



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 07:01 PM UTC

PDB ID : 1A16 / pdb_00001a16
Title : AMINOPEPTIDASE P FROM E. COLI WITH THE INHIBITOR PRO-LEU
Authors : Wilce, M.C.; Bond, C.S.; Lilley, P.E.; Dixon, N.E.; Freeman, H.C.; Guss, J.M.
Deposited on : 1997-12-22
Resolution : 2.30 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 ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

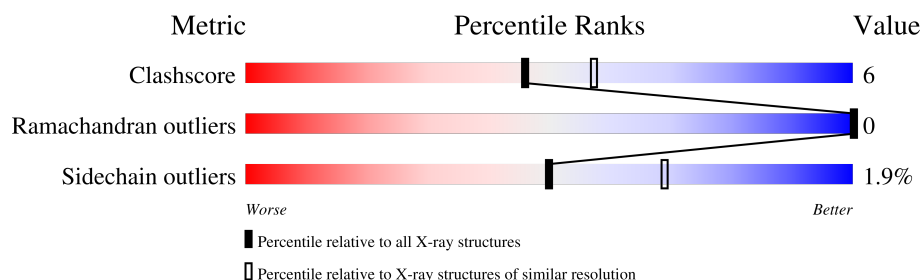
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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))
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	440	 82% 15% .

2 Entry composition [i](#)

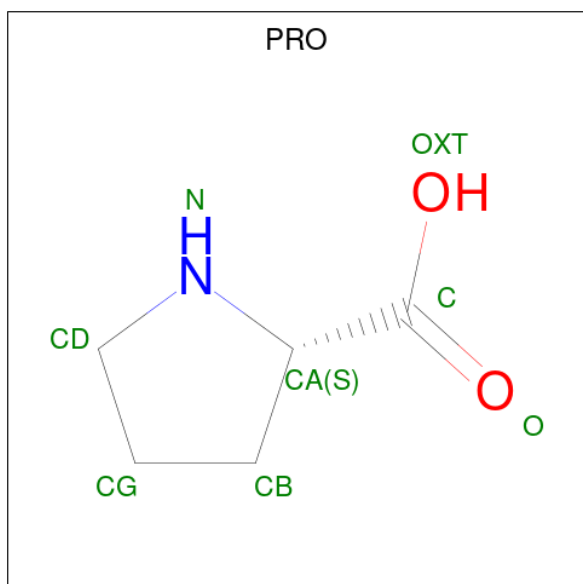
There are 5 unique types of molecules in this entry. The entry contains 3858 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 AMINOPEPTIDASE P.

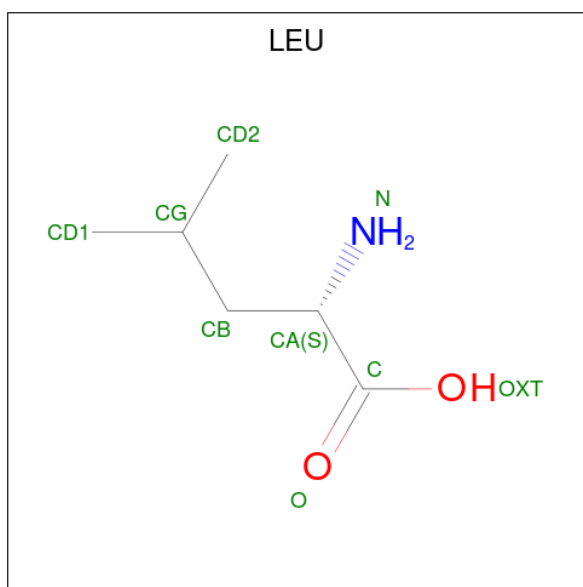
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	439	3494	2199	617	663	15	3	0	0

- Molecule 2 is PROLINE (CCD ID: PRO) (formula: $C_5H_9NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	7	5	1	1	0	0

- Molecule 3 is LEUCINE (CCD ID: LEU) (formula: $C_6H_{13}NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			9	6	1	2		

- Molecule 4 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mn	0	0
			2	2		

- Molecule 5 is water.

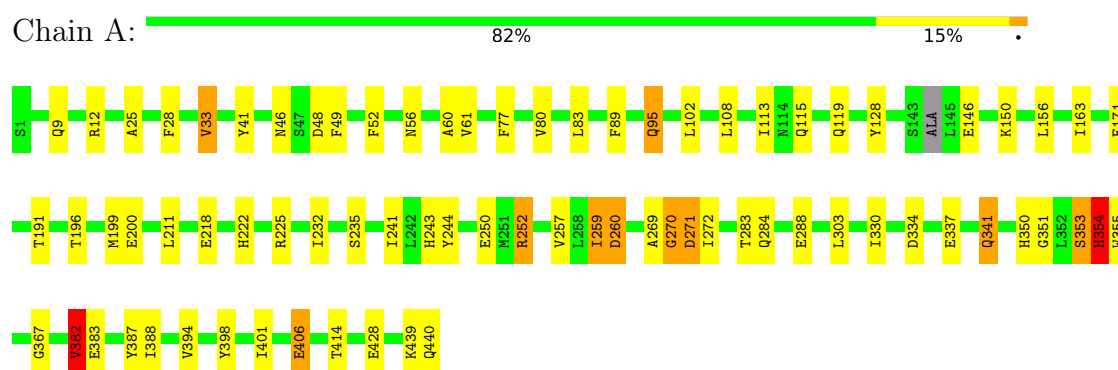
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	346	Total	O	0	0
			346	346		

3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

• Molecule 1: AMINOPEPTIDASE P



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants a, b, c, α , β , γ	177.30Å 177.30Å 96.50Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.30	Depositor
% Data completeness (in resolution range)	91.9 (50.00-2.30)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.184 , 0.226	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3858	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

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: MN

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	1.34	32/3565 (0.9%)	1.25	43/4825 (0.9%)

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	A	0	2

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	354	HIS	CA-C	28.66	1.87	1.52
1	A	353	SER	C-N	24.20	1.63	1.33
1	A	383	GLU	CA-CB	-21.95	1.24	1.53
1	A	383	GLU	C-O	-19.34	1.01	1.24
1	A	354	HIS	CG-CD2	-15.57	1.18	1.35
1	A	383	GLU	C-N	14.97	1.52	1.33
1	A	354	HIS	CD2-NE2	14.65	1.53	1.37
1	A	271	ASP	N-CA	-13.73	1.28	1.47
1	A	383	GLU	N-CA	12.48	1.60	1.46
1	A	383	GLU	CD-OE1	12.47	1.49	1.25
1	A	354	HIS	CB-CG	11.99	1.67	1.50
1	A	354	HIS	C-N	11.94	1.49	1.33
1	A	383	GLU	CA-C	11.62	1.65	1.52
1	A	260	ASP	C-O	11.57	1.37	1.24
1	A	260	ASP	N-CA	-10.12	1.33	1.46
1	A	354	HIS	CA-CB	9.98	1.71	1.53
1	A	260	ASP	CA-C	-9.81	1.40	1.53
1	A	271	ASP	C-N	-9.74	1.19	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	406	GLU	CD-OE2	9.50	1.43	1.25
1	A	382	VAL	C-N	8.96	1.44	1.33
1	A	259	ILE	C-N	8.59	1.46	1.33
1	A	271	ASP	CG-OD1	8.46	1.41	1.25
1	A	260	ASP	CA-CB	-7.90	1.39	1.53
1	A	270	GLY	C-N	7.43	1.43	1.33
1	A	260	ASP	CG-OD1	7.24	1.39	1.25
1	A	354	HIS	ND1-CE1	6.95	1.39	1.32
1	A	406	GLU	CD-OE1	6.74	1.38	1.25
1	A	383	GLU	CB-CG	6.28	1.71	1.52
1	A	406	GLU	C-O	6.01	1.30	1.24
1	A	271	ASP	CA-C	-5.83	1.48	1.53
1	A	199	MET	SD-CE	-5.70	1.65	1.79
1	A	406	GLU	CG-CD	-5.10	1.39	1.52

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	354	HIS	O-C-N	20.69	146.83	123.25
1	A	382	VAL	CA-C-N	-15.25	108.19	122.37
1	A	382	VAL	C-N-CA	-15.25	108.19	122.37
1	A	353	SER	CA-C-N	-13.41	98.55	121.75
1	A	353	SER	C-N-CA	-13.41	98.55	121.75
1	A	354	HIS	N-CA-C	-12.38	90.25	109.07
1	A	270	GLY	O-C-N	-11.54	99.42	123.07
1	A	271	ASP	CB-CA-C	10.77	126.37	111.51
1	A	383	GLU	CG-CD-OE2	9.85	141.05	118.40
1	A	354	HIS	CA-C-O	-9.39	111.19	121.33
1	A	260	ASP	CB-CA-C	9.10	126.52	111.23
1	A	259	ILE	CA-C-N	8.12	135.40	123.13
1	A	259	ILE	C-N-CA	8.12	135.40	123.13
1	A	259	ILE	O-C-N	-7.87	113.89	123.10
1	A	271	ASP	O-C-N	-7.65	116.46	123.50
1	A	383	GLU	OE1-CD-OE2	-7.13	105.79	122.90
1	A	354	HIS	CB-CG-ND1	-6.95	112.27	122.70
1	A	367	GLY	N-CA-C	-6.82	103.79	112.65
1	A	353	SER	O-C-N	6.71	131.29	123.17
1	A	83	LEU	N-CA-C	6.45	118.31	111.28
1	A	414	THR	N-CA-C	6.37	120.69	112.41
1	A	354	HIS	CA-C-N	-6.29	110.23	120.87
1	A	354	HIS	C-N-CA	-6.29	110.23	120.87
1	A	383	GLU	O-C-N	6.27	125.76	121.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	383	GLU	CA-CB-CG	6.18	126.46	114.10
1	A	269	ALA	N-CA-C	6.14	119.65	109.46
1	A	270	GLY	N-CA-C	-6.10	103.33	113.02
1	A	271	ASP	CA-CB-CG	6.04	118.64	112.60
1	A	383	GLU	CA-C-N	-5.94	113.57	119.92
1	A	383	GLU	C-N-CA	-5.94	113.57	119.92
1	A	113	ILE	N-CA-C	5.90	117.35	110.62
1	A	95	GLN	N-CA-C	5.88	118.92	111.69
1	A	354	HIS	CE1-NE2-CD2	-5.78	103.22	109.00
1	A	225	ARG	N-CA-C	5.73	117.52	111.28
1	A	89	PHE	N-CA-C	5.64	121.06	113.72
1	A	383	GLU	CB-CA-C	-5.48	104.35	111.15
1	A	303	LEU	N-CA-C	5.45	118.39	111.69
1	A	283	THR	N-CA-C	-5.40	102.76	110.59
1	A	260	ASP	OD1-CG-OD2	-5.34	110.09	122.90
1	A	56	ASN	N-CA-C	5.18	118.43	112.57
1	A	171	GLU	N-CA-C	-5.12	105.88	111.82
1	A	260	ASP	CA-C-O	-5.04	114.22	120.57
1	A	252	ARG	N-CA-C	5.03	118.41	109.96

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	353	SER	Mainchain
1	A	382	VAL	Mainchain

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	3494	0	3423	40	0
2	A	7	0	7	2	0
3	A	9	0	11	0	0
4	A	2	0	0	0	0
5	A	346	0	0	0	0
All	All	3858	0	3441	40	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 (40) 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:A:354:HIS:CA	1:A:354:HIS:C	1.87	1.44
1:A:46:ASN:HD22	1:A:49:PHE:H	1.21	0.88
1:A:354:HIS:C	1:A:354:HIS:N	2.38	0.82
1:A:250:GLU:HB3	1:A:252:ARG:HH21	1.55	0.71
1:A:350:HIS:HE1	1:A:387:TYR:OH	1.80	0.65
1:A:354:HIS:CA	1:A:355:TRP:N	2.60	0.65
1:A:354:HIS:N	1:A:354:HIS:O	2.35	0.60
1:A:284:GLN:O	1:A:288:GLU:HG3	2.03	0.58
1:A:241:ILE:HB	1:A:244:TYR:HB2	1.89	0.54
1:A:146:GLU:OE1	1:A:150:LYS:HE2	2.10	0.52
1:A:243:HIS:CD2	2:A:441:PRO:HD2	2.46	0.51
1:A:77:PHE:HA	1:A:108:LEU:O	2.11	0.51
1:A:46:ASN:ND2	1:A:49:PHE:H	2.01	0.49
1:A:211:LEU:HD13	1:A:259:ILE:HD11	1.96	0.48
1:A:350:HIS:HD2	1:A:351:GLY:O	1.97	0.48
1:A:60:ALA:HA	1:A:77:PHE:O	2.14	0.48
1:A:191:THR:HG22	1:A:272:ILE:HD12	1.96	0.48
1:A:270:GLY:O	1:A:271:ASP:HB2	2.15	0.47
1:A:337:GLU:O	1:A:341:GLN:HG2	2.15	0.46
1:A:128:TYR:HA	1:A:163:ILE:O	2.15	0.46
1:A:334:ASP:HB3	1:A:337:GLU:HB2	1.99	0.45
1:A:28:PHE:CD2	1:A:61:VAL:HG22	2.52	0.45
1:A:80:VAL:HG23	1:A:95:GLN:HE22	1.82	0.45
1:A:354:HIS:C	1:A:354:HIS:CB	2.84	0.45
1:A:388:ILE:HG21	1:A:394:VAL:HG21	1.99	0.45
1:A:115:GLN:O	1:A:119:GLN:NE2	2.48	0.44
1:A:330:ILE:HG23	1:A:394:VAL:HG11	1.98	0.44
1:A:398:TYR:HA	1:A:401:ILE:HD12	2.00	0.44
1:A:33:VAL:HG12	1:A:41:TYR:HD2	1.83	0.44
1:A:218:GLU:OE2	1:A:222:HIS:HE1	2.01	0.44
1:A:243:HIS:NE2	2:A:441:PRO:HD2	2.34	0.43
1:A:232:ILE:HB	1:A:260:ASP:HB3	2.00	0.43
1:A:46:ASN:HD21	1:A:48:ASP:HB2	1.82	0.43
1:A:439:LYS:O	1:A:440:GLN:HB2	2.19	0.43
1:A:196:THR:O	1:A:200:GLU:HG3	2.19	0.42
1:A:382:VAL:O	1:A:406:GLU:HA	2.20	0.41
1:A:12:ARG:NH1	1:A:52:PHE:O	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:350:HIS:CE1	1:A:387:TYR:OH	2.68	0.41
1:A:235:SER:HA	1:A:257:VAL:HA	2.03	0.41
1:A:25:ALA:HA	1:A:128:TYR:HB2	2.03	0.41

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	435/440 (99%)	425 (98%)	10 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	371/371 (100%)	364 (98%)	7 (2%)	50	69

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	33	VAL
1	A	102	LEU

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Mol	Chain	Res	Type
1	A	156	LEU
1	A	341	GLN
1	A	354	HIS
1	A	428	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	9	GLN
1	A	46	ASN
1	A	56	ASN
1	A	95	GLN
1	A	116	GLN
1	A	222	HIS
1	A	230	ASN
1	A	284	GLN
1	A	341	GLN
1	A	350	HIS
1	A	368	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	PRO	A	441	3	6,7,8	0.50	0	7,8,10	1.02	0
3	LEU	A	442	2	6,8,8	1.18	1 (16%)	5,10,10	1.65	1 (20%)

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	PRO	A	441	3	-	0/0/9/11	0/1/1/1
3	LEU	A	442	2	-	2/8/8/8	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	442	LEU	OXT-C	-2.25	1.23	1.30

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	442	LEU	OXT-C-O	-3.42	116.32	124.08

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	442	LEU	OXT-C-CA-N
3	A	442	LEU	C-CA-CB-CG

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	441	PRO	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	353:SER	C	354:HIS	N	1.63
1	A	271:ASP	C	272:ILE	N	1.19

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.