



# Full wwPDB X-ray Structure Validation Report ⓘ

May 3, 2016 – 01:42 PM EDT

PDB ID : 5D8N  
Title : Tomato leucine aminopeptidase mutant - K354E  
Authors : DuPrez, K.T.; Scranton, M.A.; Walling, L.L.; Fan, L.  
Deposited on : 2015-08-17  
Resolution : 2.15 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

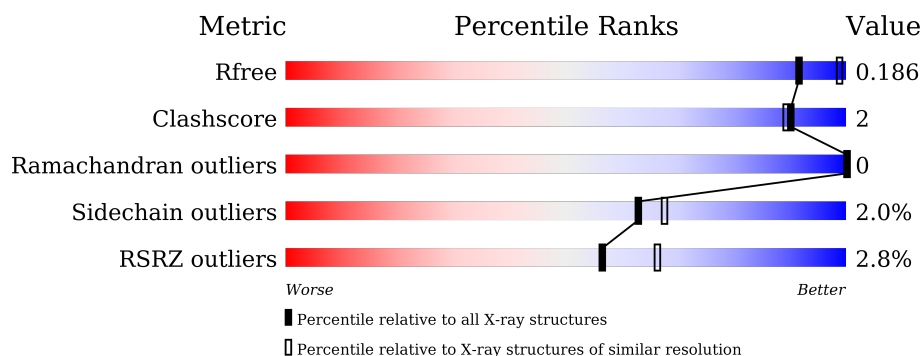
MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20027457  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20027457



**i**

## X-RAY DIFFRACTION

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	1045 (2.16-2.16)
Clashscore	102246	1152 (2.16-2.16)
Ramachandran outliers	100387	1131 (2.16-2.16)
Sidechain outliers	100360	1131 (2.16-2.16)
RSRZ outliers	91569	1050 (2.16-2.16)

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

Mol	Chain	Length	Quality of chain
1	A	529	<div> <div>4%</div> <div>90%</div> <div>5%</div> <div>5%</div> </div>
1	B	529	<div> <div>2%</div> <div>91%</div> <div>•</div> <div>5%</div> </div>
1	C	529	<div> <div>2%</div> <div>89%</div> <div>5%</div> <div>•</div> <div>5%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:



Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	MG	B	609	-	-	-	X
4	ACT	A	606	-	-	-	X
4	ACT	C	608	-	-	-	X
5	EDO	B	602	-	-	-	X
5	EDO	B	603	-	-	-	X
5	EDO	C	603	-	-	-	X



## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 12858 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 Leucine aminopeptidase 1, chloroplastic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	505	Total	C	N	O	S	0	19	0
			3847	2393	660	778	16			
1	B	504	Total	C	N	O	S	0	13	0
			3791	2366	648	761	16			
1	C	504	Total	C	N	O	S	0	25	0
			3889	2419	667	787	16			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	43	MET	-	initiating methionine	UNP Q10712
A	44	ARG	-	expression tag	UNP Q10712
A	45	GLY	-	expression tag	UNP Q10712
A	46	SER	-	expression tag	UNP Q10712
A	47	HIS	-	expression tag	UNP Q10712
A	48	HIS	-	expression tag	UNP Q10712
A	49	HIS	-	expression tag	UNP Q10712
A	50	HIS	-	expression tag	UNP Q10712
A	51	HIS	-	expression tag	UNP Q10712
A	52	HIS	-	expression tag	UNP Q10712
A	53	GLY	-	expression tag	UNP Q10712
A	354	GLU	LYS	engineered mutation	UNP Q10712
B	43	MET	-	initiating methionine	UNP Q10712
B	44	ARG	-	expression tag	UNP Q10712
B	45	GLY	-	expression tag	UNP Q10712
B	46	SER	-	expression tag	UNP Q10712
B	47	HIS	-	expression tag	UNP Q10712
B	48	HIS	-	expression tag	UNP Q10712
B	49	HIS	-	expression tag	UNP Q10712
B	50	HIS	-	expression tag	UNP Q10712
B	51	HIS	-	expression tag	UNP Q10712
B	52	HIS	-	expression tag	UNP Q10712
B	53	GLY	-	expression tag	UNP Q10712

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Chain	Residue	Modelled	Actual	Comment	Reference
B	354	GLU	LYS	engineered mutation	UNP Q10712
C	43	MET	-	initiating methionine	UNP Q10712
C	44	ARG	-	expression tag	UNP Q10712
C	45	GLY	-	expression tag	UNP Q10712
C	46	SER	-	expression tag	UNP Q10712
C	47	HIS	-	expression tag	UNP Q10712
C	48	HIS	-	expression tag	UNP Q10712
C	49	HIS	-	expression tag	UNP Q10712
C	50	HIS	-	expression tag	UNP Q10712
C	51	HIS	-	expression tag	UNP Q10712
C	52	HIS	-	expression tag	UNP Q10712
C	53	GLY	-	expression tag	UNP Q10712
C	354	GLU	LYS	engineered mutation	UNP Q10712

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

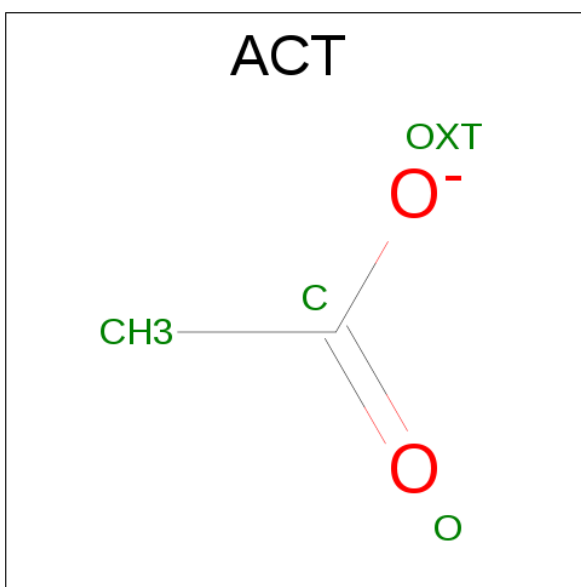
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	3	Total Cl 3 3	0	0
2	A	3	Total Cl 3 3	0	0
2	C	2	Total Cl 2 2	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	3	Total Mg 3 3	0	0
3	A	2	Total Mg 2 2	0	0
3	C	2	Total Mg 2 2	0	0

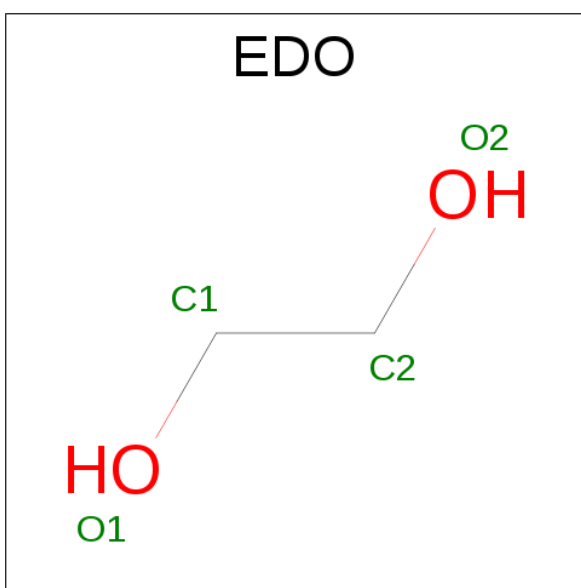
- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: C<sub>2</sub>H<sub>3</sub>O<sub>2</sub>).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	C	1	Total 4	C 2	O 2	0	0
5	C	1	Total 4	C 2	O 2	0	0
5	C	1	Total 4	C 2	O 2	0	0

- Molecule 6 is water.

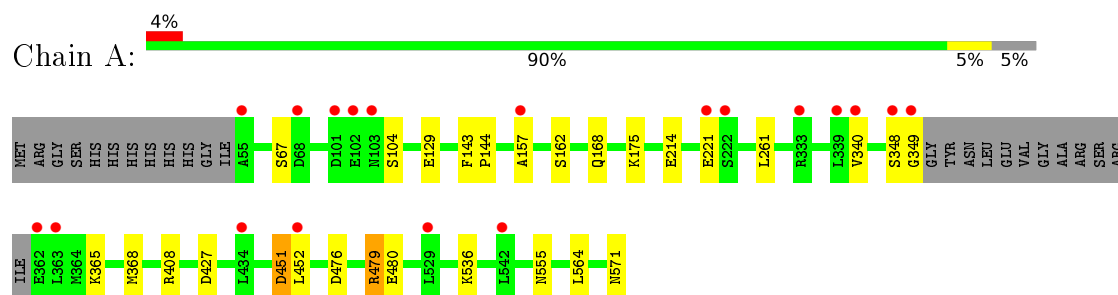
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	402	Total 402	O 402	0	0
6	B	461	Total 461	O 461	0	0
6	C	417	Total 417	O 417	0	0



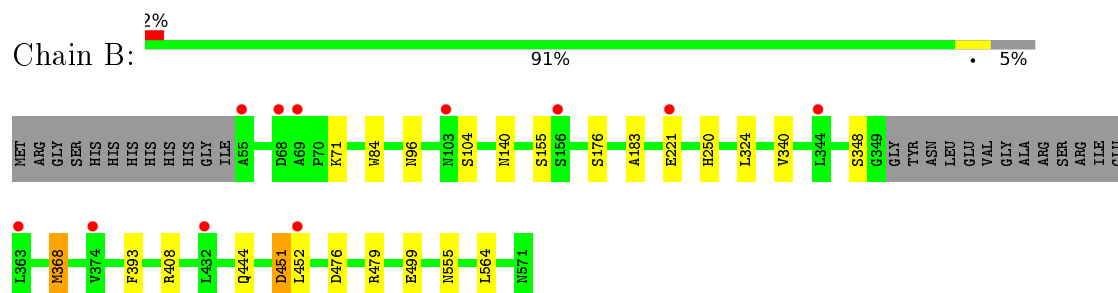
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

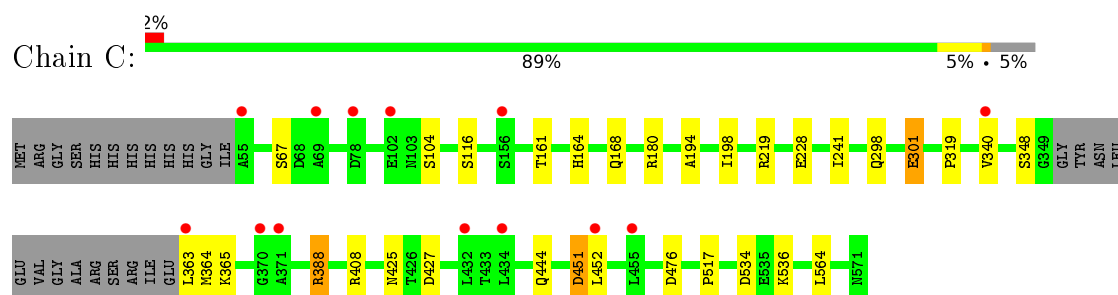
- Molecule 1: Leucine aminopeptidase 1, chloroplastic



- Molecule 1: Leucine aminopeptidase 1, chloroplastic



- Molecule 1: Leucine aminopeptidase 1, chloroplastic





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	127.11Å 165.19Å 168.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.15 29.89 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.9 (30.00-2.15) 100.0 (29.89-2.15)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.79 (at 2.16Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.153 , 0.179 0.162 , 0.186	Depositor DCC
$R_{free}$ test set	2000 reflections (2.13%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.4	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 39.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	12858	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.55% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: MG, CL, EDO, ACT

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	0.40	0/3900	0.68	0/5284
1	B	0.42	0/3844	0.67	1/5211 (0.0%)
1	C	0.41	0/3942	0.70	3/5339 (0.1%)
All	All	0.41	0/11686	0.68	4/15834 (0.0%)

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
1	B	0	1
1	C	0	1
All	All	0	4

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	C	219	ARG	NE-CZ-NH2	-5.81	117.39	120.30
1	B	368	MET	CB-CG-SD	-5.32	96.44	112.40
1	C	534	ASP	CB-CG-OD2	-5.22	113.60	118.30
1	C	476	ASP	CB-CG-OD2	-5.13	113.68	118.30

There are no chirality outliers.

All (4) planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	A	408	ARG	Sidechain
1	A	479[A]	ARG	Sidechain
1	B	408	ARG	Sidechain
1	C	408	ARG	Sidechain

## 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	3847	0	3799	14	0
1	B	3791	0	3749	12	0
1	C	3889	0	3841	25	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
2	C	2	0	0	0	0
3	A	2	0	0	0	0
3	B	3	0	0	0	0
3	C	2	0	0	0	0
4	A	4	0	3	0	0
4	C	4	0	3	1	0
5	B	16	0	24	2	0
5	C	12	0	18	2	0
6	A	402	0	0	4	0
6	B	461	0	0	6	0
6	C	417	0	0	13	0
All	All	12858	0	11437	53	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 2.

All (53) 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:96[B]:ASN:ND2	6:B:701:HOH:O	1.70	1.14
1:C:180:ARG:NH1	1:C:228:GLU:OE1	1.90	1.04
1:B:499:GLU:OE1	6:B:702:HOH:O	1.78	0.99
1:A:476[A]:ASP:OD1	1:A:479[A]:ARG:NH2	2.01	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:298:GLN:O	1:C:301[A]:GLU:HG3	1.77	0.84
4:C:608:ACT:H1	6:C:871:HOH:O	1.82	0.79
1:B:476[B]:ASP:OD1	1:B:479[B]:ARG:NH1	2.18	0.76
1:C:388[A]:ARG:HG2	6:C:1024:HOH:O	1.88	0.72
1:C:180:ARG:NH1	1:C:228:GLU:CD	2.43	0.71
5:C:601:EDO:O1	6:C:701:HOH:O	2.12	0.67
1:C:161[B]:THR:HG22	6:C:833:HOH:O	1.95	0.67
1:B:444:GLN:OE1	6:B:704:HOH:O	2.14	0.66
1:C:444:GLN:OE1	6:C:702:HOH:O	2.14	0.65
1:C:161[B]:THR:HG21	6:C:933:HOH:O	1.95	0.64
1:C:198:ILE:HD12	1:C:241:ILE:HB	1.87	0.56
1:C:364:MET:CE	6:C:1021:HOH:O	2.54	0.55
1:C:164:HIS:O	1:C:168[B]:GLN:HG2	2.07	0.55
1:C:180:ARG:HD2	6:C:889:HOH:O	2.07	0.54
1:A:157:ALA:HB1	1:A:162:SER:OG	2.08	0.53
1:B:140[B]:ASN:ND2	6:B:709:HOH:O	2.40	0.53
1:A:365:LYS:HA	6:A:843:HOH:O	2.07	0.53
1:C:364:MET:HE1	6:C:1021:HOH:O	2.09	0.51
1:C:168[A]:GLN:NE2	6:C:716:HOH:O	2.43	0.50
1:A:571:ASN:OXT	6:A:701:HOH:O	2.19	0.49
1:C:194:ALA:O	1:C:198:ILE:HG12	2.13	0.49
1:A:129:GLU:HB3	5:B:604:EDO:H22	1.93	0.48
1:A:175:LYS:HE3	1:A:214[B]:GLU:OE2	2.14	0.47
1:A:261:LEU:HB3	1:A:368:MET:HE1	1.96	0.47
1:B:250:HIS:HD2	6:B:1083:HOH:O	1.97	0.47
1:C:198:ILE:HD12	1:C:241:ILE:CB	2.45	0.46
1:C:319:PRO:HD3	6:C:753:HOH:O	2.15	0.46
1:A:168[B]:GLN:NE2	6:A:711:HOH:O	2.48	0.45
1:B:155:SER:HA	6:B:756:HOH:O	2.17	0.45
1:B:452:LEU:HD11	1:B:564:LEU:HD21	1.98	0.45
1:A:452:LEU:HD11	1:A:564:LEU:HD21	1.99	0.45
1:C:198:ILE:CD1	1:C:241:ILE:HB	2.47	0.45
1:B:176:SER:HB2	5:B:604:EDO:H21	1.99	0.44
1:C:425:ASN:OD1	1:C:427:ASP:HB2	2.18	0.44
1:C:180:ARG:HB3	1:C:180:ARG:NH1	2.33	0.44
1:B:340:VAL:O	1:B:451:ASP:HA	2.17	0.43
1:B:84:TRP:CE2	1:B:183:ALA:HB2	2.53	0.43
1:C:517:PRO:HG2	5:C:603:EDO:H11	1.99	0.43
1:A:340:VAL:O	1:A:451:ASP:HA	2.19	0.43
1:C:452:LEU:HD11	1:C:564:LEU:HD21	2.00	0.43
1:C:340:VAL:O	1:C:451:ASP:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:364:MET:HE2	6:C:798:HOH:O	2.20	0.42
1:A:349:GLY:HA2	1:A:427:ASP:OD1	2.20	0.42
1:A:157:ALA:CB	1:A:162:SER:OG	2.67	0.41
1:B:324:LEU:HB2	1:B:393:PHE:HB2	2.02	0.41
1:A:143:PHE:HB2	1:A:144:PRO:HD2	2.01	0.41
1:C:365:LYS:HG2	6:C:804:HOH:O	2.21	0.41
1:C:180:ARG:HH11	1:C:228:GLU:CD	2.23	0.40
1:A:480:GLU:OE2	6:A:702:HOH:O	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	520/529 (98%)	510 (98%)	10 (2%)	0	100	100
1	B	513/529 (97%)	507 (99%)	6 (1%)	0	100	100
1	C	525/529 (99%)	518 (99%)	7 (1%)	0	100	100
All	All	1558/1587 (98%)	1535 (98%)	23 (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	403/409 (98%)	395 (98%)	8 (2%)	63	67
1	B	393/409 (96%)	384 (98%)	9 (2%)	58	62
1	C	408/409 (100%)	396 (97%)	12 (3%)	50	49
All	All	1204/1227 (98%)	1175 (98%)	29 (2%)	63	60

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	SER
1	A	104	SER
1	A	221[A]	GLU
1	A	221[B]	GLU
1	A	348	SER
1	A	451	ASP
1	A	536	LYS
1	A	555	ASN
1	B	71	LYS
1	B	104[A]	SER
1	B	104[B]	SER
1	B	221[A]	GLU
1	B	221[B]	GLU
1	B	348	SER
1	B	368	MET
1	B	451	ASP
1	B	555	ASN
1	C	67	SER
1	C	104	SER
1	C	116[A]	SER
1	C	116[B]	SER
1	C	301[A]	GLU
1	C	301[B]	GLU
1	C	348	SER
1	C	363	LEU
1	C	388[A]	ARG
1	C	388[B]	ARG
1	C	451	ASP
1	C	536	LYS

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



Mol	Chain	Res	Type
1	B	178	GLN
1	C	250	HIS
1	C	538	GLN
1	C	555	ASN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 24 ligands modelled in this entry, 15 are monoatomic - leaving 9 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	$\# Z  > 2$
4	ACT	A	606	-	0,3,3	0.00	-	0,3,3	0.00	-
5	EDO	B	601	-	3,3,3	0.42	0	2,2,2	0.63	0
5	EDO	B	602	-	3,3,3	0.38	0	2,2,2	0.73	0
5	EDO	B	603	-	3,3,3	0.44	0	2,2,2	0.40	0
5	EDO	B	604	-	3,3,3	0.49	0	2,2,2	0.26	0
5	EDO	C	601	-	3,3,3	0.43	0	2,2,2	0.73	0
5	EDO	C	602	-	3,3,3	0.40	0	2,2,2	0.50	0
5	EDO	C	603	-	3,3,3	0.45	0	2,2,2	0.34	0
4	ACT	C	608	-	0,3,3	0.00	-	0,3,3	0.00	-

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
4	ACT	A	606	-	-	0/0/0/0	0/0/0/0
5	EDO	B	601	-	-	0/1/1/1	0/0/0/0
5	EDO	B	602	-	-	0/1/1/1	0/0/0/0
5	EDO	B	603	-	-	0/1/1/1	0/0/0/0
5	EDO	B	604	-	-	0/1/1/1	0/0/0/0
5	EDO	C	601	-	-	0/1/1/1	0/0/0/0
5	EDO	C	602	-	-	0/1/1/1	0/0/0/0
5	EDO	C	603	-	-	0/1/1/1	0/0/0/0
4	ACT	C	608	-	-	0/0/0/0	0/0/0/0

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.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	604	EDO	2	0
5	C	601	EDO	1	0
5	C	603	EDO	1	0
4	C	608	ACT	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, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	505/529 (95%)	-0.10	19 (3%) 44 54	17, 26, 64, 147	0
1	B	504/529 (95%)	-0.25	11 (2%) 65 73	16, 24, 55, 114	1 (0%)
1	C	504/529 (95%)	-0.18	13 (2%) 59 68	17, 27, 56, 102	0
All	All	1513/1587 (95%)	-0.18	43 (2%) 56 66	16, 26, 58, 147	1 (0%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	363	LEU	6.2
1	A	68	ASP	4.8
1	B	363	LEU	4.3
1	B	103	ASN	3.8
1	B	68	ASP	3.6
1	B	55	ALA	3.5
1	A	103	ASN	3.4
1	A	362	GLU	3.4
1	A	102[A]	GLU	3.3
1	A	452	LEU	3.2
1	A	434	LEU	3.1
1	A	55	ALA	3.1
1	C	432[A]	LEU	3.0
1	B	432[A]	LEU	3.0
1	C	434	LEU	2.9
1	C	340	VAL	2.8
1	A	529	LEU	2.8
1	C	102	GLU	2.7
1	C	156	SER	2.7
1	A	349	GLY	2.7
1	A	542	LEU	2.7
1	A	340	VAL	2.6
1	C	78	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	333	ARG	2.5
1	B	221[A]	GLU	2.4
1	A	221[A]	GLU	2.4
1	A	101	ASP	2.3
1	A	222	SER	2.2
1	C	371	ALA	2.2
1	B	374	VAL	2.2
1	B	344	LEU	2.2
1	B	156	SER	2.2
1	C	69	ALA	2.2
1	C	455	LEU	2.1
1	C	370	GLY	2.1
1	C	55	ALA	2.1
1	C	363	LEU	2.1
1	C	452	LEU	2.1
1	A	339	LEU	2.0
1	B	452	LEU	2.0
1	B	69	ALA	2.0
1	A	157	ALA	2.0
1	A	348	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, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
5	EDO	B	602	4/4	0.87	0.18	10.37	50,53,54,58	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
5	EDO	C	603	4/4	0.91	0.22	9.29	54,54,62,62	0
3	MG	B	609	1/1	0.90	0.16	6.57	49,49,49,49	0
4	ACT	A	606	4/4	0.88	0.22	6.43	38,42,50,51	0
4	ACT	C	608	4/4	0.91	0.20	5.53	40,43,44,46	0
5	EDO	B	603	4/4	0.92	0.32	4.87	47,58,64,68	0
3	MG	C	606	1/1	0.90	0.10	1.41	47,47,47,47	0
3	MG	A	604	1/1	0.97	0.12	0.52	47,47,47,47	0
5	EDO	C	602	4/4	0.94	0.15	0.27	40,42,51,60	0
2	CL	B	605	1/1	0.99	0.15	-0.49	32,32,32,32	0
2	CL	B	606	1/1	0.99	0.12	-0.89	34,34,34,34	0
3	MG	B	608	1/1	0.92	0.07	-0.90	37,37,37,37	0
2	CL	A	603	1/1	0.95	0.09	-0.95	34,34,34,34	0
2	CL	A	601	1/1	1.00	0.12	-0.98	31,31,31,31	0
2	CL	C	605	1/1	0.98	0.12	-1.02	38,38,38,38	0
3	MG	A	605	1/1	0.90	0.08	-1.11	50,50,50,50	0
2	CL	C	604	1/1	0.99	0.10	-1.54	32,32,32,32	0
3	MG	C	607	1/1	0.95	0.05	-1.87	43,43,43,43	0
3	MG	B	610	1/1	0.94	0.18	-	47,47,47,47	0
5	EDO	B	604	4/4	0.82	0.15	-	58,64,66,68	0
5	EDO	B	601	4/4	0.92	0.21	-	51,52,66,69	0
2	CL	B	607	1/1	0.95	0.09	-	59,59,59,59	0
5	EDO	C	601	4/4	0.87	0.24	-	37,42,44,44	0
2	CL	A	602	1/1	0.94	0.09	-	54,54,54,54	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.