1.52
1	B	140	GLU	CD-OE1	-6.33	1.18	1.25
1	A	73	TYR	CG-CD2	-6.29	1.30	1.39
1	B	7	VAL	CB-CG1	-6.28	1.39	1.52
1	B	73	TYR	CG-CD2	-6.28	1.30	1.39
1	A	118	ARG	CZ-NH1	-6.05	1.25	1.33
1	B	84	ARG	CZ-NH1	-6.05	1.25	1.33
1	A	59	LEU	CG-CD2	-5.75	1.30	1.51
2	E	3	SER	CA-CB	-5.58	1.44	1.52

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	85	ARG	NE-CZ-NH1	21.11	130.85	120.30
1	A	85	ARG	NE-CZ-NH2	-16.01	112.29	120.30
1	A	118	ARG	NE-CZ-NH2	15.57	128.08	120.30
1	B	118	ARG	NE-CZ-NH1	14.52	127.56	120.30
1	A	170	ARG	NE-CZ-NH1	-13.85	113.38	120.30
1	A	170	ARG	NE-CZ-NH2	13.70	127.15	120.30
1	B	170	ARG	NE-CZ-NH2	-13.01	113.79	120.30
1	B	170	ARG	NE-CZ-NH1	12.81	126.71	120.30
1	A	140	GLU	OE1-CD-OE2	-12.09	108.80	123.30
1	B	140	GLU	OE1-CD-OE2	-11.52	109.47	123.30
1	A	85	ARG	CD-NE-CZ	9.88	137.43	123.60
1	B	118	ARG	NH1-CZ-NH2	-8.91	109.59	119.40
1	A	118	ARG	NH1-CZ-NH2	-8.59	109.95	119.40
1	A	70	TYR	CB-CG-CD1	8.24	125.94	121.00
1	B	84	ARG	CG-CD-NE	7.98	128.56	111.80
1	B	85	ARG	NE-CZ-NH2	7.97	124.28	120.30
1	A	84	ARG	CG-CD-NE	7.65	127.87	111.80
1	A	59	LEU	CB-CG-CD2	7.23	123.28	111.00
1	B	70	TYR	CZ-CE2-CD2	7.11	126.20	119.80
1	A	170	ARG	CD-NE-CZ	7.04	133.46	123.60
1	B	59	LEU	CB-CG-CD2	6.91	122.74	111.00
1	B	70	TYR	CD1-CG-CD2	-6.88	110.33	117.90
1	A	70	TYR	CZ-CE2-CD2	6.85	125.96	119.80

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	70	TYR	CD1-CG-CD2	-6.83	110.39	117.90
1	B	170	ARG	CD-NE-CZ	6.82	133.15	123.60
1	A	70	TYR	CD1-CE1-CZ	6.64	125.78	119.80
1	B	70	TYR	CD1-CE1-CZ	6.58	125.72	119.80
2	E	2	ALA	O-C-N	6.49	133.08	122.70
1	B	70	TYR	CB-CG-CD2	6.39	124.83	121.00
1	B	85	ARG	CG-CD-NE	6.35	125.13	111.80
1	B	68	LYS	CD-CE-NZ	-6.25	97.33	111.70
1	B	70	TYR	CB-CG-CD1	6.23	124.74	121.00
1	B	84	ARG	NH1-CZ-NH2	-6.20	112.58	119.40
1	B	146	TRP	CE2-CD2-CG	6.12	112.19	107.30
1	B	84	ARG	NE-CZ-NH2	6.06	123.33	120.30
1	A	73	TYR	CB-CG-CD1	6.00	124.60	121.00
1	A	70	TYR	CE1-CZ-CE2	-5.96	110.27	119.80
1	B	70	TYR	CE1-CZ-CE2	-5.88	110.39	119.80
1	B	73	TYR	CB-CG-CD1	5.87	124.52	121.00
1	A	45	ALA	N-CA-C	-5.82	95.28	111.00
1	B	84	ARG	NE-CZ-NH1	5.68	123.14	120.30
1	A	68	LYS	CD-CE-NZ	-5.67	98.65	111.70
1	A	60	ARG	NE-CZ-NH1	-5.66	117.47	120.30
1	A	84	ARG	NE-CZ-NH1	5.55	123.07	120.30
1	A	84	ARG	NH1-CZ-NH2	-5.46	113.40	119.40
1	B	146	TRP	CD2-CE2-CZ2	5.39	128.76	122.30
1	B	45	ALA	N-CA-C	-5.27	96.78	111.00
1	A	59	LEU	CD1-CG-CD2	-5.17	94.98	110.50
1	B	59	LEU	CD1-CG-CD2	-5.11	95.17	110.50
1	A	84	ARG	NE-CZ-NH2	5.01	122.81	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	118	ARG	Sidechain

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	1478	0	1463	128	0
1	B	1470	0	1451	122	0
2	D	20	0	21	7	0
2	E	20	0	21	5	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	30	0	0	3	0
4	B	35	0	0	1	0
4	D	1	0	0	0	0
All	All	3056	0	2956	239	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 40.

All (239) 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:B:7:VAL:CG2	1:B:14:LEU:HD11	1.82	1.09
1:B:7:VAL:HG21	1:B:14:LEU:HD11	1.51	0.92
1:B:65:GLU:HB2	1:B:89:LEU:HA	1.53	0.89
1:A:130:THR:HG22	1:A:140:GLU:HG3	1.58	0.86
1:A:78:GLU:HG2	1:A:82:GLN:HE21	1.41	0.85
1:B:7:VAL:HG23	1:B:14:LEU:HD11	1.56	0.84
1:A:177:ASN:HB3	1:A:180:LEU:HD12	1.59	0.84
1:B:177:ASN:HB3	1:B:180:LEU:HD12	1.58	0.83
1:A:10:GLY:H	1:A:15:HIS:HE1	1.26	0.83
1:B:150:ILE:HG12	2:E:1:MET:CE	2.09	0.82
1:B:109:GLU:HG2	2:E:1:MET:SD	2.22	0.80
1:A:7:VAL:HB	1:A:14:LEU:HD11	1.63	0.80
1:B:10:GLY:H	1:B:15:HIS:HE1	1.29	0.80
1:A:134:ARG:HG2	1:A:134:ARG:HH11	1.47	0.79
1:A:21:VAL:HG21	1:A:31:ILE:HD13	1.62	0.79
1:A:134:ARG:HH11	1:A:134:ARG:CG	1.95	0.79
1:A:39:ILE:HD13	1:A:89:LEU:HD21	1.65	0.77
1:B:130:THR:HG22	1:B:140:GLU:HG3	1.66	0.77
1:A:71:ILE:HG12	1:A:84:ARG:HD3	1.67	0.76
1:A:26:ILE:HD13	1:A:93:VAL:HG21	1.67	0.74
1:B:93:VAL:CG2	1:B:132:TYR:HB2	2.17	0.74
1:A:105:ALA:HB2	1:A:125:LEU:CD2	2.18	0.73
1:A:48:VAL:HG13	2:D:3:SER:HB2	1.70	0.73
1:A:21:VAL:CG2	1:A:31:ILE:HD13	2.20	0.72
1:B:192:HIS:ND1	1:B:192:HIS:C	2.42	0.72

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:MET:HE3	1:B:63:VAL:HG13	1.71	0.72
1:B:60:ARG:NH1	1:B:159:ASP:OD1	2.22	0.72
1:B:156:ASP:CG	1:B:166:LYS:HZ1	1.92	0.72
1:A:48:VAL:HG13	2:D:3:SER:CB	2.20	0.72
1:A:78:GLU:HG2	1:A:82:GLN:NE2	2.04	0.72
1:B:156:ASP:OD2	1:B:166:LYS:NZ	2.23	0.71
1:A:105:ALA:HB2	1:A:125:LEU:HD22	1.73	0.71
1:B:3:LEU:HD11	1:B:40:LYS:HB3	1.72	0.71
1:B:105:ALA:HB2	1:B:125:LEU:CD2	2.21	0.71
1:B:66:ASP:HB2	1:B:87:PHE:CE1	2.26	0.70
1:A:156:ASP:OD2	1:A:166:LYS:NZ	2.27	0.68
1:B:3:LEU:HD11	1:B:40:LYS:CB	2.23	0.68
1:B:105:ALA:HB2	1:B:125:LEU:HD22	1.74	0.67
1:B:93:VAL:HG23	1:B:132:TYR:HB2	1.77	0.67
1:A:65:GLU:HB2	1:A:89:LEU:HA	1.76	0.66
1:B:156:ASP:CG	1:B:166:LYS:NZ	2.48	0.66
1:B:174:THR:HG22	1:B:176:ASP:H	1.60	0.66
1:A:93:VAL:CG2	1:A:132:TYR:HB2	2.25	0.66
1:A:156:ASP:CG	1:A:166:LYS:NZ	2.50	0.66
1:A:10:GLY:H	1:A:15:HIS:CE1	2.12	0.65
1:A:32:GLN:OE1	1:A:134:ARG:NH1	2.28	0.65
1:A:174:THR:HG22	1:A:176:ASP:H	1.62	0.65
1:B:77:GLU:O	1:B:80:LEU:HB3	1.96	0.65
1:B:10:GLY:H	1:B:15:HIS:CE1	2.15	0.64
1:B:150:ILE:HG12	2:E:1:MET:HE2	1.77	0.64
1:A:15:HIS:HB2	1:B:183:ALA:HB2	1.79	0.63
1:B:48:VAL:HG13	2:E:3:SER:HB2	1.80	0.63
1:A:167:MET:CE	1:A:172:PHE:HB2	2.29	0.62
1:B:38:MET:CE	1:B:63:VAL:HG13	2.28	0.62
1:A:156:ASP:CG	1:A:166:LYS:HZ1	2.03	0.61
1:B:153:HIS:CG	2:E:1:MET:HG3	2.35	0.61
1:B:68:LYS:HG3	1:B:86:HIS:CG	2.35	0.61
1:A:109:GLU:CG	2:D:1:MET:HE1	2.31	0.60
1:A:60:ARG:NH1	1:A:159:ASP:OD1	2.33	0.60
1:A:183:ALA:HB1	1:B:16:GLU:HG2	1.82	0.60
1:B:57:VAL:HG12	1:B:59:LEU:HG	1.83	0.60
1:A:9:SER:HB3	1:B:173:ARG:HA	1.84	0.59
1:A:25:GLU:O	1:A:28:SER:HB3	2.01	0.59
1:B:7:VAL:HG21	1:B:14:LEU:CD1	2.30	0.59
1:B:167:MET:HE3	1:B:172:PHE:HB2	1.83	0.59
1:A:7:VAL:CB	1:A:14:LEU:HD11	2.32	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:ASP:HB2	1:A:87:PHE:CE1	2.37	0.59
1:A:93:VAL:HG23	1:A:132:TYR:HB2	1.85	0.59
1:B:39:ILE:HD13	1:B:89:LEU:HD21	1.82	0.59
1:B:62:ILE:HG23	1:B:158:LEU:HD11	1.84	0.59
1:B:21:VAL:HG12	1:B:94:ASN:OD1	2.03	0.58
1:A:161:ASN:OD1	1:B:170:ARG:NH1	2.36	0.58
1:B:177:ASN:CB	1:B:180:LEU:HD12	2.31	0.58
1:A:134:ARG:CG	1:A:134:ARG:NH1	2.63	0.58
1:A:177:ASN:CB	1:A:180:LEU:HD12	2.31	0.57
1:A:67:THR:C	1:A:69:GLU:H	2.07	0.57
1:A:154:GLU:OE2	2:D:1:MET:N	2.35	0.57
1:B:26:ILE:HD13	1:B:93:VAL:HG21	1.85	0.57
1:B:97:LEU:HG	1:B:127:VAL:HG21	1.86	0.57
1:A:66:ASP:O	1:A:86:HIS:HB2	2.04	0.57
1:B:19:ARG:HG3	1:B:58:PRO:O	2.04	0.57
1:A:26:ILE:HD12	1:A:136:GLY:HA2	1.86	0.57
1:A:75:PRO:O	1:A:79:ILE:HG23	2.05	0.56
1:A:100:ARG:HE	1:A:126:GLU:HB3	1.70	0.56
1:B:74:ALA:HB3	1:B:79:ILE:HD11	1.87	0.56
1:B:167:MET:CE	1:B:172:PHE:HB2	2.34	0.56
1:A:110:GLY:HA3	1:A:118:ARG:HE	1.70	0.56
1:A:168:VAL:HB	1:A:171:THR:HG21	1.87	0.56
1:B:3:LEU:HD22	1:B:4:PRO:HD2	1.87	0.56
1:B:68:LYS:HD2	1:B:86:HIS:CE1	2.41	0.56
1:A:66:ASP:OD1	1:A:84:ARG:NH1	2.39	0.56
1:A:21:VAL:CG2	1:A:31:ILE:CD1	2.84	0.55
1:B:108:PHE:CE1	1:B:189:LEU:HD11	2.42	0.55
1:A:44:LEU:HD12	1:A:44:LEU:O	2.07	0.55
1:B:78:GLU:O	1:B:82:GLN:HG3	2.07	0.55
1:B:89:LEU:HD12	1:B:90:MET:H	1.72	0.55
1:B:168:VAL:HB	1:B:171:THR:HG21	1.89	0.55
1:A:52:ALA:HB3	1:A:53:PRO:HD3	1.88	0.55
1:A:67:THR:HB	1:A:69:GLU:H	1.72	0.55
1:B:67:THR:HB	1:B:69:GLU:H	1.72	0.55
1:A:94:ASN:O	4:A:217:HOH:O	2.18	0.54
1:B:52:ALA:HB3	1:B:53:PRO:HD3	1.89	0.54
1:A:109:GLU:HG2	2:D:1:MET:HE1	1.88	0.54
1:A:133:ASP:OD1	1:A:137:LYS:N	2.39	0.54
1:A:173:ARG:HA	1:B:9:SER:HB3	1.89	0.54
1:B:71:ILE:O	1:B:79:ILE:HD11	2.08	0.53
1:A:97:LEU:HG	1:A:127:VAL:HG21	1.88	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:97:LEU:HG	1:B:127:VAL:CG2	2.38	0.53
1:A:57:VAL:CG1	1:A:59:LEU:HD23	2.39	0.53
1:A:112:LEU:HG	4:A:199:HOH:O	2.09	0.53
1:A:66:ASP:CG	1:A:84:ARG:HH11	2.12	0.53
1:B:55:ILE:HG13	1:B:57:VAL:HG23	1.91	0.53
1:A:13:VAL:CG1	1:A:14:LEU:HD12	2.39	0.53
1:A:71:ILE:HG12	1:A:84:ARG:CD	2.37	0.53
1:A:109:GLU:HG2	2:D:1:MET:CE	2.39	0.53
1:A:89:LEU:HD12	1:A:90:MET:H	1.74	0.52
1:B:97:LEU:CG	1:B:127:VAL:HG21	2.39	0.52
1:A:160:GLY:O	1:B:170:ARG:HD3	2.10	0.52
1:A:13:VAL:HG13	1:A:14:LEU:HD12	1.91	0.52
1:A:21:VAL:HG23	1:A:31:ILE:CD1	2.40	0.52
1:B:84:ARG:NH1	1:B:146:TRP:NE1	2.58	0.52
1:A:24:GLY:C	1:A:26:ILE:N	2.63	0.52
1:A:167:MET:HE1	1:A:172:PHE:HB2	1.91	0.52
1:A:10:GLY:O	4:A:227:HOH:O	2.19	0.52
1:B:57:VAL:CG1	1:B:59:LEU:HG	2.40	0.52
1:A:38:MET:CE	1:A:63:VAL:HG13	2.40	0.51
1:B:24:GLY:C	1:B:26:ILE:N	2.63	0.51
1:A:84:ARG:HG3	1:A:146:TRP:CD1	2.46	0.51
1:B:97:LEU:HD23	1:B:98:LYS:N	2.25	0.51
1:A:168:VAL:O	1:A:171:THR:HG23	2.11	0.50
1:B:64:LEU:HD23	1:B:64:LEU:N	2.26	0.50
1:B:44:LEU:O	1:B:44:LEU:HD12	2.12	0.50
1:A:64:LEU:N	1:A:64:LEU:HD23	2.26	0.50
1:B:13:VAL:HG13	1:B:14:LEU:HD12	1.93	0.50
1:A:173:ARG:HD3	1:A:178:LEU:HA	1.94	0.50
1:B:192:HIS:CE1	4:B:219:HOH:O	2.65	0.50
1:A:167:MET:HE3	1:A:172:PHE:HB2	1.93	0.49
1:B:21:VAL:CG2	1:B:31:ILE:HD13	2.42	0.49
1:B:71:ILE:HD12	1:B:86:HIS:HB3	1.94	0.49
1:B:24:GLY:C	1:B:26:ILE:H	2.15	0.49
1:A:55:ILE:HG13	1:A:57:VAL:HG23	1.95	0.49
1:B:67:THR:OG1	1:B:70:TYR:CD1	2.65	0.49
1:A:28:SER:O	1:A:32:GLN:HG2	2.12	0.49
1:A:24:GLY:C	1:A:26:ILE:H	2.15	0.48
1:B:97:LEU:HD23	1:B:97:LEU:C	2.34	0.48
1:A:97:LEU:HG	1:A:127:VAL:CG2	2.43	0.48
1:B:67:THR:C	1:B:69:GLU:H	2.15	0.48
1:A:29:GLU:HG2	1:A:29:GLU:O	2.14	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:ILE:O	1:A:39:ILE:HG13	2.14	0.47
1:B:35:ILE:O	1:B:39:ILE:HG13	2.15	0.47
1:B:13:VAL:CG1	1:B:14:LEU:HD12	2.44	0.47
1:B:164:VAL:O	1:B:167:MET:HG2	2.15	0.47
1:B:173:ARG:HD3	1:B:178:LEU:HA	1.96	0.47
1:A:97:LEU:CG	1:A:127:VAL:HG21	2.45	0.47
1:A:149:ARG:HG2	2:D:1:MET:CE	2.45	0.47
1:B:3:LEU:HD11	1:B:40:LYS:HB2	1.95	0.47
1:B:71:ILE:HG22	1:B:79:ILE:HD13	1.96	0.47
1:A:84:ARG:HG3	1:A:146:TRP:CG	2.50	0.47
1:A:9:SER:CB	1:B:173:ARG:HA	2.44	0.47
1:A:21:VAL:HG12	1:A:94:ASN:OD1	2.15	0.47
1:A:67:THR:C	1:A:69:GLU:N	2.67	0.46
1:A:82:GLN:OE1	1:A:146:TRP:CZ3	2.68	0.46
1:A:50:LEU:HD22	1:A:54:GLN:OE1	2.15	0.46
1:B:168:VAL:O	1:B:171:THR:HG23	2.15	0.46
1:A:39:ILE:CD1	1:A:89:LEU:HD21	2.40	0.46
1:A:22:ASP:HB3	1:A:25:GLU:HG3	1.97	0.46
1:A:45:ALA:O	1:A:47:GLY:N	2.49	0.46
1:A:61:ILE:HG12	1:A:93:VAL:HG12	1.98	0.46
1:A:133:ASP:HB3	1:A:139:ILE:HG21	1.97	0.45
1:A:38:MET:HE3	1:A:63:VAL:HG13	1.97	0.45
1:B:52:ALA:HB1	1:B:57:VAL:HB	1.98	0.45
1:A:173:ARG:HA	1:B:9:SER:CB	2.46	0.45
1:B:67:THR:C	1:B:69:GLU:N	2.69	0.45
1:A:118:ARG:NH2	1:A:175:VAL:HG12	2.31	0.45
1:B:71:ILE:CG2	1:B:79:ILE:HD13	2.47	0.45
1:B:17:LYS:HE2	1:B:60:ARG:HH11	1.82	0.45
1:B:66:ASP:O	1:B:86:HIS:HB2	2.16	0.45
1:B:107:PHE:CZ	1:B:190:GLY:HA2	2.52	0.44
1:A:146:TRP:O	1:A:150:ILE:HG13	2.18	0.44
1:B:34:ILE:HD13	1:B:57:VAL:HG11	1.98	0.44
1:B:50:LEU:HD22	1:B:54:GLN:OE1	2.17	0.44
1:A:173:ARG:CD	1:A:178:LEU:HA	2.47	0.44
1:A:170:ARG:NH1	1:B:161:ASN:OD1	2.50	0.44
1:B:50:LEU:HD13	1:B:51:ALA:N	2.32	0.44
1:A:68:LYS:HG3	1:A:86:HIS:CG	2.53	0.44
1:A:11:ASP:OD1	1:A:13:VAL:HG12	2.18	0.44
1:B:25:GLU:O	1:B:28:SER:HB3	2.18	0.44
1:B:98:LYS:O	1:B:127:VAL:HG23	2.18	0.44
1:B:25:GLU:O	1:B:31:ILE:HD12	2.18	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:VAL:CG2	1:A:14:LEU:HD11	2.47	0.43
1:A:14:LEU:O	1:A:160:GLY:HA2	2.18	0.43
1:A:52:ALA:HB1	1:A:57:VAL:HB	2.00	0.43
1:B:123:ARG:HH12	1:B:166:LYS:HZ3	1.66	0.43
1:A:62:ILE:HG23	1:A:158:LEU:HD11	2.01	0.43
1:A:111:CYS:CB	1:A:163:TYR:HE2	2.31	0.43
1:B:173:ARG:CD	1:B:178:LEU:HA	2.49	0.43
1:A:50:LEU:HD13	1:A:51:ALA:N	2.33	0.43
1:B:37:ASP:O	1:B:41:VAL:HG23	2.18	0.43
1:B:3:LEU:HD13	1:B:41:VAL:HG22	2.01	0.43
1:B:14:LEU:O	1:B:160:GLY:HA2	2.18	0.43
1:B:97:LEU:HD21	1:B:127:VAL:HG21	2.01	0.43
1:B:116:GLY:C	1:B:175:VAL:HG13	2.40	0.43
1:A:164:VAL:O	1:A:167:MET:HG2	2.18	0.42
1:B:26:ILE:HD12	1:B:136:GLY:HA2	2.01	0.42
1:B:146:TRP:O	1:B:150:ILE:HG13	2.19	0.42
1:A:15:HIS:HB2	1:B:183:ALA:CB	2.47	0.42
1:B:11:ASP:OD1	1:B:13:VAL:HG12	2.20	0.42
1:A:110:GLY:HA3	1:A:118:ARG:NE	2.33	0.42
1:A:169:PRO:HA	1:B:165:ASP:OD1	2.20	0.42
1:A:97:LEU:HD23	1:A:98:LYS:N	2.34	0.42
1:B:45:ALA:O	1:B:47:GLY:N	2.52	0.42
1:A:34:ILE:HD13	1:A:57:VAL:HG11	2.01	0.42
1:B:12:PRO:O	1:B:16:GLU:HG3	2.19	0.42
1:B:22:ASP:HB3	1:B:25:GLU:HG3	2.00	0.42
1:A:146:TRP:CZ3	1:A:149:ARG:HD2	2.55	0.42
1:A:127:VAL:HG22	1:A:128:VAL:N	2.34	0.41
1:A:106:LEU:HD11	1:A:187:PRO:HG2	2.02	0.41
1:B:3:LEU:HD21	1:B:37:ASP:HA	2.02	0.41
1:B:31:ILE:O	1:B:35:ILE:HG13	2.20	0.41
1:A:37:ASP:O	1:A:41:VAL:HG23	2.20	0.41
1:B:150:ILE:O	1:B:154:GLU:HG2	2.20	0.41
1:A:26:ILE:CD1	1:A:93:VAL:HG21	2.43	0.41
1:A:30:ARG:O	1:A:33:LYS:HB3	2.20	0.41
1:A:78:GLU:CG	1:A:82:GLN:NE2	2.80	0.41
1:B:97:LEU:CD2	1:B:127:VAL:HG21	2.51	0.41
1:B:127:VAL:HG22	1:B:128:VAL:N	2.36	0.41
1:A:3:LEU:HD12	1:A:4:PRO:HD2	2.02	0.41
1:B:3:LEU:CD2	1:B:4:PRO:HD2	2.50	0.41
1:A:63:VAL:HG12	1:A:91:VAL:HG22	2.02	0.40
1:A:89:LEU:HD12	1:A:90:MET:N	2.36	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:ASP:HB3	1:B:86:HIS:HA	2.03	0.40
1:A:97:LEU:HD23	1:A:97:LEU:C	2.42	0.40
1:A:150:ILE:O	1:A:154:GLU:HG2	2.21	0.40
1:B:89:LEU:HD12	1:B:90:MET:N	2.35	0.40
1:A:138:ARG:H	1:A:138:ARG:HG2	1.69	0.40
1:B:103:LYS:HE3	1:B:103:LYS:HB2	1.80	0.40
1:A:118:ARG:HH21	1:A:118:ARG:CG	2.35	0.40
1:B:103:LYS:NZ	1:B:126:GLU:OE1	2.50	0.40
1:B:178:LEU:HD12	1:B:178:LEU:O	2.21	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	189/197 (96%)	171 (90%)	16 (8%)	2 (1%)	14 50
1	B	189/197 (96%)	170 (90%)	18 (10%)	1 (0%)	29 68
2	D	1/3 (33%)	1 (100%)	0	0	100 100
2	E	1/3 (33%)	1 (100%)	0	0	100 100
All	All	380/400 (95%)	343 (90%)	34 (9%)	3 (1%)	19 57

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28	SER
1	B	46	PRO
1	A	46	PRO

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	157/170 (92%)	140 (89%)	17 (11%)	6 26
1	B	155/170 (91%)	142 (92%)	13 (8%)	11 38
2	D	2/2 (100%)	2 (100%)	0	100 100
2	E	2/2 (100%)	0	2 (100%)	0 0
All	All	316/344 (92%)	284 (90%)	32 (10%)	7 29

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	VAL
1	A	50	LEU
1	A	59	LEU
1	A	64	LEU
1	A	65	GLU
1	A	67	THR
1	A	79	ILE
1	A	82	GLN
1	A	88	ASP
1	A	106	LEU
1	A	118	ARG
1	A	121	VAL
1	A	125	LEU
1	A	134	ARG
1	A	135	GLN
1	A	138	ARG
1	A	184	GLU
1	B	19	ARG
1	B	50	LEU
1	B	64	LEU
1	B	65	GLU
1	B	67	THR
1	B	85	ARG
1	B	88	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	106	LEU
1	B	121	VAL
1	B	125	LEU
1	B	134	ARG
1	B	184	GLU
1	B	192	HIS
2	E	1	MET
2	E	3	SER

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

Mol	Chain	Res	Type
1	A	15	HIS
1	A	82	GLN
1	A	135	GLN
1	B	15	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)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

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 (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	191/197 (96%)	-0.20	0	100	100	29, 48, 70, 89	0
1	B	191/197 (96%)	-0.15	0	100	100	32, 49, 76, 90	0
2	D	3/3 (100%)	0.31	0	100	100	42, 42, 44, 45	0
2	E	3/3 (100%)	0.53	0	100	100	56, 56, 62, 66	0
All	All	388/400 (97%)	-0.17	0	100	100	29, 48, 75, 90	0

There are no RSRZ outliers to report.

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
3	ZN	A	198	1/1	0.96	0.11	38,38,38,38	0
3	ZN	B	198	1/1	0.98	0.11	42,42,42,42	0

6.5 Other polymers (i)

There are no such residues in this entry.