



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 03:16 am GMT

PDB ID : 3PYB  
Title : Crystal structure of ent-copalyl diphosphate synthase from *Arabidopsis thaliana* in complex with 13-aza-13,14-dihydrocopalyl diphosphate  
Authors : Koksai, M.; Christianson, D.W.  
Deposited on : 2010-12-12  
Resolution : 2.76 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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

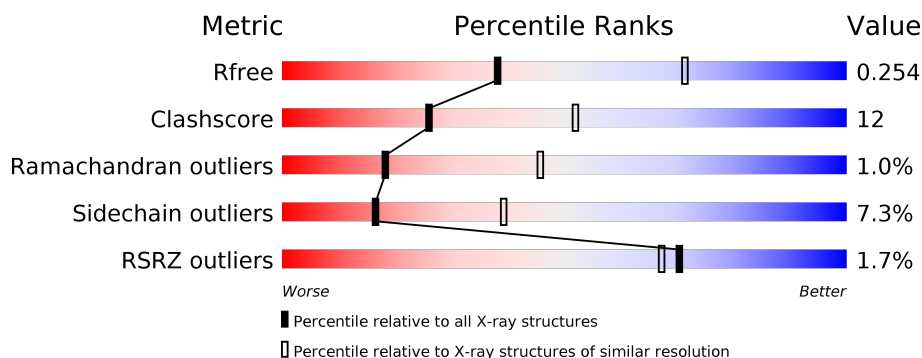
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.76 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	100719	3666 (2.80-2.72)
Clashscore	112137	4174 (2.80-2.72)
Ramachandran outliers	110173	4103 (2.80-2.72)
Sidechain outliers	110143	4106 (2.80-2.72)
RSRZ outliers	101464	3697 (2.80-2.72)

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	727	<div> <div style="width: 3%;"></div> <div style="width: 68%; background-color: green;"></div> <div style="width: 24%; background-color: yellow;"></div> <div style="width: 6%; background-color: orange;"></div> <div style="width: 1%; background-color: red;"></div> <div style="width: 1%; background-color: grey;"></div> </div> <div>68% 24% • 6%</div>
1	B	727	<div> <div style="width: 3%;"></div> <div style="width: 66%; background-color: green;"></div> <div style="width: 26%; background-color: yellow;"></div> <div style="width: 5%; background-color: orange;"></div> <div style="width: 1%; background-color: red;"></div> <div style="width: 1%; background-color: grey;"></div> </div> <div>66% 26% • 5%</div>
1	C	727	<div> <div style="width: 3%;"></div> <div style="width: 67%; background-color: green;"></div> <div style="width: 24%; background-color: yellow;"></div> <div style="width: 6%; background-color: orange;"></div> <div style="width: 1%; background-color: red;"></div> <div style="width: 1%; background-color: grey;"></div> </div> <div>67% 24% • 6%</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	1PE	A	921	-	-	-	X
3	1PE	A	931	-	-	-	X
3	1PE	B	921	-	-	-	X
3	1PE	B	931	-	-	-	X
3	1PE	C	911	-	-	-	X
3	1PE	C	931	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 17263 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 Ent-copalyl diphosphate synthase, chloroplastic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	685	Total	C	N	O	S	0	4	0
			5633	3618	948	1038	29			
1	B	689	Total	C	N	O	S	0	1	0
			5649	3622	952	1046	29			
1	C	685	Total	C	N	O	S	0	0	0
			5607	3598	943	1037	29			

There are 27 discrepancies between the modelled and reference sequences:

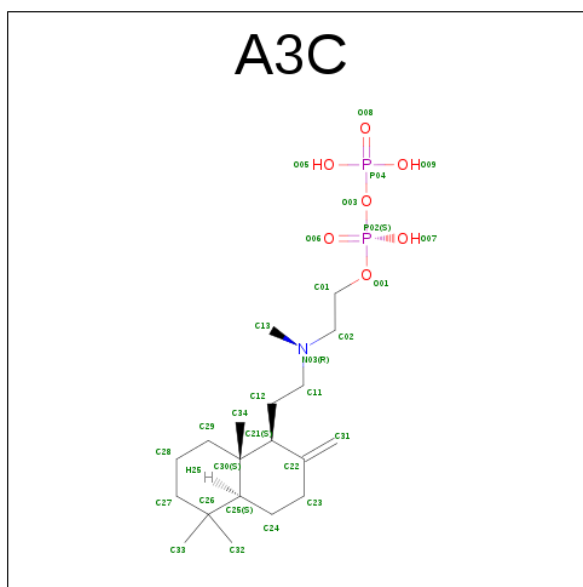
Chain	Residue	Modelled	Actual	Comment	Reference
A	84	MET	-	EXPRESSION TAG	UNP Q38802
A	803	GLY	-	EXPRESSION TAG	UNP Q38802
A	804	SER	-	EXPRESSION TAG	UNP Q38802
A	805	HIS	-	EXPRESSION TAG	UNP Q38802
A	806	HIS	-	EXPRESSION TAG	UNP Q38802
A	807	HIS	-	EXPRESSION TAG	UNP Q38802
A	808	HIS	-	EXPRESSION TAG	UNP Q38802
A	809	HIS	-	EXPRESSION TAG	UNP Q38802
A	810	HIS	-	EXPRESSION TAG	UNP Q38802
B	84	MET	-	EXPRESSION TAG	UNP Q38802
B	803	GLY	-	EXPRESSION TAG	UNP Q38802
B	804	SER	-	EXPRESSION TAG	UNP Q38802
B	805	HIS	-	EXPRESSION TAG	UNP Q38802
B	806	HIS	-	EXPRESSION TAG	UNP Q38802
B	807	HIS	-	EXPRESSION TAG	UNP Q38802
B	808	HIS	-	EXPRESSION TAG	UNP Q38802
B	809	HIS	-	EXPRESSION TAG	UNP Q38802
B	810	HIS	-	EXPRESSION TAG	UNP Q38802
C	84	MET	-	EXPRESSION TAG	UNP Q38802
C	803	GLY	-	EXPRESSION TAG	UNP Q38802
C	804	SER	-	EXPRESSION TAG	UNP Q38802
C	805	HIS	-	EXPRESSION TAG	UNP Q38802
C	806	HIS	-	EXPRESSION TAG	UNP Q38802

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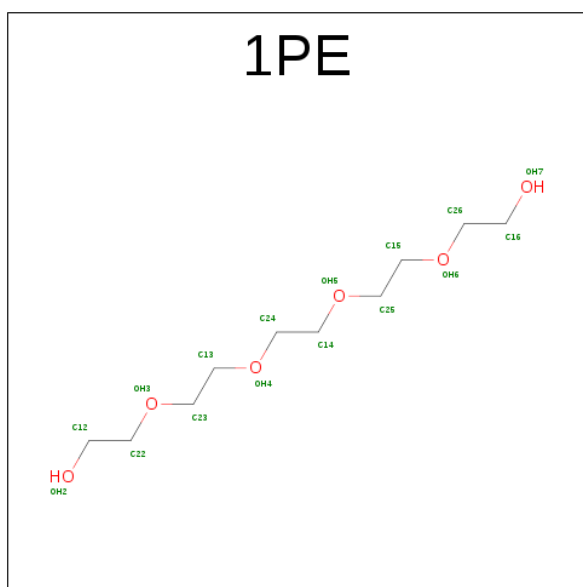
Chain	Residue	Modelled	Actual	Comment	Reference
C	807	HIS	-	EXPRESSION TAG	UNP Q38802
C	808	HIS	-	EXPRESSION TAG	UNP Q38802
C	809	HIS	-	EXPRESSION TAG	UNP Q38802
C	810	HIS	-	EXPRESSION TAG	UNP Q38802

- Molecule 2 is 2-(METHYL{2-[(1S,4AS,8AS)-5,5,8A-TRIMETHYL-2-METHYLIDE NEDECAHYDRONAPHTHALEN-1-YL]ETHYL}AMINO)ETHYL TRIHYDROGEN DIPHOSPHATE (three-letter code: A3C) (formula: C<sub>19</sub>H<sub>37</sub>NO<sub>7</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			29	19	1	7	2		

- Molecule 3 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: C<sub>10</sub>H<sub>22</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			16	10	6		
3	A	1	Total	C	O	0	0
			16	10	6		
3	A	1	Total	C	O	0	0
			16	10	6		
3	A	1	Total	C	O	0	0
			16	10	6		
3	B	1	Total	C	O	0	0
			16	10	6		
3	B	1	Total	C	O	0	0
			16	10	6		
3	B	1	Total	C	O	0	0
			16	10	6		
3	C	1	Total	C	O	0	0
			16	10	6		
3	C	1	Total	C	O	0	0
			16	10	6		

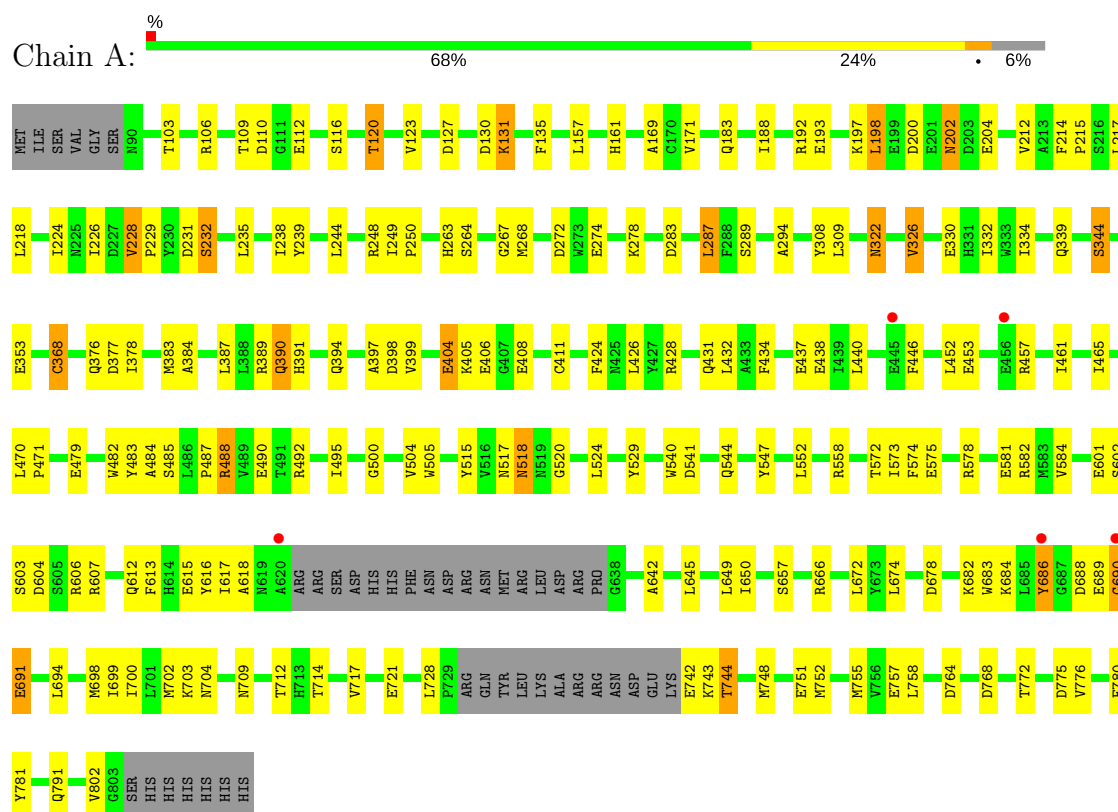
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	84	Total	O	0	0
			84	84		
4	B	68	Total	O	0	0
			68	68		
4	C	49	Total	O	0	0
			49	49		

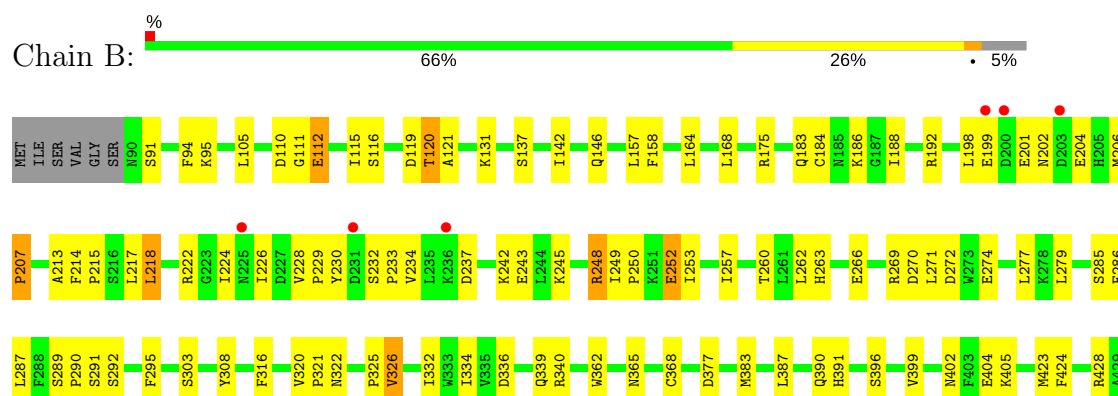
### 3 Residue-property plots

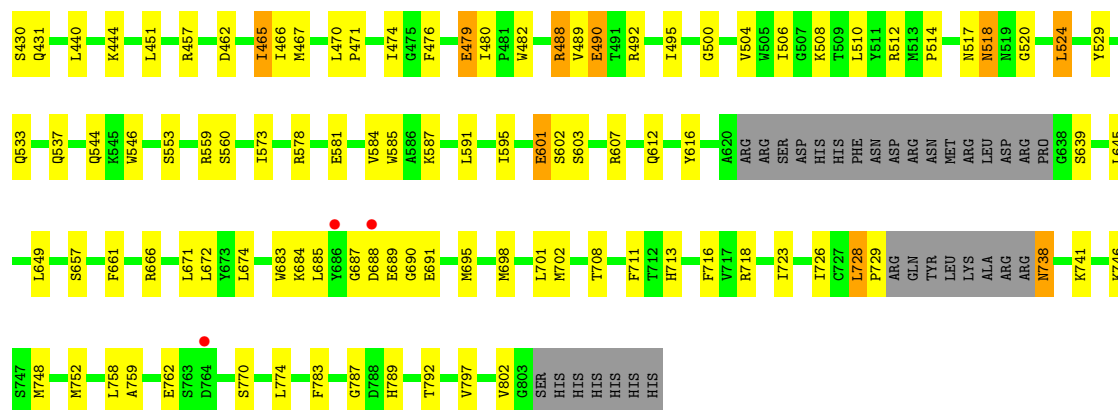
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: Ent-copalyl diphosphate synthase, chloroplastic

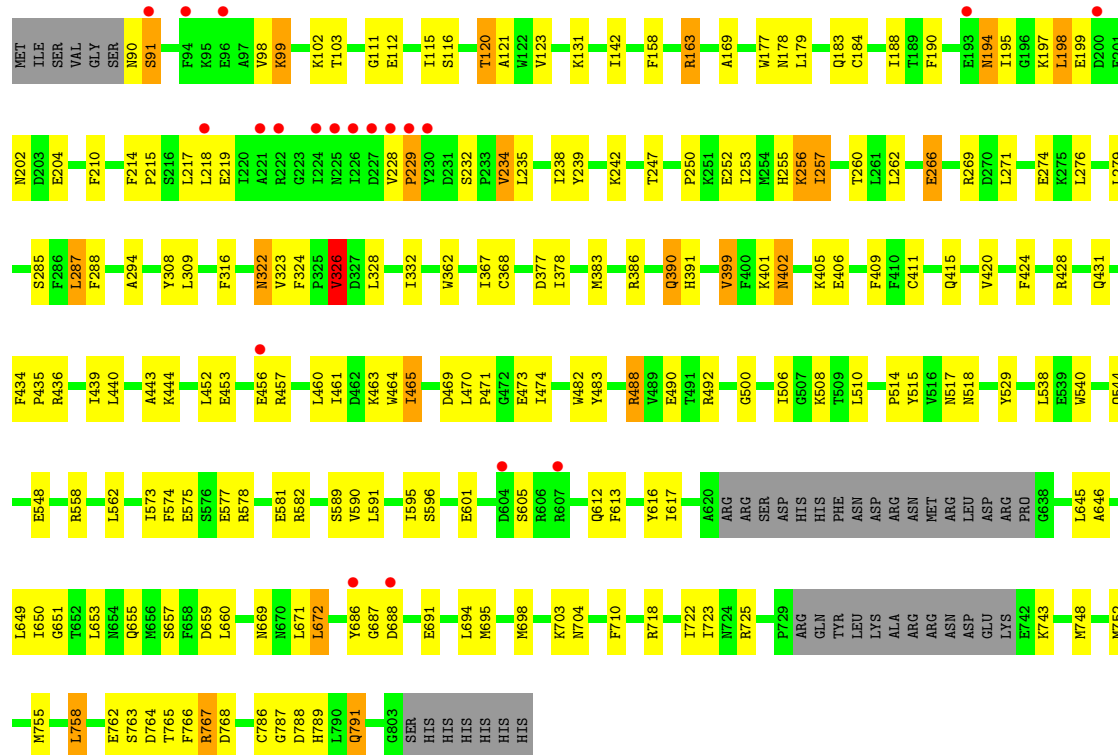


- Molecule 1: Ent-copalyl diphosphate synthase, chloroplastic





● Molecule 1: Ent-copalyl diphosphate synthase, chloroplastic





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	63.74Å 188.55Å 229.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.52 – 2.76 44.52 – 2.76	Depositor EDS
% Data completeness (in resolution range)	86.1 (44.52-2.76) 86.1 (44.52-2.76)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	0.14	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.09 (at 2.77Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, $R_{free}$	0.177 , 0.257 0.170 , 0.254	Depositor DCC
$R_{free}$ test set	1860 reflections (2.96%)	DCC
Wilson B-factor (Å <sup>2</sup> )	46.4	Xtrriage
Anisotropy	0.741	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 59.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	17263	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.88% 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, 0.333 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: A3C, 1PE

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	0.29	0/5785	0.44	0/7817
1	B	0.27	0/5790	0.42	0/7823
1	C	0.27	0/5745	0.42	0/7764
All	All	0.28	0/17320	0.43	0/23404

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	5633	0	5486	123	0
1	B	5649	0	5494	136	0
1	C	5607	0	5452	133	0
2	A	29	0	34	6	0
3	A	64	0	88	2	0
3	B	48	0	66	2	0
3	C	32	0	44	4	0
4	A	84	0	0	0	0
4	B	68	0	0	0	0
4	C	49	0	0	1	0
All	All	17263	0	16664	397	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 12.

All (397) 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:C:163:ARG:HH11	1:C:163:ARG:HG2	1.23	1.04
1:C:431:GLN:HE22	1:C:482:TRP:H	1.10	0.98
1:B:728:LEU:HB3	1:B:729:PRO:HD2	1.44	0.97
1:C:453:GLU:HB3	1:C:457:ARG:HH12	1.27	0.94
1:A:431:GLN:HE22	1:A:482:TRP:H	1.17	0.90
1:B:215:PRO:HB2	1:B:242:LYS:HD2	1.54	0.88
1:B:504:VAL:HG22	1:B:512:ARG:HH12	1.38	0.88
1:C:163:ARG:HH11	1:C:163:ARG:CG	1.88	0.86
1:C:431:GLN:NE2	1:C:482:TRP:H	1.71	0.86
1:A:431:GLN:NE2	1:A:482:TRP:H	1.81	0.79
1:A:686:TYR:HD2	1:A:689:GLU:HG3	1.48	0.78
1:B:518:ASN:ND2	1:B:520:GLY:H	1.81	0.77
1:C:253:ILE:HG23	1:C:257:ILE:HD11	1.66	0.77
1:C:452:LEU:O	1:C:456:GLU:HG2	1.84	0.77
1:A:601:GLU:HG2	1:A:602:SER:H	1.49	0.76
1:A:103:THR:HG22	1:A:106[A]:ARG:HH22	1.50	0.76
1:B:120:THR:HG22	1:B:320:VAL:H	1.52	0.74
1:B:728:LEU:HB3	1:B:729:PRO:CD	2.17	0.73
1:C:562:LEU:HD12	1:C:590:VAL:HG21	1.71	0.72
1:C:453:GLU:HB3	1:C:457:ARG:NH1	2.05	0.71
1:A:112:GLU:HG3	1:A:515:TYR:CE1	2.25	0.71
1:C:215:PRO:HB2	1:C:242:LYS:HD2	1.71	0.71
1:C:195:ILE:HD11	1:C:235:LEU:HD21	1.71	0.71
1:A:272:ASP:OD1	1:A:274:GLU:HB3	1.92	0.70
1:C:169:ALA:HB2	1:C:217:LEU:HD21	1.74	0.70
1:A:518:ASN:ND2	1:A:520:GLY:H	1.90	0.69
1:A:518:ASN:HD22	1:A:520:GLY:N	1.89	0.69
2:A:901:A3C:H11	2:A:901:A3C:H31	1.73	0.68
1:B:683:TRP:HD1	1:B:688:ASP:HA	1.58	0.68
1:C:112:GLU:HG3	1:C:515:TYR:CE1	2.28	0.68
1:C:111:GLY:HA3	1:C:326:VAL:HG12	1.75	0.68
1:A:666:ARG:HG2	1:A:666:ARG:HH11	1.57	0.68
1:C:90:ASN:O	1:C:91:SER:HB2	1.94	0.68
1:A:212:VAL:HG12	1:A:267:GLY:HA3	1.74	0.67
1:B:512:ARG:HG2	1:B:512:ARG:HH11	1.60	0.67
1:C:431:GLN:HE22	1:C:482:TRP:N	1.88	0.67
1:B:207:PRO:HB3	1:B:508:LYS:O	1.94	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:362:TRP:CZ2	1:C:399:VAL:HG21	2.30	0.66
1:A:390:GLN:HG2	1:A:391:HIS:CD2	2.31	0.66
1:A:405:LYS:HG2	1:A:406:GLU:HG2	1.77	0.66
1:B:116:SER:O	1:B:120:THR:HG23	1.96	0.66
1:B:500:GLY:O	1:B:517:ASN:HB3	1.96	0.65
1:C:163:ARG:NH1	1:C:163:ARG:HG2	2.00	0.65
1:A:244:LEU:HD21	1:A:248:ARG:NH2	2.12	0.65
1:B:726:ILE:HG23	1:B:748:MET:HG2	1.78	0.65
1:C:131:LYS:HD2	1:C:131:LYS:N	2.13	0.64
1:C:190:PHE:O	1:C:194:ASN:HB2	1.99	0.62
1:C:362:TRP:HZ2	1:C:399:VAL:HG21	1.64	0.62
1:B:431:GLN:NE2	1:B:482:TRP:H	1.97	0.62
1:A:601:GLU:HG2	1:A:602:SER:N	2.14	0.62
1:A:431:GLN:HE22	1:A:482:TRP:N	1.95	0.62
1:B:691:GLU:O	1:B:695:MET:HG2	2.00	0.62
1:B:430:SER:HA	1:B:440:LEU:HD22	1.81	0.62
1:A:214:PHE:HB3	1:A:215:PRO:HD3	1.80	0.61
1:B:105:LEU:HD21	1:B:334:ILE:HG21	1.81	0.61
1:A:202:ASN:ND2	1:A:202:ASN:H	1.98	0.61
1:A:229:PRO:O	1:A:235:LEU:HD12	1.99	0.61
1:B:326:VAL:O	1:B:326:VAL:HG13	1.99	0.61
1:C:238:ILE:HD12	1:C:239:TYR:N	2.16	0.61
1:A:465:ILE:HD13	1:A:505:TRP:CD2	2.36	0.61
1:B:111:GLY:HA3	1:B:326:VAL:HG12	1.82	0.61
1:C:285:SER:HB3	1:C:308:TYR:CE2	2.36	0.61
1:A:294:ALA:HA	1:A:309:LEU:HD11	1.82	0.60
1:A:518:ASN:ND2	1:A:520:GLY:N	2.49	0.60
1:B:199:GLU:HB2	1:B:234:VAL:HB	1.81	0.60
1:B:518:ASN:HD22	1:B:520:GLY:N	1.99	0.60
1:A:603:SER:HA	1:A:606:ARG:NH2	2.16	0.60
1:B:489:VAL:HG22	1:B:797:VAL:HG21	1.83	0.60
1:C:401:LYS:HG2	1:C:439:ILE:HD12	1.83	0.60
1:A:326:VAL:O	1:A:326:VAL:HG13	2.01	0.60
1:B:431:GLN:HE22	1:B:482:TRP:H	1.50	0.60
1:C:378:ILE:HG22	1:C:411:CYS:HA	1.83	0.60
1:B:518:ASN:HD22	1:B:520:GLY:H	1.46	0.59
1:B:698:MET:HE2	1:B:701:LEU:HD12	1.83	0.59
1:C:266:GLU:OE2	1:C:508:LYS:HE2	2.02	0.59
1:C:657:SER:HB2	1:C:672:LEU:HD12	1.84	0.59
1:B:332:ILE:HG21	1:B:383:MET:HB3	1.84	0.59
1:B:518:ASN:ND2	1:B:520:GLY:N	2.50	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:453:GLU:O	1:A:457:ARG:HG3	2.03	0.59
1:C:218:LEU:HD21	1:C:235:LEU:HD13	1.85	0.59
1:B:232:SER:OG	1:B:233:PRO:HD2	2.03	0.59
1:B:213:ALA:O	1:B:217:LEU:HB2	2.03	0.58
1:A:116:SER:O	1:A:120:THR:HG23	2.03	0.58
1:B:336:ASP:O	1:B:340:ARG:HG3	2.03	0.58
1:A:518:ASN:HD22	1:A:520:GLY:H	1.49	0.58
1:B:533:GLN:O	1:B:537:GLN:HG3	2.03	0.58
1:C:326:VAL:O	1:C:326:VAL:HG13	2.02	0.58
1:A:686:TYR:HB2	1:A:689:GLU:OE1	2.03	0.58
1:C:500:GLY:O	1:C:517:ASN:HB3	2.04	0.58
1:C:646:ALA:O	1:C:650:ILE:HD13	2.02	0.58
1:B:601:GLU:HG2	1:B:602:SER:H	1.68	0.58
1:A:575:GLU:OE1	1:A:578:ARG:HD3	2.04	0.57
2:A:901:A3C:H13B	2:A:901:A3C:O01	2.03	0.57
1:C:752:MET:O	1:C:755:MET:HB3	2.04	0.57
1:B:465:ILE:HG13	1:B:467:MET:HG3	1.85	0.57
1:C:492:ARG:HD3	1:C:529:TYR:CE2	2.39	0.57
1:A:192:ARG:HG2	1:A:229:PRO:HD3	1.87	0.57
1:A:483:TYR:CE1	3:A:911:1PE:H162	2.40	0.57
1:B:404:GLU:HG2	1:B:405:LYS:N	2.19	0.57
1:C:514:PRO:O	1:C:518:ASN:HB2	2.05	0.57
1:A:330:GLU:O	1:A:334:ILE:HG13	2.03	0.57
1:B:214:PHE:HB3	1:B:215:PRO:HD3	1.86	0.57
1:A:263:HIS:CE1	1:A:289:SER:HB2	2.40	0.56
1:A:617:ILE:HG22	1:A:618:ALA:N	2.19	0.56
1:C:420:VAL:HG23	4:C:826:HOH:O	2.06	0.56
1:A:408:GLU:HG2	1:A:446:PHE:CE1	2.40	0.56
1:A:500:GLY:O	1:A:517:ASN:HB3	2.05	0.56
1:B:495:ILE:HG21	1:B:802:VAL:HG11	1.87	0.56
1:C:285:SER:HB3	1:C:308:TYR:CD2	2.41	0.56
1:C:198:LEU:O	1:C:198:LEU:HD22	2.06	0.56
1:C:215:PRO:O	1:C:219:GLU:HG2	2.05	0.56
1:C:718:ARG:O	1:C:722:ILE:HG13	2.05	0.55
1:B:424:PHE:CZ	1:B:428:ARG:HD2	2.41	0.55
1:B:687:GLY:O	1:B:689:GLU:HG2	2.05	0.55
1:A:484:ALA:HB2	1:A:781:TYR:HD2	1.71	0.55
1:B:479:GLU:HG2	1:B:480:ILE:HG12	1.88	0.55
1:C:402:ASN:H	1:C:402:ASN:ND2	2.03	0.55
1:B:657:SER:HB3	1:B:672:LEU:HD12	1.87	0.55
1:C:90:ASN:O	1:C:91:SER:CB	2.55	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:573:ILE:HG23	1:B:578:ARG:NH1	2.22	0.54
1:C:390:GLN:HG2	1:C:391:HIS:CD2	2.42	0.54
1:B:269:ARG:O	1:B:271:LEU:HG	2.06	0.54
1:B:683:TRP:CD1	1:B:688:ASP:HA	2.39	0.54
1:A:212:VAL:CG1	1:A:267:GLY:HA3	2.37	0.54
1:A:657:SER:HB3	1:A:672:LEU:HD12	1.88	0.54
2:A:901:A3C:C11	2:A:901:A3C:H31	2.36	0.54
1:B:424:PHE:HA	1:B:451:LEU:HD11	1.89	0.54
1:A:112:GLU:HG3	1:A:515:TYR:CZ	2.43	0.54
1:B:601:GLU:H	1:B:601:GLU:CD	2.11	0.54
1:C:691:GLU:HG3	1:C:695:MET:HE2	1.89	0.54
1:A:322:ASN:H	1:A:322:ASN:HD22	1.55	0.54
1:A:215:PRO:HG3	1:A:238:ILE:HD13	1.90	0.54
1:A:404:GLU:HG2	1:A:408:GLU:O	2.08	0.54
1:C:322:ASN:HD22	1:C:322:ASN:C	2.11	0.54
1:A:465:ILE:HD12	1:A:504:VAL:O	2.07	0.54
1:B:444:LYS:HE2	3:B:911:1PE:OH6	2.08	0.54
1:C:115:ILE:HD11	1:C:316:PHE:HB3	1.89	0.54
1:A:686:TYR:CD2	1:A:689:GLU:HG3	2.37	0.53
1:A:717:VAL:O	1:A:721:GLU:HG3	2.07	0.53
1:B:168:LEU:HD21	1:B:218:LEU:HD13	1.90	0.53
1:C:271:LEU:HD22	1:C:276:LEU:HD11	1.89	0.53
1:C:390:GLN:HG2	1:C:391:HIS:HD2	1.73	0.53
1:B:285:SER:HB3	1:B:308:TYR:CE2	2.43	0.53
1:A:744:THR:O	1:A:748:MET:HG3	2.09	0.53
1:A:540:TRP:O	1:A:544:GLN:HG3	2.08	0.53
1:B:228:VAL:O	1:B:230:TYR:N	2.43	0.52
1:C:488:ARG:HD3	1:C:574:PHE:CG	2.45	0.52
1:B:404:GLU:HG2	1:B:405:LYS:H	1.75	0.52
1:C:483:TYR:CE1	3:C:911:1PE:H262	2.43	0.52
1:B:506:ILE:HG13	1:B:510:LEU:CD2	2.39	0.52
1:C:288:PHE:HB3	1:C:324:PHE:O	2.08	0.52
1:A:378:ILE:HG22	1:A:411:CYS:HA	1.91	0.52
1:C:444:LYS:HE2	3:C:911:1PE:H241	1.92	0.52
1:A:700:ILE:O	1:A:703:LYS:HG3	2.10	0.52
1:B:726:ILE:O	1:B:728:LEU:HG	2.10	0.52
1:A:202:ASN:HD22	1:A:202:ASN:H	1.58	0.51
1:B:115:ILE:HD11	1:B:316:PHE:HB3	1.91	0.51
1:C:405:LYS:HG2	1:C:406:GLU:HG3	1.92	0.51
1:B:146:GLN:OE1	1:B:186:LYS:HE2	2.10	0.51
1:A:397:ALA:HB3	1:A:437:GLU:HG2	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:484:ALA:HB2	1:A:781:TYR:CD2	2.45	0.51
1:B:672:LEU:HD23	1:B:701:LEU:CD1	2.41	0.51
1:C:99:LYS:O	1:C:103:THR:HG23	2.11	0.51
1:C:163:ARG:NH1	1:C:163:ARG:CG	2.58	0.51
1:A:283:ASP:O	1:A:308:TYR:HB2	2.11	0.50
1:B:462:ASP:HB3	1:B:465:ILE:HG23	1.92	0.50
1:C:575:GLU:OE1	1:C:578:ARG:HD3	2.11	0.50
1:B:603:SER:O	1:B:607:ARG:HG3	2.12	0.50
1:C:322:ASN:HD22	1:C:323:VAL:N	2.10	0.50
1:C:378:ILE:CG2	1:C:411:CYS:HA	2.40	0.50
1:A:378:ILE:HD11	1:A:426:LEU:HD13	1.93	0.50
1:A:700:ILE:HA	1:A:703:LYS:HG3	1.94	0.50
1:B:476:PHE:CE2	1:B:490:GLU:HG2	2.46	0.50
1:C:791:GLN:HA	1:C:791:GLN:HE21	1.76	0.50
1:B:759:ALA:HA	1:B:770:SER:HB2	1.92	0.50
2:A:901:A3C:C31	2:A:901:A3C:C11	2.89	0.50
1:A:198:LEU:HD22	1:A:198:LEU:O	2.12	0.50
1:B:672:LEU:HD23	1:B:701:LEU:HD13	1.94	0.50
1:B:512:ARG:NH1	1:B:512:ARG:HG2	2.25	0.49
1:C:612:GLN:O	1:C:616:TYR:HB2	2.12	0.49
1:B:368:CYS:HA	1:B:377:ASP:OD1	2.11	0.49
1:B:783:PHE:O	1:B:787:GLY:HA3	2.12	0.49
1:A:171:VAL:HG11	1:A:188:ILE:HG12	1.92	0.49
1:B:164:LEU:CD1	1:B:198:LEU:HB2	2.42	0.49
1:C:332:ILE:HG21	1:C:383:MET:HB3	1.93	0.49
1:A:492:ARG:HD3	1:A:529:TYR:CZ	2.47	0.49
1:B:544:GLN:HE21	1:B:559:ARG:HE	1.60	0.49
1:B:451:LEU:HD22	1:B:474:ILE:HG21	1.93	0.49
1:C:748:MET:O	1:C:752:MET:HB2	2.13	0.49
1:C:465:ILE:HG13	1:C:465:ILE:O	2.12	0.49
1:A:495:ILE:HG21	1:A:802:VAL:HG11	1.95	0.48
1:C:463:LYS:HE3	1:C:464:TRP:CZ2	2.48	0.48
1:C:791:GLN:CA	1:C:791:GLN:HE21	2.26	0.48
1:B:91:SER:O	1:B:95:LYS:HG3	2.12	0.48
1:A:485:SER:OG	1:A:490:GLU:HG3	2.13	0.48
1:B:224:ILE:HD11	1:B:226:ILE:HD11	1.95	0.48
1:C:698:MET:HA	1:C:698:MET:CE	2.44	0.48
1:B:332:ILE:CG2	1:B:383:MET:HB3	2.43	0.48
1:C:260:THR:HG23	1:C:415:GLN:HE22	1.78	0.48
1:C:591:LEU:HD13	1:C:649:LEU:HD23	1.95	0.48
1:A:424:PHE:CZ	1:A:428:ARG:HD2	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:581:GLU:HB3	1:A:699:ILE:HG23	1.95	0.48
1:A:686:TYR:HD2	1:A:689:GLU:CG	2.25	0.48
1:B:607:ARG:HG2	1:B:684:LYS:HG2	1.95	0.48
1:A:332:ILE:CD1	1:A:384:ALA:HB2	2.44	0.48
1:B:266:GLU:HA	1:B:295:PHE:CG	2.49	0.48
1:C:184:CYS:O	1:C:188:ILE:HG13	2.14	0.48
1:C:613:PHE:CE1	1:C:650:ILE:HD11	2.49	0.48
1:A:689:GLU:O	1:A:691:GLU:N	2.44	0.47
1:C:158:PHE:CD2	1:C:158:PHE:C	2.88	0.47
1:A:287:LEU:HA	1:A:287:LEU:HD23	1.69	0.47
1:B:728:LEU:CB	1:B:729:PRO:CD	2.89	0.47
1:C:424:PHE:CZ	1:C:428:ARG:HD2	2.49	0.47
1:C:544:GLN:O	1:C:548:GLU:HG3	2.14	0.47
1:A:683:TRP:HD1	1:A:688:ASP:HA	1.80	0.47
1:C:177:TRP:O	1:C:179:LEU:HG	2.14	0.47
1:B:112:GLU:O	1:B:514:PRO:HD2	2.14	0.47
1:C:764:ASP:O	1:C:767:ARG:HB3	2.15	0.47
1:A:215:PRO:HG3	1:A:238:ILE:CD1	2.45	0.47
1:A:434:PHE:CE1	1:A:757:GLU:HA	2.49	0.47
1:A:666:ARG:HG2	1:A:666:ARG:NH1	2.29	0.47
1:B:466:ILE:HG13	1:B:506:ILE:HD11	1.96	0.47
1:B:488:ARG:HD2	1:B:488:ARG:HA	1.49	0.47
1:B:518:ASN:HD22	1:B:518:ASN:C	2.18	0.47
1:B:689:GLU:O	1:B:691:GLU:N	2.35	0.47
1:A:109:THR:OG1	1:A:110:ASP:N	2.47	0.47
1:A:332:ILE:HG23	1:A:387:LEU:HD12	1.97	0.47
1:B:252:GLU:HG3	1:B:253:ILE:N	2.29	0.47
1:B:332:ILE:HG23	1:B:387:LEU:HD12	1.95	0.47
1:C:255:HIS:C	1:C:256:LYS:HD2	2.36	0.47
1:A:389:ARG:HG2	1:A:432:LEU:HD23	1.97	0.46
1:A:575:GLU:H	1:A:575:GLU:CD	2.17	0.46
1:B:222:ARG:NH2	1:B:230:TYR:HB3	2.30	0.46
1:B:290:PRO:HG2	1:B:321:PRO:O	2.15	0.46
1:B:120:THR:CG2	1:B:320:VAL:H	2.24	0.46
1:B:110:ASP:HB3	1:B:325:PRO:HB3	1.97	0.46
1:C:368:CYS:HA	1:C:377:ASP:OD1	2.15	0.46
1:C:651:GLY:O	1:C:655:GLN:HG3	2.15	0.46
1:A:584:VAL:HG11	1:A:702:MET:HG3	1.97	0.46
1:B:470:LEU:N	1:B:471:PRO:CD	2.78	0.46
1:B:119:ASP:OD2	1:B:291:SER:HB3	2.15	0.46
1:B:584:VAL:HB	1:B:702:MET:HG3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:234:VAL:O	1:B:237:ASP:HB3	2.16	0.46
1:B:546:TRP:CE2	1:B:587:LYS:HE2	2.51	0.46
1:A:470:LEU:N	1:A:471:PRO:CD	2.79	0.46
1:A:488:ARG:HD3	1:A:574:PHE:CD2	2.50	0.46
1:C:131:LYS:HD2	1:C:131:LYS:H	1.79	0.46
1:A:390:GLN:HG2	1:A:391:HIS:HD2	1.77	0.46
1:B:365:ASN:HD22	1:B:402:ASN:HD21	1.63	0.46
1:C:198:LEU:HB3	1:C:234:VAL:HG21	1.98	0.46
1:C:215:PRO:CB	1:C:242:LYS:HD2	2.41	0.46
1:C:469:ASP:O	1:C:473:GLU:HG3	2.16	0.46
1:A:235:LEU:HA	1:A:235:LEU:HD23	1.68	0.45
1:A:714:THR:HG23	1:B:204:GLU:OE1	2.16	0.45
1:A:465:ILE:HD13	1:A:505:TRP:CE3	2.51	0.45
1:B:201:GLU:HG2	1:B:202:ASN:N	2.31	0.45
1:C:723:ILE:C	1:C:725:ARG:H	2.19	0.45
2:A:901:A3C:H01	2:A:901:A3C:H11A	1.64	0.45
1:B:716:PHE:C	1:B:716:PHE:CD2	2.89	0.45
1:C:409:PHE:CE2	1:C:443:ALA:HA	2.51	0.45
1:C:704:ASN:N	1:C:704:ASN:HD22	2.14	0.45
1:B:263:HIS:CE1	1:B:289:SER:HB2	2.52	0.45
1:A:607:ARG:HG2	1:A:684:LYS:HD3	1.98	0.45
1:B:546:TRP:CZ2	1:B:587:LYS:HG2	2.51	0.45
1:C:786:CYS:HB3	1:C:789:HIS:CE1	2.52	0.45
1:C:116:SER:O	1:C:120:THR:HG23	2.16	0.45
1:A:332:ILE:HD13	1:A:384:ALA:HB2	1.99	0.45
1:B:272:ASP:OD1	1:B:274:GLU:HB3	2.17	0.45
1:B:492:ARG:HD3	1:B:529:TYR:CZ	2.51	0.45
1:A:488:ARG:HD2	1:A:488:ARG:HA	1.54	0.45
1:B:184:CYS:O	1:B:188:ILE:HG13	2.16	0.45
1:B:728:LEU:CB	1:B:729:PRO:HD2	2.31	0.45
1:B:269:ARG:HH11	1:B:269:ARG:HG2	1.82	0.45
1:C:613:PHE:CD1	1:C:650:ILE:HD11	2.53	0.44
1:A:169:ALA:HB2	1:A:217:LEU:HD21	1.99	0.44
1:A:368:CYS:HA	1:A:377:ASP:OD1	2.18	0.44
1:A:682:LYS:HD2	1:A:690:GLY:H	1.83	0.44
1:C:98:VAL:O	1:C:102:LYS:HB2	2.17	0.44
1:C:199:GLU:OE2	1:C:234:VAL:HA	2.17	0.44
1:C:214:PHE:HB3	1:C:215:PRO:HD3	1.98	0.44
1:C:686:TYR:HB3	1:C:687:GLY:H	1.49	0.44
1:A:492:ARG:HD3	1:A:529:TYR:CE2	2.52	0.44
1:A:776:VAL:O	1:A:780[B]:PHE:HD2	2.01	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:287:LEU:HD23	1:C:287:LEU:HA	1.57	0.44
3:C:931:1PE:H241	3:C:931:1PE:H231	1.79	0.44
1:A:326:VAL:CG1	1:A:326:VAL:O	2.66	0.44
1:B:285:SER:HB3	1:B:308:TYR:CD2	2.52	0.44
1:C:229:PRO:O	1:C:235:LEU:HD12	2.17	0.44
1:C:434:PHE:HB3	1:C:435:PRO:HD2	2.00	0.44
1:C:460:LEU:HD23	1:C:460:LEU:HA	1.74	0.44
1:B:158:PHE:C	1:B:158:PHE:CD2	2.91	0.44
1:B:168:LEU:CD2	1:B:218:LEU:HD13	2.48	0.44
1:C:506:ILE:HG12	1:C:510:LEU:CD2	2.48	0.44
1:C:121:ALA:HA	1:C:142:ILE:HD11	1.98	0.44
1:C:748:MET:O	1:C:752:MET:N	2.51	0.44
1:B:201:GLU:HG2	1:B:202:ASN:ND2	2.33	0.43
1:C:328:LEU:HD13	1:C:367:ILE:HD12	2.00	0.43
1:A:657:SER:HB3	1:A:672:LEU:CD1	2.47	0.43
1:A:228:VAL:HG22	1:A:229:PRO:HD2	2.00	0.43
1:B:206:MET:O	1:B:206:MET:HG3	2.18	0.43
1:B:661:PHE:HA	1:B:666:ARG:H	1.83	0.43
1:C:460:LEU:HD22	1:C:474:ILE:HD12	1.99	0.43
1:B:396:SER:O	1:B:399:VAL:HG23	2.18	0.43
1:B:698:MET:CE	1:B:698:MET:HA	2.48	0.43
1:A:249:ILE:HA	1:A:250:PRO:HD3	1.80	0.43
1:A:613:PHE:CE1	1:A:650:ILE:HD11	2.53	0.43
1:B:121:ALA:HA	1:B:142:ILE:HD11	2.00	0.43
1:B:585:TRP:C	1:B:585:TRP:CD1	2.92	0.43
1:B:215:PRO:CB	1:B:242:LYS:HD2	2.37	0.43
1:B:245:LYS:O	1:B:249:ILE:HG12	2.18	0.43
1:C:332:ILE:CG2	1:C:383:MET:HB3	2.49	0.43
1:C:470:LEU:N	1:C:471:PRO:CD	2.81	0.43
1:C:488:ARG:HD3	1:C:574:PHE:CD2	2.54	0.43
1:C:710:PHE:CD1	1:C:766:PHE:HD2	2.37	0.43
1:A:440:LEU:HD23	1:A:440:LEU:HA	1.83	0.43
1:B:277:LEU:HD23	1:B:286:PHE:HZ	1.84	0.43
1:C:492:ARG:HD3	1:C:529:TYR:CZ	2.54	0.43
3:B:911:1PE:H252	3:B:911:1PE:H262	1.61	0.43
1:C:613:PHE:O	1:C:617:ILE:HG13	2.18	0.43
1:B:262:LEU:HD11	1:B:279:LEU:HB3	2.00	0.43
1:C:440:LEU:HA	1:C:440:LEU:HD23	1.83	0.43
1:C:540:TRP:O	1:C:544:GLN:HG3	2.19	0.43
1:A:612:GLN:HG3	1:A:642:ALA:HB3	1.99	0.42
1:A:742:GLU:HG2	1:A:743:LYS:N	2.33	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:339:GLN:NE2	1:B:391:HIS:NE2	2.63	0.42
1:C:786:CYS:O	1:C:788:ASP:N	2.51	0.42
1:A:339:GLN:HA	1:A:344:SER:HB3	2.02	0.42
1:A:704:ASN:HD22	1:A:704:ASN:N	2.17	0.42
2:A:901:A3C:H12	2:A:901:A3C:H02	1.79	0.42
1:A:483:TYR:CZ	3:A:911:1PE:H162	2.55	0.42
1:B:198:LEU:C	1:B:198:LEU:HD13	2.40	0.42
1:A:224:ILE:HG13	1:A:226:ILE:HG13	2.01	0.42
1:B:601:GLU:HG2	1:B:602:SER:N	2.35	0.42
1:C:573:ILE:O	1:C:582:ARG:HD3	2.20	0.42
1:A:584:VAL:HG22	1:A:698:MET:HB3	2.02	0.42
1:B:105:LEU:CD2	1:B:334:ILE:HG21	2.50	0.42
1:C:116:SER:O	1:C:120:THR:CG2	2.67	0.42
1:C:228:VAL:HG13	1:C:229:PRO:HD2	2.01	0.42
1:A:231:ASP:O	1:A:232:SER:C	2.58	0.42
1:B:105:LEU:HD21	1:B:334:ILE:CG2	2.49	0.42
1:B:215:PRO:HG2	1:B:242:LYS:HB2	2.02	0.42
1:C:210:PHE:C	1:C:210:PHE:CD2	2.93	0.42
1:A:572:THR:HG21	1:A:775:ASP:HB3	2.01	0.42
1:A:742:GLU:OE1	1:A:742:GLU:N	2.53	0.41
1:B:112:GLU:OE2	1:B:112:GLU:HA	2.20	0.41
1:B:738:ASN:O	1:B:741:LYS:HB3	2.19	0.41
1:C:123:VAL:HG22	1:C:309:LEU:HD22	2.02	0.41
1:C:595:ILE:HG12	1:C:645:LEU:HD11	2.02	0.41
1:A:244:LEU:HD23	1:A:244:LEU:C	2.40	0.41
1:C:257:ILE:N	1:C:257:ILE:HD13	2.35	0.41
1:A:127:ASP:OD1	1:A:131:LYS:HD3	2.19	0.41
1:A:238:ILE:HD12	1:A:239:TYR:N	2.35	0.41
1:C:694:LEU:O	1:C:698:MET:HG2	2.20	0.41
1:C:589:SER:HA	1:C:695:MET:CE	2.51	0.41
1:A:674:LEU:HD13	1:B:457:ARG:HG3	2.02	0.41
1:B:340:ARG:HD2	1:B:428:ARG:NH2	2.35	0.41
1:B:612:GLN:O	1:B:616:TYR:HB2	2.20	0.41
1:C:326:VAL:CG1	1:C:326:VAL:O	2.68	0.41
1:A:694:LEU:O	1:A:698:MET:HG2	2.21	0.41
1:C:252:GLU:OE1	1:C:252:GLU:HA	2.21	0.41
1:C:655:GLN:O	1:C:659:ASP:HB2	2.20	0.41
1:A:130:ASP:OD1	1:A:130:ASP:N	2.54	0.41
1:A:547:TYR:HA	1:A:552:LEU:HD12	2.02	0.41
1:A:748:MET:CE	1:A:780[A]:PHE:HE1	2.33	0.41
1:C:202:ASN:OD1	1:C:204:GLU:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:660:LEU:O	1:C:660:LEU:HG	2.20	0.41
1:A:573:ILE:O	1:A:582:ARG:HD3	2.20	0.41
1:B:249:ILE:HA	1:B:250:PRO:HD3	1.87	0.41
1:B:269:ARG:NH1	1:B:269:ARG:HG2	2.35	0.41
1:B:581:GLU:O	1:B:584:VAL:HG12	2.21	0.41
1:C:190:PHE:CD2	1:C:190:PHE:C	2.94	0.41
1:C:386:ARG:NH2	1:C:428:ARG:NH2	2.69	0.41
1:B:188:ILE:O	1:B:192:ARG:HB2	2.21	0.41
1:A:332:ILE:HG21	1:A:383:MET:HB3	2.02	0.41
1:B:362:TRP:CZ2	1:B:399:VAL:HG21	2.56	0.41
1:B:774:LEU:HA	1:B:774:LEU:HD23	1.83	0.41
1:C:294:ALA:HA	1:C:309:LEU:HD11	2.01	0.41
1:A:488:ARG:HD3	1:A:574:PHE:CG	2.56	0.41
1:A:123:VAL:HG13	1:A:135:PHE:CD2	2.57	0.40
1:B:591:LEU:O	1:B:595:ILE:HG13	2.21	0.40
1:C:444:LYS:NZ	3:C:911:1PE:H122	2.35	0.40
1:A:772:THR:O	1:A:776:VAL:HG23	2.21	0.40
1:B:248:ARG:HD2	1:B:248:ARG:HA	1.95	0.40
1:C:601:GLU:HG2	1:C:605:SER:OG	2.21	0.40
1:A:161:HIS:CD2	1:A:198:LEU:HD21	2.56	0.40
1:B:175:ARG:HD3	1:B:175:ARG:O	2.22	0.40
1:A:678:ASP:OD2	1:B:457:ARG:NH2	2.55	0.40
1:B:758:LEU:HA	1:B:758:LEU:HD23	1.87	0.40
1:A:231:ASP:O	1:A:232:SER:O	2.40	0.40
1:A:368:CYS:HB3	1:A:376:GLN:HA	2.03	0.40
1:B:115:ILE:CD1	1:B:316:PHE:HB3	2.51	0.40
1:C:581:GLU:OE2	1:C:703:LYS:HG2	2.22	0.40
1:B:94:PHE:CE2	1:B:524:LEU:CD2	3.05	0.40
1:C:262:LEU:HD11	1:C:279:LEU:HB3	2.04	0.40
1:C:645:LEU:HD22	1:C:645:LEU:O	2.22	0.40
1:C:649:LEU:O	1:C:653:LEU:HG	2.21	0.40
1:C:758:LEU:O	1:C:762:GLU:HG2	2.21	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries

of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	683/727 (94%)	651 (95%)	29 (4%)	3 (0%)	38	69
1	B	684/727 (94%)	628 (92%)	50 (7%)	6 (1%)	20	49
1	C	679/727 (93%)	615 (91%)	52 (8%)	12 (2%)	10	28
All	All	2046/2181 (94%)	1894 (93%)	131 (6%)	21 (1%)	18	46

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	232	SER
1	B	601	GLU
1	C	91	SER
1	C	763	SER
1	A	690	GLY
1	B	229	PRO
1	B	690	GLY
1	B	728	LEU
1	C	194	ASN
1	C	326	VAL
1	C	229	PRO
1	C	787	GLY
1	C	197	LYS
1	C	399	VAL
1	B	207	PRO
1	B	711	PHE
1	C	232	SER
1	C	266	GLU
1	C	688	ASP
1	C	250	PRO
1	A	487	PRO

### 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	608/644 (94%)	555 (91%)	53 (9%)	12	30
1	B	609/644 (95%)	565 (93%)	44 (7%)	17	40
1	C	604/644 (94%)	569 (94%)	35 (6%)	23	52
All	All	1821/1932 (94%)	1689 (93%)	132 (7%)	16	40

All (132) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	120	THR
1	A	131	LYS
1	A	157	LEU
1	A	183	GLN
1	A	193	GLU
1	A	197	LYS
1	A	198	LEU
1	A	200	ASP
1	A	202	ASN
1	A	204	GLU
1	A	218	LEU
1	A	228	VAL
1	A	264	SER
1	A	268	MET
1	A	278	LYS
1	A	287	LEU
1	A	322	ASN
1	A	326	VAL
1	A	344	SER
1	A	353	GLU
1	A	368	CYS
1	A	390	GLN
1	A	394	GLN
1	A	398	ASP
1	A	399	VAL
1	A	404	GLU
1	A	438	GLU
1	A	452	LEU
1	A	461	ILE
1	A	479	GLU
1	A	488	ARG
1	A	518	ASN
1	A	524	LEU
1	A	541	ASP

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Mol	Chain	Res	Type
1	A	558	ARG
1	A	604	ASP
1	A	615	GLU
1	A	616	TYR
1	A	645	LEU
1	A	649	LEU
1	A	686	TYR
1	A	691	GLU
1	A	709	ASN
1	A	712	THR
1	A	728	LEU
1	A	744	THR
1	A	751	GLU
1	A	752	MET
1	A	755	MET
1	A	758	LEU
1	A	764	ASP
1	A	768	ASP
1	A	791	GLN
1	B	112	GLU
1	B	120	THR
1	B	131	LYS
1	B	137	SER
1	B	157	LEU
1	B	183	GLN
1	B	218	LEU
1	B	243	GLU
1	B	248	ARG
1	B	252	GLU
1	B	257	ILE
1	B	260	THR
1	B	270	ASP
1	B	287	LEU
1	B	292	SER
1	B	303	SER
1	B	322	ASN
1	B	326	VAL
1	B	390	GLN
1	B	423	MET
1	B	465	ILE
1	B	479	GLU
1	B	488	ARG

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Mol	Chain	Res	Type
1	B	490	GLU
1	B	518	ASN
1	B	524	LEU
1	B	553	SER
1	B	560	SER
1	B	639	SER
1	B	645	LEU
1	B	649	LEU
1	B	671	LEU
1	B	674	LEU
1	B	685	LEU
1	B	708	THR
1	B	713	HIS
1	B	718	ARG
1	B	723	ILE
1	B	738	ASN
1	B	746	LYS
1	B	752	MET
1	B	762	GLU
1	B	789	HIS
1	B	792	THR
1	C	99	LYS
1	C	120	THR
1	C	163	ARG
1	C	178	ASN
1	C	183	GLN
1	C	198	LEU
1	C	234	VAL
1	C	247	THR
1	C	256	LYS
1	C	257	ILE
1	C	269	ARG
1	C	274	GLU
1	C	287	LEU
1	C	322	ASN
1	C	326	VAL
1	C	390	GLN
1	C	402	ASN
1	C	436	ARG
1	C	461	ILE
1	C	465	ILE
1	C	488	ARG

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Mol	Chain	Res	Type
1	C	490	GLU
1	C	538	LEU
1	C	558	ARG
1	C	577	GLU
1	C	596	SER
1	C	669	ASN
1	C	671	LEU
1	C	672	LEU
1	C	743	LYS
1	C	758	LEU
1	C	765	THR
1	C	767	ARG
1	C	768	ASP
1	C	791	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	145	ASN
1	A	202	ASN
1	A	322	ASN
1	A	339	GLN
1	A	391	HIS
1	A	394	GLN
1	A	431	GLN
1	A	502	ASN
1	A	518	ASN
1	A	535	GLN
1	A	655	GLN
1	A	664	HIS
1	A	669	ASN
1	A	704	ASN
1	A	713	HIS
1	B	282	GLN
1	B	365	ASN
1	B	390	GLN
1	B	415	GLN
1	B	431	GLN
1	B	449	ASN
1	B	518	ASN
1	B	527	GLN
1	B	535	GLN

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Mol	Chain	Res	Type
1	B	544	GLN
1	B	580	HIS
1	B	655	GLN
1	B	664	HIS
1	B	791	GLN
1	C	183	GLN
1	C	299	GLN
1	C	322	ASN
1	C	390	GLN
1	C	402	ASN
1	C	415	GLN
1	C	431	GLN
1	C	518	ASN
1	C	612	GLN
1	C	614	HIS
1	C	655	GLN
1	C	669	ASN
1	C	704	ASN
1	C	789	HIS
1	C	791	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 carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

10 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length

(or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	1PE	A	811	-	15,15,15	0.54	0	14,14,14	0.77	0
2	A3C	A	901	-	29,30,30	0.90	0	39,47,47	2.63	11 (28%)
3	1PE	A	911	-	15,15,15	0.57	0	14,14,14	0.76	0
3	1PE	A	921	-	15,15,15	0.57	0	14,14,14	0.77	0
3	1PE	A	931	-	15,15,15	0.62	0	14,14,14	0.61	0
3	1PE	B	911	-	15,15,15	0.56	0	14,14,14	0.80	0
3	1PE	B	921	-	15,15,15	0.57	0	14,14,14	0.74	0
3	1PE	B	931	-	15,15,15	0.61	0	14,14,14	0.69	0
3	1PE	C	911	-	15,15,15	0.57	0	14,14,14	0.74	0
3	1PE	C	931	-	15,15,15	0.55	0	14,14,14	0.82	0

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
3	1PE	A	811	-	-	0/13/13/13	0/0/0/0
2	A3C	A	901	-	-	0/18/53/53	0/2/2/2
3	1PE	A	911	-	-	0/13/13/13	0/0/0/0
3	1PE	A	921	-	-	0/13/13/13	0/0/0/0
3	1PE	A	931	-	-	0/13/13/13	0/0/0/0
3	1PE	B	911	-	-	0/13/13/13	0/0/0/0
3	1PE	B	921	-	-	0/13/13/13	0/0/0/0
3	1PE	B	931	-	-	0/13/13/13	0/0/0/0
3	1PE	C	911	-	-	0/13/13/13	0/0/0/0
3	1PE	C	931	-	-	0/13/13/13	0/0/0/0

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	901	A3C	C23-C22-C21	-6.76	104.85	113.54
2	A	901	A3C	C23-C22-C31	-4.09	115.48	122.19
2	A	901	A3C	C34-C30-C29	-3.72	102.13	108.24
2	A	901	A3C	C28-C29-C30	-2.86	109.54	113.19
2	A	901	A3C	C24-C25-C26	-2.05	111.27	114.27
2	A	901	A3C	C30-C21-C22	2.80	112.93	109.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	901	A3C	C23-C24-C25	2.95	115.63	110.91
2	A	901	A3C	C29-C30-C25	2.98	111.80	108.04
2	A	901	A3C	C34-C30-C25	4.78	121.78	112.94
2	A	901	A3C	C24-C23-C22	5.19	119.03	110.96
2	A	901	A3C	C21-C22-C31	9.04	139.63	124.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	901	A3C	6	0
3	A	911	1PE	2	0
3	B	911	1PE	2	0
3	C	911	1PE	3	0
3	C	931	1PE	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	685/727 (94%)	-0.22	5 (0%) 87 85	26, 46, 94, 133	0
1	B	689/727 (94%)	-0.17	9 (1%) 77 74	28, 55, 100, 177	0
1	C	685/727 (94%)	-0.05	20 (2%) 52 46	30, 57, 106, 173	0
All	All	2059/2181 (94%)	-0.15	34 (1%) 70 67	26, 53, 101, 177	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	228	VAL	5.1
1	B	199	GLU	4.5
1	C	229	PRO	4.2
1	B	200	ASP	3.7
1	C	686	TYR	3.6
1	B	686	TYR	3.4
1	C	222	ARG	3.3
1	B	236	LYS	3.3
1	C	688	ASP	3.2
1	A	686	TYR	3.2
1	C	224	ILE	3.2
1	B	688	ASP	3.1
1	C	230	TYR	3.1
1	B	764	ASP	3.1
1	B	203	ASP	2.9
1	C	226	ILE	2.9
1	C	221	ALA	2.9
1	C	200	ASP	2.7
1	C	456	GLU	2.7
1	A	620	ALA	2.6
1	C	604	ASP	2.3
1	A	690	GLY	2.3
1	C	225	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	218	LEU	2.2
1	C	227	ASP	2.2
1	A	456	GLU	2.2
1	B	231	ASP	2.2
1	C	607	ARG	2.2
1	A	445	GLU	2.1
1	B	225	ASN	2.1
1	C	91	SER	2.1
1	C	193	GLU	2.1
1	C	94	PHE	2.0
1	C	96	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, 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
3	1PE	A	931	16/16	0.86	0.51	16.86	61,78,96,98	0
3	1PE	B	931	16/16	0.86	0.31	7.66	55,77,98,102	0
3	1PE	A	921	16/16	0.94	0.23	3.97	42,58,77,92	0
3	1PE	B	921	16/16	0.95	0.21	2.97	54,69,89,97	0
3	1PE	C	931	16/16	0.89	0.23	2.73	61,80,95,97	0
3	1PE	C	911	16/16	0.89	0.26	2.10	54,79,93,95	0
3	1PE	A	811	16/16	0.95	0.24	0.90	41,60,79,87	0
3	1PE	A	911	16/16	0.90	0.26	0.72	60,72,90,90	0
3	1PE	B	911	16/16	0.94	0.17	0.54	61,72,95,96	0
2	A3C	A	901	29/29	0.95	0.15	0.25	39,52,112,143	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.