



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:37 PM GMT

PDB ID : 4A0H  
Title : Structure of bifunctional DAPA aminotransferase-DTB synthetase from *Arabidopsis thaliana* bound to 7-keto 8-amino pelargonic acid (KAPA)  
Authors : Cobessi, D.; Dumas, R.; Pautre, V.; Meinguet, C.; Ferrer, J.L.; Alban, C.  
Deposited on : 2011-09-09  
Resolution : 2.81 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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

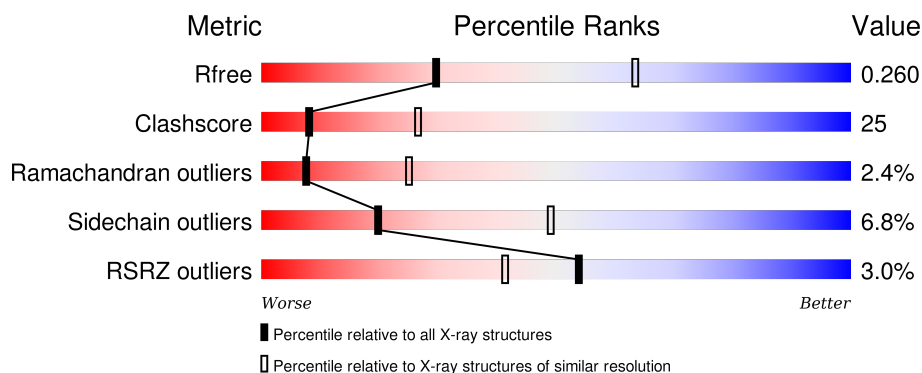
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.81 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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	831	<div> <div>2%</div> <div>52%</div> <div>34%</div> <div>10%</div> </div>
1	B	831	<div> <div>3%</div> <div>50%</div> <div>34%</div> <div>11%</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	KAP	A	1809	-	-	-	X
3	KAP	B	1808	-	-	-	X
4	TLA	A	1810	-	-	X	-
4	TLA	B	1810	-	-	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11459 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 ADENOSYLMETHIONINE-8-AMINO-7-OXONONANOATE AMINOTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	748	Total	C	N	O	S	0	1	0
			5732	3672	964	1065	31			
1	B	738	Total	C	N	O	S	0	1	0
			5625	3608	947	1039	31			

There are 40 discrepancies between the modelled and reference sequences:

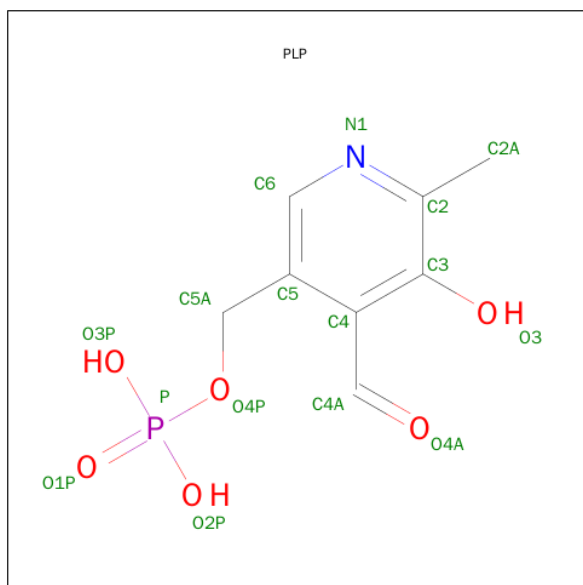
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	GLY	-	EXPRESSION TAG	UNP B0F481
A	-18	SER	-	EXPRESSION TAG	UNP B0F481
A	-17	SER	-	EXPRESSION TAG	UNP B0F481
A	-16	HIS	-	EXPRESSION TAG	UNP B0F481
A	-15	HIS	-	EXPRESSION TAG	UNP B0F481
A	-14	HIS	-	EXPRESSION TAG	UNP B0F481
A	-13	HIS	-	EXPRESSION TAG	UNP B0F481
A	-12	HIS	-	EXPRESSION TAG	UNP B0F481
A	-11	HIS	-	EXPRESSION TAG	UNP B0F481
A	-10	SER	-	EXPRESSION TAG	UNP B0F481
A	-9	SER	-	EXPRESSION TAG	UNP B0F481
A	-8	GLY	-	EXPRESSION TAG	UNP B0F481
A	-7	LEU	-	EXPRESSION TAG	UNP B0F481
A	-6	VAL	-	EXPRESSION TAG	UNP B0F481
A	-5	PRO	-	EXPRESSION TAG	UNP B0F481
A	-4	ARG	-	EXPRESSION TAG	UNP B0F481
A	-3	GLY	-	EXPRESSION TAG	UNP B0F481
A	-2	SER	-	EXPRESSION TAG	UNP B0F481
A	-1	HIS	-	EXPRESSION TAG	UNP B0F481
A	0	MET	-	EXPRESSION TAG	UNP B0F481
B	-19	GLY	-	EXPRESSION TAG	UNP B0F481
B	-18	SER	-	EXPRESSION TAG	UNP B0F481
B	-17	SER	-	EXPRESSION TAG	UNP B0F481
B	-16	HIS	-	EXPRESSION TAG	UNP B0F481

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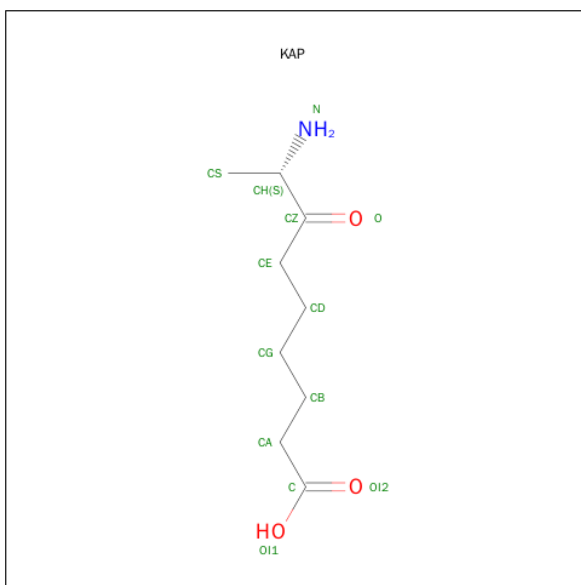
Chain	Residue	Modelled	Actual	Comment	Reference
B	-15	HIS	-	EXPRESSION TAG	UNP B0F481
B	-14	HIS	-	EXPRESSION TAG	UNP B0F481
B	-13	HIS	-	EXPRESSION TAG	UNP B0F481
B	-12	HIS	-	EXPRESSION TAG	UNP B0F481
B	-11	HIS	-	EXPRESSION TAG	UNP B0F481
B	-10	SER	-	EXPRESSION TAG	UNP B0F481
B	-9	SER	-	EXPRESSION TAG	UNP B0F481
B	-8	GLY	-	EXPRESSION TAG	UNP B0F481
B	-7	LEU	-	EXPRESSION TAG	UNP B0F481
B	-6	VAL	-	EXPRESSION TAG	UNP B0F481
B	-5	PRO	-	EXPRESSION TAG	UNP B0F481
B	-4	ARG	-	EXPRESSION TAG	UNP B0F481
B	-3	GLY	-	EXPRESSION TAG	UNP B0F481
B	-2	SER	-	EXPRESSION TAG	UNP B0F481
B	-1	HIS	-	EXPRESSION TAG	UNP B0F481
B	0	MET	-	EXPRESSION TAG	UNP B0F481

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula:  $C_8H_{10}NO_6P$ ).



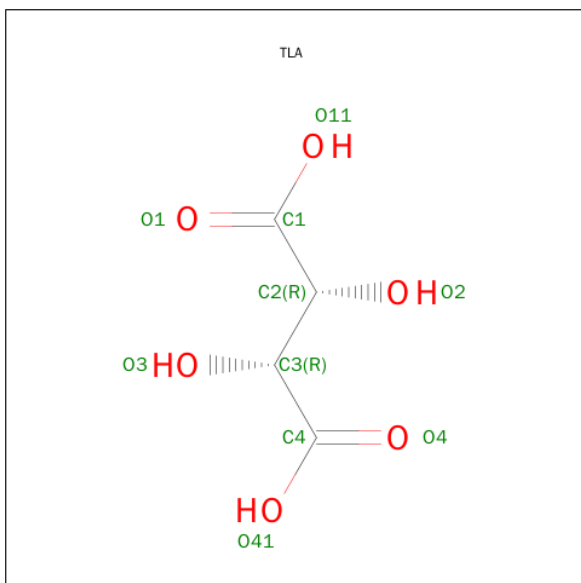
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 3 is 7-KETO-8-AMINOPELARGONIC ACID (three-letter code: KAP) (formula:  $C_9H_{17}NO_3$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			13	9	1	3		
3	A	1	Total	C	N	O	0	0
			13	9	1	3		
3	B	1	Total	C	N	O	0	0
			13	9	1	3		
3	B	1	Total	C	N	O	0	0
			13	9	1	3		

- Molecule 4 is L(+)-TARTARIC ACID (three-letter code: TLA) (formula: C<sub>4</sub>H<sub>6</sub>O<sub>6</sub>).

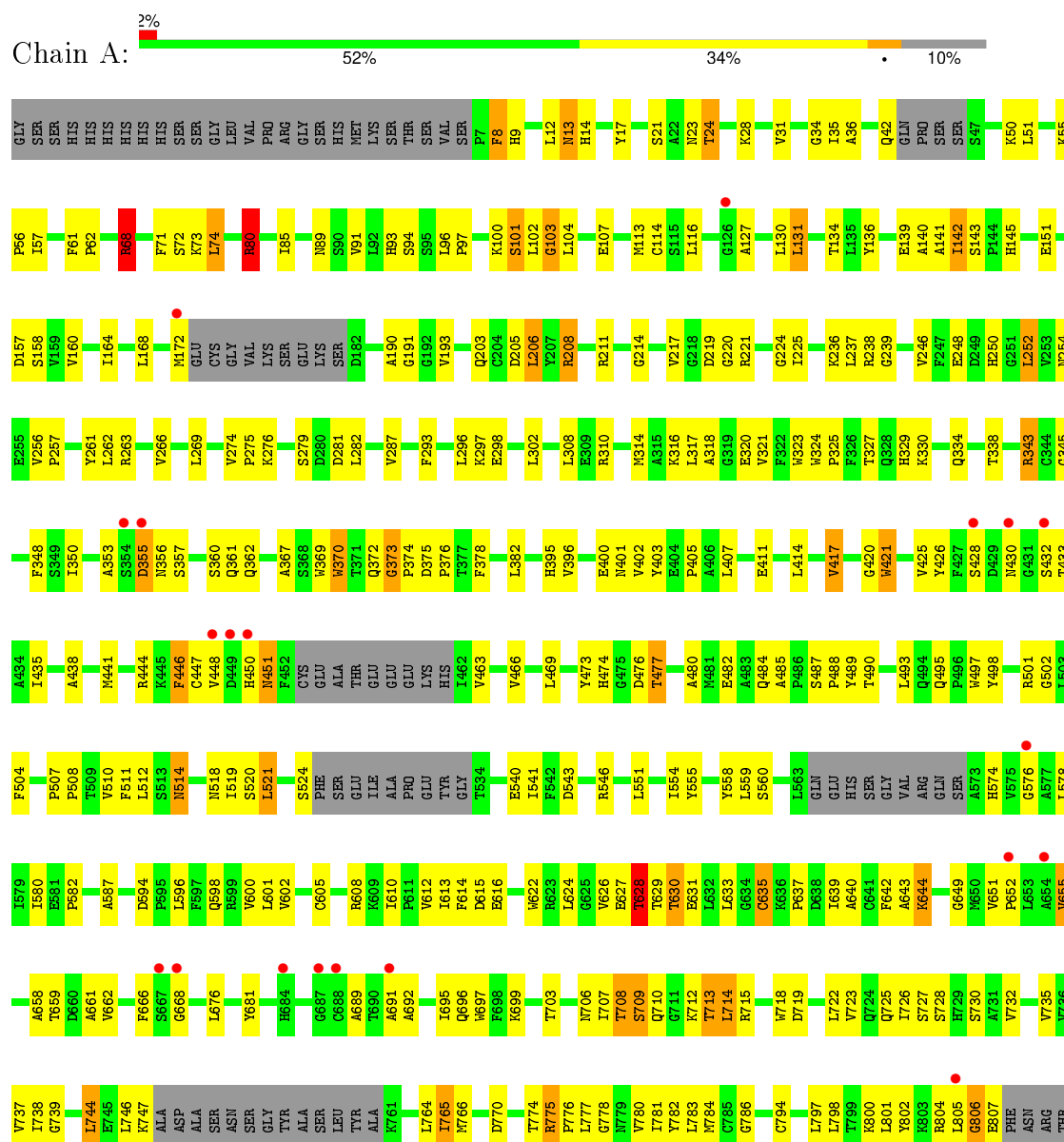


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	4	6		
4	B	1	Total	C	O	0	0
			10	4	6		

### 3 Residue-property plots

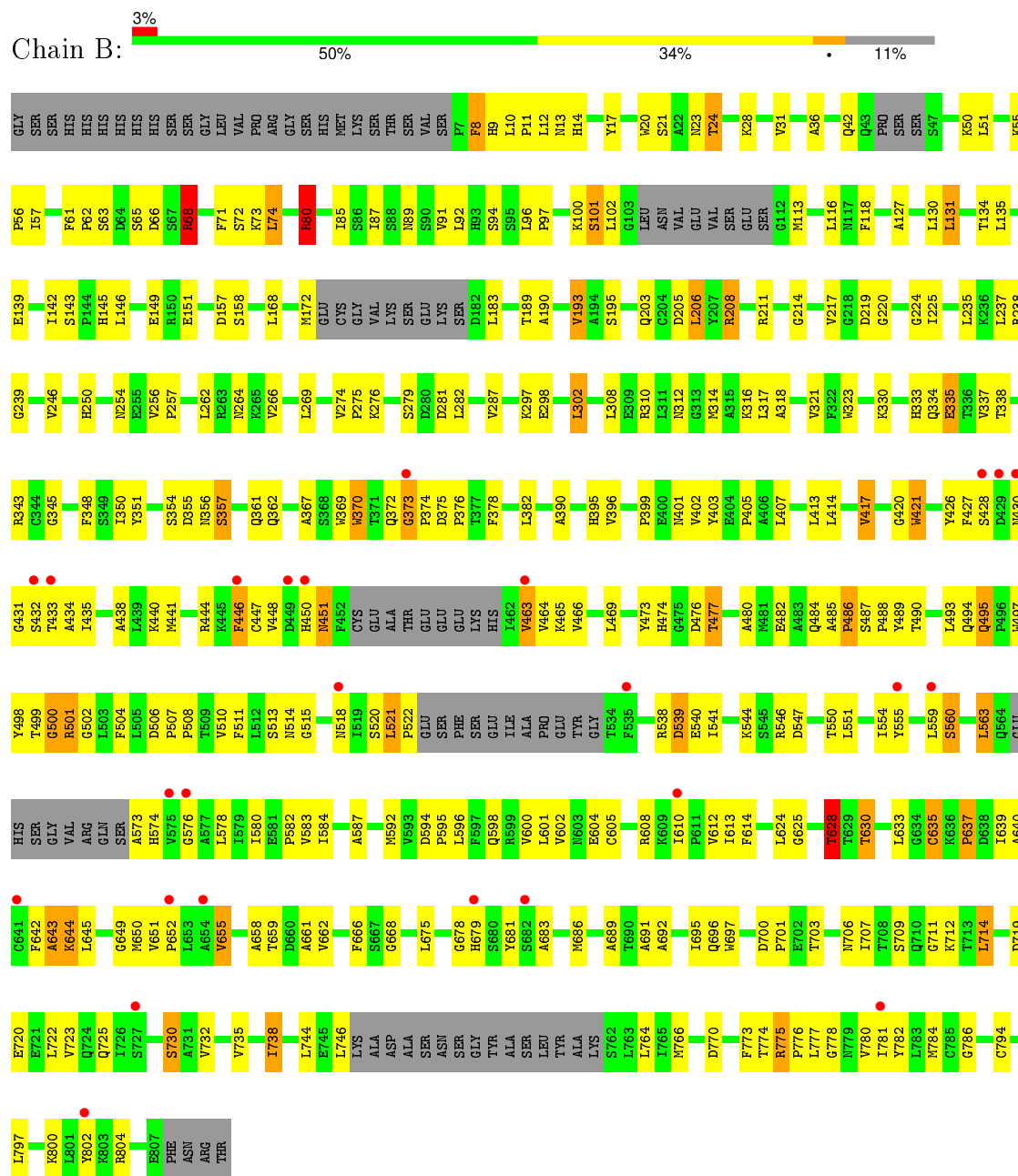
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

- Molecule 1: ADENOSYLMETHIONINE-8-AMINO-7-OXONONANOATE AMINOTRANSFERASE



- Molecule 1: ADENOSYLMETHIONINE-8-AMINO-7-OXONONANOATE AMINOTRANSFERASE





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	235.32Å 76.94Å 89.22Å 90.00° 109.89° 90.00°	Depositor
Resolution (Å)	44.57 – 2.81 44.57 – 2.81	Depositor EDS
% Data completeness (in resolution range)	99.1 (44.57-2.81) 99.2 (44.57-2.81)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.88 (at 2.81Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.193 , 0.262 0.194 , 0.260	Depositor DCC
$R_{free}$ test set	1822 reflections (5.24%)	DCC
Wilson B-factor (Å <sup>2</sup> )	52.7	Xtriage
Anisotropy	0.589	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 51.9	EDS
Estimated twinning fraction	0.021 for -h-2*k,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 36581 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	11459	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% 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: TLA, KAP, PLP

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	0.52	1/5866 (0.0%)	0.69	7/7975 (0.1%)
1	B	0.51	1/5757 (0.0%)	0.66	3/7832 (0.0%)
All	All	0.52	2/11623 (0.0%)	0.67	10/15807 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	628	THR	CB-CG2	-9.08	1.22	1.52
1	B	628	THR	CB-CG2	-7.13	1.28	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	80	ARG	NE-CZ-NH2	8.97	124.79	120.30
1	B	80	ARG	CG-CD-NE	8.25	129.13	111.80
1	A	68	ARG	CG-CD-NE	7.54	127.64	111.80
1	B	628	THR	OG1-CB-CG2	-7.43	92.92	110.00
1	A	80	ARG	CD-NE-CZ	-7.29	113.39	123.60
1	A	80	ARG	CG-CD-NE	6.48	125.42	111.80
1	A	80	ARG	NE-CZ-NH1	-6.35	117.12	120.30
1	A	628	THR	OG1-CB-CG2	-6.21	95.71	110.00
1	B	68	ARG	CG-CD-NE	5.34	123.01	111.80
1	A	628	THR	CB-CA-C	-5.14	97.72	111.60

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5732	0	5648	308	0
1	B	5625	0	5524	293	1
2	A	15	0	6	3	0
2	B	15	0	6	4	0
3	A	26	0	32	5	0
3	B	26	0	32	7	0
4	A	10	0	4	5	0
4	B	10	0	4	4	0
All	All	11459	0	11256	559	1

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 25.

All (559) 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:80:ARG:NH1	1:A:287:VAL:HG21	1.66	1.11
1:A:659:THR:HG22	1:A:661:ALA:H	1.15	1.10
1:B:659:THR:HG22	1:B:661:ALA:H	1.18	1.04
1:B:65:SER:HB3	1:B:68:ARG:NH1	1.77	0.99
1:A:80:ARG:HH11	1:A:287:VAL:HG21	1.26	0.99
1:A:220:GLY:O	1:A:252[A]:LEU:HD21	1.63	0.97
1:A:252[A]:LEU:HD12	1:A:252[A]:LEU:H	1.28	0.95
1:B:446:PHE:CD2	1:B:576:GLY:HA2	2.02	0.94
1:A:446:PHE:CD2	1:A:576:GLY:HA2	2.05	0.92
1:B:208:ARG:HH12	1:B:211:ARG:HD3	1.31	0.92
1:A:493:LEU:HB3	1:B:444:ARG:HH21	1.36	0.91
1:A:55:LYS:NZ	4:A:1810:TLA:O4	2.04	0.90
1:A:208:ARG:HH12	1:A:211:ARG:HD3	1.36	0.89
1:B:68:ARG:HB3	1:B:68:ARG:HH11	1.35	0.88
1:A:521:LEU:CD1	1:A:521:LEU:N	2.35	0.88
1:B:560:SER:HA	1:B:563:LEU:HD12	1.56	0.88
1:A:708:THR:HG21	1:A:710:GLN:NE2	1.89	0.86
1:A:262:LEU:HD13	1:A:266:VAL:HB	1.59	0.85
1:A:510:VAL:HG11	1:A:541:ILE:HD13	1.59	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:521:LEU:H	1:B:521:LEU:HD12	1.41	0.84
1:A:401:ASN:OD1	1:B:338:THR:HB	1.79	0.83
1:A:417:VAL:HG22	1:A:628:THR:HG23	1.60	0.82
1:A:497:TRP:HZ3	1:B:482:GLU:HB2	1.45	0.80
1:B:36:ALA:HA	1:B:51:LEU:HD13	1.64	0.79
1:A:446:PHE:HD2	1:A:576:GLY:HA2	1.48	0.79
1:A:103:GLY:O	1:A:104:LEU:HG	1.83	0.79
1:A:36:ALA:HA	1:A:51:LEU:HD13	1.65	0.79
1:B:28:LYS:NZ	4:B:1810:TLA:H2	1.98	0.78
1:B:101:SER:O	1:B:102:LEU:HD23	1.83	0.78
1:B:428:SER:HB2	1:B:433:THR:OG1	1.84	0.78
1:A:21:SER:HB2	1:A:28:LYS:HD3	1.66	0.78
1:A:338:THR:HB	1:B:401:ASN:OD1	1.85	0.77
1:B:65:SER:HB3	1:B:68:ARG:HH12	1.47	0.77
1:A:325:PRO:O	1:A:327:THR:HG23	1.85	0.77
1:B:446:PHE:HD2	1:B:576:GLY:HA2	1.44	0.77
1:A:521:LEU:HD12	1:A:521:LEU:N	1.98	0.77
1:B:262:LEU:HD13	1:B:266:VAL:HB	1.68	0.76
1:B:405:PRO:HB3	1:B:689:ALA:HB2	1.67	0.76
1:A:520:SER:C	1:A:521:LEU:HD12	2.05	0.76
1:A:446:PHE:CE2	1:A:576:GLY:HA2	2.23	0.73
1:A:723:VAL:HG11	1:A:737:VAL:HG11	1.70	0.73
1:B:446:PHE:CE2	1:B:576:GLY:HA2	2.23	0.73
1:B:68:ARG:CB	1:B:68:ARG:HH11	2.00	0.73
1:A:428:SER:HB2	1:A:433:THR:OG1	1.87	0.73
1:B:21:SER:HB2	1:B:28:LYS:HD3	1.71	0.73
1:B:770:ASP:HB3	1:B:800:LYS:HD2	1.71	0.72
1:B:521:LEU:HD12	1:B:521:LEU:N	2.04	0.72
1:A:497:TRP:CZ3	1:B:482:GLU:HB2	2.23	0.72
1:B:608:ARG:HG3	1:B:608:ARG:HH11	1.54	0.72
1:A:373:GLY:H	1:A:374:PRO:CD	2.03	0.71
1:A:96:LEU:HB3	1:A:97:PRO:HD3	1.73	0.71
1:B:373:GLY:H	1:B:374:PRO:CD	2.04	0.71
1:A:770:ASP:HB3	1:A:800:LYS:HD2	1.71	0.71
1:A:726:ILE:HG12	1:A:802:TYR:HD1	1.55	0.71
1:A:208:ARG:NH1	1:A:211:ARG:HD3	2.06	0.70
1:A:497:TRP:HB3	1:B:501:ARG:HE	1.56	0.69
1:A:708:THR:HG23	1:A:709:SER:N	2.07	0.69
1:B:94:SER:O	1:B:113:MET:HE3	1.93	0.69
1:A:441:MET:HE1	1:A:658:ALA:HB2	1.74	0.69
3:A:1809:KAP:HA2	1:B:195:SER:OG	1.91	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:608:ARG:HH11	1:A:608:ARG:HG3	1.56	0.69
1:A:730:SER:O	1:A:747:LYS:HD2	1.93	0.69
1:B:8:PHE:CE2	1:B:361:GLN:HG3	2.27	0.68
1:A:8:PHE:CE2	1:A:361:GLN:HG3	2.29	0.68
1:B:208:ARG:NH1	1:B:211:ARG:HD3	2.06	0.68
1:A:8:PHE:CD2	1:A:361:GLN:HG3	2.28	0.68
1:B:485:ALA:HB2	1:B:504:PHE:CG	2.29	0.68
1:A:74:LEU:HD13	1:A:130:LEU:HD22	1.75	0.67
1:B:766:MET:HE2	1:B:804:ARG:HB3	1.76	0.67
1:A:477:THR:HG21	1:B:477:THR:HG21	1.77	0.67
1:A:405:PRO:HB3	1:A:689:ALA:HB2	1.76	0.67
1:B:8:PHE:CD2	1:B:361:GLN:HG3	2.30	0.66
1:A:659:THR:HG22	1:A:661:ALA:N	2.00	0.66
1:B:578:LEU:HD23	1:B:612:VAL:HG13	1.77	0.65
1:A:485:ALA:HB2	1:A:504:PHE:CG	2.32	0.65
1:A:101:SER:O	1:A:102:LEU:HD23	1.95	0.65
1:B:143:SER:HB3	1:B:145:HIS:ND1	2.10	0.65
1:A:13:ASN:ND2	1:A:356:ASN:HB3	2.12	0.65
1:B:551:LEU:HA	1:B:554:ILE:HD12	1.79	0.65
1:B:96:LEU:HB3	1:B:97:PRO:HD3	1.79	0.64
1:A:766:MET:HE2	1:A:804:ARG:HB3	1.80	0.64
1:A:8:PHE:C	1:A:8:PHE:CD1	2.71	0.64
1:A:521:LEU:N	1:A:521:LEU:HD13	2.13	0.64
1:A:80:ARG:HH12	1:A:287:VAL:HG21	1.62	0.64
1:A:375:ASP:HB2	1:A:378:PHE:H	1.63	0.64
1:A:9:HIS:ND1	1:A:360:SER:HB3	2.13	0.63
1:B:430:ASN:HB2	2:B:1644:PLP:O3P	1.98	0.63
1:B:484:GLN:HE22	1:B:489:TYR:HB2	1.63	0.63
1:A:580:ILE:O	1:A:582:PRO:HD3	1.97	0.63
1:B:62:PRO:HG3	1:B:102:LEU:HD21	1.79	0.63
1:B:474:HIS:O	1:B:480:ALA:HB1	1.99	0.62
1:A:191:GLY:H	4:A:1810:TLA:H3	1.64	0.62
1:B:61:PHE:O	1:B:139:GLU:HA	1.98	0.62
1:B:350:ILE:HD12	1:B:362:GLN:HB3	1.81	0.62
1:A:616:GLU:OE2	1:A:628:THR:HG22	2.00	0.62
1:A:546:ARG:NH2	1:A:596:LEU:HD23	2.15	0.61
1:B:692:ALA:O	1:B:696:GLN:HG3	2.00	0.61
1:A:367:ALA:HB1	1:A:372:GLN:HB2	1.82	0.61
1:A:321:VAL:HG13	1:B:414:LEU:CD1	2.30	0.61
1:B:373:GLY:H	1:B:374:PRO:HD3	1.65	0.61
1:B:580:ILE:O	1:B:582:PRO:HD3	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:718:TRP:CE3	1:A:794:CYS:HB3	2.34	0.61
1:A:484:GLN:HE22	1:A:489:TYR:HB2	1.66	0.61
1:B:651:VAL:HG22	1:B:652:PRO:HD2	1.82	0.61
1:B:441:MET:HE2	1:B:639:ILE:HG12	1.80	0.61
1:B:310:ARG:O	1:B:314:MET:HG3	2.01	0.61
1:B:8:PHE:CD1	1:B:8:PHE:C	2.73	0.60
1:B:601:LEU:O	1:B:601:LEU:HD12	2.00	0.60
1:A:252[A]:LEU:HD22	1:A:254:ASN:OD1	2.01	0.60
1:A:708:THR:HB	1:A:713:THR:O	2.02	0.60
1:B:143:SER:HB3	1:B:145:HIS:CE1	2.36	0.60
1:A:62:PRO:HG3	1:A:102:LEU:HD21	1.83	0.60
1:B:74:LEU:HD13	1:B:130:LEU:HD22	1.84	0.60
1:B:546:ARG:NH2	1:B:596:LEU:HD23	2.17	0.59
1:B:65:SER:HB3	1:B:68:ARG:HH11	1.66	0.59
1:B:539:ASP:C	1:B:541:ILE:H	2.04	0.59
1:A:692:ALA:O	1:A:696:GLN:HG3	2.02	0.59
1:B:28:LYS:HZ1	4:B:1810:TLA:H2	1.68	0.59
1:A:719:ASP:O	1:A:723:VAL:HG23	2.02	0.59
1:B:510:VAL:HG11	1:B:541:ILE:HD13	1.83	0.59
1:B:66:ASP:OD2	4:B:1810:TLA:O1	2.20	0.59
1:A:497:TRP:CZ3	1:B:482:GLU:CB	2.85	0.59
1:A:157:ASP:HB3	1:A:206:LEU:HD12	1.84	0.59
1:A:430:ASN:HB2	2:A:1644:PLP:O3P	2.03	0.59
1:A:310:ARG:O	1:A:314:MET:HG3	2.02	0.59
1:A:707:ILE:HD13	1:A:714:LEU:HD12	1.83	0.59
1:A:353:ALA:HA	1:A:357:SER:HA	1.85	0.59
1:B:659:THR:HG22	1:B:661:ALA:N	2.03	0.59
1:B:485:ALA:HB2	1:B:504:PHE:CD1	2.37	0.59
1:A:766:MET:HE3	1:A:807:GLU:OE1	2.02	0.58
1:A:68:ARG:HB3	1:A:68:ARG:HH11	1.66	0.58
1:A:220:GLY:C	1:A:252[A]:LEU:HD21	2.23	0.58
1:A:578:LEU:HD23	1:A:612:VAL:HG13	1.84	0.58
1:A:441:MET:HE2	1:A:639:ILE:HG12	1.84	0.58
1:A:140:ALA:HB1	1:A:142:ILE:HD13	1.85	0.58
4:A:1810:TLA:O41	3:B:1809:KAP:N	2.36	0.58
1:A:519:ILE:HG12	1:A:521:LEU:CD1	2.33	0.58
1:B:725:GLN:HB3	1:B:802:TYR:CE1	2.37	0.58
1:A:211:ARG:HD2	1:A:239:GLY:O	2.04	0.57
1:B:80:ARG:NH1	1:B:287:VAL:HG21	2.20	0.57
1:A:100:LYS:C	1:A:102:LEU:H	2.08	0.57
1:A:370:TRP:N	1:A:370:TRP:CD1	2.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:485:ALA:HB2	1:A:504:PHE:CD1	2.39	0.57
1:B:14:HIS:HB2	1:B:172:MET:CE	2.34	0.57
1:B:91:VAL:HG22	1:B:134:THR:HB	1.86	0.57
1:A:598:GLN:O	1:A:602:VAL:HG23	2.05	0.57
1:A:225:ILE:HD13	1:A:257:PRO:HG2	1.86	0.57
1:A:373:GLY:H	1:A:374:PRO:HD3	1.68	0.57
1:B:31:VAL:HG21	1:B:217:VAL:HG22	1.86	0.57
1:B:446:PHE:CD1	1:B:446:PHE:C	2.77	0.57
1:B:335[B]:GLU:OE1	1:B:335[B]:GLU:HA	2.05	0.57
1:B:208:ARG:HH12	1:B:211:ARG:CD	2.11	0.57
1:B:508:PRO:HA	1:B:520:SER:O	2.05	0.57
1:A:220:GLY:O	1:A:250:HIS:HB2	2.05	0.56
1:A:220:GLY:O	1:A:252[A]:LEU:CD2	2.47	0.56
1:A:508:PRO:HA	1:A:520:SER:O	2.06	0.56
1:B:651:VAL:CG2	1:B:652:PRO:HD2	2.36	0.56
1:A:473:TYR:OH	1:A:476:ASP:OD1	2.16	0.56
1:B:732:VAL:HG22	1:B:744:LEU:HD22	1.87	0.56
1:B:466:VAL:HG12	1:B:502:GLY:HA3	1.87	0.56
1:A:61:PHE:O	1:A:139:GLU:HA	2.05	0.56
1:A:601:LEU:O	1:A:601:LEU:HD12	2.05	0.56
1:B:417:VAL:HG23	1:B:628:THR:HG22	1.86	0.56
1:A:719:ASP:HB3	1:A:722:LEU:HD12	1.87	0.56
1:A:323:TRP:HB3	1:B:426:TYR:HD2	1.71	0.56
1:B:539:ASP:C	1:B:541:ILE:N	2.59	0.56
1:A:446:PHE:CD1	1:A:446:PHE:C	2.76	0.56
1:A:493:LEU:HB3	1:B:444:ARG:NH2	2.14	0.56
1:B:370:TRP:N	1:B:370:TRP:CD1	2.70	0.55
1:A:706:ASN:HA	1:A:715:ARG:HE	1.71	0.55
1:A:417:VAL:CG2	1:A:628:THR:HG23	2.33	0.55
1:B:23:ASN:OD1	1:B:24:THR:N	2.32	0.55
1:A:396:VAL:HG12	1:B:348:PHE:HE1	1.70	0.55
1:B:644:LYS:NZ	2:B:1644:PLP:O3	2.40	0.55
1:B:375:ASP:HB2	1:B:378:PHE:H	1.70	0.55
1:B:711:GLY:O	1:B:712:LYS:HB2	2.05	0.55
1:B:254:ASN:O	1:B:257:PRO:HD2	2.06	0.55
1:B:441:MET:HE1	1:B:658:ALA:HB2	1.88	0.55
1:B:100:LYS:C	1:B:102:LEU:H	2.10	0.55
3:A:1808:KAP:H2	1:B:679:HIS:HA	1.72	0.55
1:A:551:LEU:HA	1:A:554:ILE:HD12	1.89	0.55
1:A:723:VAL:CG1	1:A:737:VAL:HG11	2.36	0.55
1:B:432:SER:HB3	1:B:477:THR:OG1	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:414:LEU:CD1	1:B:321:VAL:HG13	2.35	0.55
1:A:512:LEU:HD23	1:A:727:SER:HB3	1.88	0.55
1:B:417:VAL:HG21	1:B:642:PHE:CE1	2.41	0.54
1:A:441:MET:CE	1:A:658:ALA:HB2	2.37	0.54
1:A:521:LEU:H	1:A:521:LEU:HD13	1.70	0.54
1:A:732:VAL:HG22	1:A:744:LEU:HD22	1.88	0.54
1:A:97:PRO:HG2	1:A:151:GLU:HB3	1.90	0.54
2:A:1644:PLP:C4A	3:A:1808:KAP:O	2.56	0.54
1:B:628:THR:OG1	1:B:630:THR:HB	2.07	0.54
1:A:805:LEU:C	1:A:807:GLU:H	2.11	0.54
1:A:546:ARG:HB3	1:A:596:LEU:HD22	1.88	0.54
1:A:490:THR:HB	1:A:498:TYR:CE2	2.43	0.54
1:B:746:LEU:HD11	1:B:781:ILE:HD12	1.90	0.54
1:A:708:THR:CG2	1:A:710:GLN:CD	2.76	0.54
1:B:506:ASP:HB3	1:B:522:PRO:HG2	1.89	0.54
1:B:208:ARG:NH1	1:B:211:ARG:CD	2.71	0.54
1:A:100:LYS:O	1:A:102:LEU:N	2.40	0.54
1:B:719:ASP:HB3	1:B:722:LEU:HB2	1.90	0.54
1:B:12:LEU:O	1:B:308:LEU:HD11	2.08	0.53
1:A:474:HIS:O	1:A:480:ALA:HB1	2.08	0.53
1:A:190:ALA:O	1:A:203:GLN:NE2	2.42	0.53
1:A:330:LYS:HD3	1:B:668:GLY:O	2.07	0.53
1:A:350:ILE:HD12	1:A:362:GLN:HB3	1.88	0.53
1:A:681:TYR:OH	1:B:649:GLY:HA2	2.09	0.53
1:A:14:HIS:HB2	1:A:172:MET:CE	2.38	0.53
1:B:625:GLY:HA2	1:B:714:LEU:HD12	1.91	0.53
1:A:746:LEU:HD11	1:A:781:ILE:HD12	1.90	0.53
1:B:473:TYR:OH	1:B:476:ASP:OD1	2.19	0.52
1:A:622:TRP:CE3	1:A:714:LEU:HD13	2.44	0.52
1:A:668:GLY:O	1:B:330:LYS:HD3	2.10	0.52
1:A:497:TRP:HE1	1:B:464:VAL:HG23	1.74	0.52
1:A:706:ASN:O	1:A:715:ARG:HG2	2.09	0.52
1:A:252[A]:LEU:HD22	1:A:254:ASN:CG	2.30	0.52
1:B:211:ARG:HD2	1:B:239:GLY:O	2.10	0.52
1:A:469:LEU:HD13	1:A:507:PRO:HB3	1.90	0.52
1:B:544:LYS:O	1:B:547:ASP:HB2	2.09	0.52
1:A:28:LYS:HZ2	4:A:1810:TLA:H2	1.75	0.52
1:B:405:PRO:HB3	1:B:689:ALA:CB	2.39	0.52
2:B:1644:PLP:C4A	3:B:1808:KAP:O	2.58	0.52
1:A:432:SER:HB3	1:A:477:THR:OG1	2.09	0.52
1:A:644:LYS:NZ	2:A:1644:PLP:O3	2.43	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:225:ILE:HD13	1:B:257:PRO:HG2	1.92	0.51
1:A:708:THR:HG21	1:A:710:GLN:CD	2.31	0.51
1:A:91:VAL:HG22	1:A:134:THR:HB	1.91	0.51
1:A:708:THR:CG2	1:A:709:SER:N	2.72	0.51
1:A:396:VAL:HG12	1:B:348:PHE:CE1	2.45	0.51
1:B:782:TYR:CE2	1:B:784:MET:HE3	2.46	0.51
1:B:487:SER:HB2	1:B:488:PRO:HD2	1.92	0.51
1:A:71:PHE:CD2	1:A:89:ASN:HB2	2.45	0.51
1:B:446:PHE:HD1	1:B:446:PHE:C	2.14	0.51
1:A:208:ARG:NH1	1:A:211:ARG:CD	2.72	0.51
1:B:720:GLU:O	1:B:723:VAL:HG22	2.11	0.51
1:B:71:PHE:CD2	1:B:89:ASN:HB2	2.46	0.51
1:A:501:ARG:HD2	1:B:497:TRP:CD2	2.46	0.51
1:A:777:LEU:HB2	1:A:780:VAL:HB	1.91	0.51
1:A:446:PHE:HD1	1:A:446:PHE:C	2.13	0.51
1:A:321:VAL:HG13	1:B:414:LEU:HD12	1.93	0.51
1:A:93:HIS:HB2	1:A:113:MET:HE2	1.92	0.51
1:B:298:GLU:O	1:B:302:LEU:HB2	2.09	0.51
1:A:514:ASN:HD22	1:A:728:SER:HA	1.75	0.51
1:A:68:ARG:NH1	1:A:68:ARG:HB3	2.25	0.51
1:A:256:VAL:HB	1:A:257:PRO:HD3	1.92	0.51
1:B:441:MET:CE	1:B:658:ALA:HB2	2.40	0.51
1:B:42:GLN:HA	1:B:42:GLN:OE1	2.10	0.51
1:B:563:LEU:HD13	1:B:608:ARG:NH1	2.26	0.50
1:B:85:ILE:HA	1:B:127:ALA:HB1	1.93	0.50
1:A:511:PHE:CE1	1:A:518:ASN:HB2	2.46	0.50
1:B:469:LEU:HD13	1:B:507:PRO:HB3	1.93	0.50
1:B:28:LYS:CE	4:B:1810:TLA:H2	2.40	0.50
1:B:605:CYS:HB3	1:B:610:ILE:HB	1.93	0.50
1:A:356:ASN:O	1:A:357:SER:HB2	2.12	0.50
1:B:80:ARG:HG3	1:B:80:ARG:HH11	1.77	0.50
1:A:482:GLU:HB3	1:A:501:ARG:HB3	1.94	0.50
1:B:608:ARG:CG	1:B:608:ARG:HH11	2.24	0.50
1:A:321:VAL:HG13	1:B:414:LEU:HD11	1.94	0.50
1:A:735:VAL:HG22	1:A:744:LEU:HB3	1.94	0.50
1:B:157:ASP:HB3	1:B:206:LEU:HD12	1.92	0.50
1:B:345:GLY:HA3	1:B:376:PRO:HD3	1.94	0.50
1:B:367:ALA:HB1	1:B:372:GLN:HB2	1.92	0.50
1:A:93:HIS:HA	1:A:114:CYS:O	2.12	0.49
1:B:777:LEU:HB2	1:B:780:VAL:HB	1.94	0.49
1:A:85:ILE:HA	1:A:127:ALA:HB1	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:252[A]:LEU:H	1:A:252[A]:LEU:CD1	2.04	0.49
1:B:633:LEU:HB3	1:B:635:CYS:SG	2.53	0.49
1:A:605:CYS:HB3	1:A:610:ILE:HB	1.94	0.49
1:B:563:LEU:HD11	1:B:604:GLU:HB3	1.94	0.49
1:A:373:GLY:HA2	1:A:786:GLY:HA3	1.94	0.49
1:A:805:LEU:O	1:A:807:GLU:N	2.37	0.49
1:B:190:ALA:O	1:B:203:GLN:NE2	2.43	0.49
1:A:766:MET:HB3	1:A:804:ARG:NE	2.27	0.49
1:B:438:ALA:HB1	1:B:613:ILE:HD13	1.93	0.49
1:B:390:ALA:HB2	1:B:686:MET:HE1	1.93	0.49
1:B:511:PHE:CE1	1:B:518:ASN:HB2	2.48	0.49
1:A:519:ILE:HG12	1:A:521:LEU:HD12	1.94	0.49
1:B:97:PRO:HG2	1:B:151:GLU:HB3	1.93	0.49
1:B:237:LEU:HA	1:B:237:LEU:HD12	1.57	0.49
1:A:220:GLY:O	1:A:252[A]:LEU:HD11	2.13	0.49
1:A:254:ASN:O	1:A:257:PRO:HD2	2.12	0.49
1:A:446:PHE:C	1:A:448:VAL:H	2.16	0.48
1:A:775:ARG:HG3	1:A:776:PRO:HD2	1.95	0.48
1:A:782:TYR:CE2	1:A:784:MET:HE3	2.48	0.48
1:A:407:LEU:HD11	1:B:317:LEU:HG	1.95	0.48
1:A:487:SER:HB2	1:A:488:PRO:HD2	1.94	0.48
1:B:559:LEU:HD12	1:B:600:VAL:HG12	1.96	0.48
1:B:316:LYS:HB2	1:B:334:GLN:OE1	2.12	0.48
1:A:463:VAL:O	1:A:574:HIS:HB2	2.13	0.48
1:B:50:LYS:HE3	1:B:131:LEU:HD12	1.95	0.48
1:B:594:ASP:OD1	1:B:596:LEU:HB3	2.13	0.48
1:B:735:VAL:HG22	1:B:744:LEU:HB3	1.95	0.48
1:B:598:GLN:O	1:B:602:VAL:HG23	2.13	0.48
1:B:420:GLY:O	1:B:421:TRP:HB3	2.13	0.48
1:B:446:PHE:C	1:B:448:VAL:H	2.16	0.48
1:B:473:TYR:C	1:B:474:HIS:CD2	2.87	0.48
1:B:446:PHE:HD2	1:B:576:GLY:CA	2.21	0.48
1:B:775:ARG:NE	3:B:1808:KAP:OI1	2.32	0.48
1:B:465:LYS:NZ	1:B:573:ALA:HB1	2.28	0.48
1:B:312:ASN:ND2	1:B:351:TYR:OH	2.46	0.48
1:A:21:SER:CB	1:A:28:LYS:HD3	2.41	0.48
1:B:498:TYR:CE1	1:B:500:GLY:N	2.82	0.48
1:A:708:THR:CG2	1:A:710:GLN:NE2	2.72	0.48
1:B:546:ARG:HB3	1:B:596:LEU:HD22	1.95	0.48
1:B:625:GLY:HA2	1:B:714:LEU:HB2	1.95	0.48
1:A:160:VAL:O	1:A:164:ILE:HG13	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:400:GLU:O	1:B:337:VAL:HA	2.13	0.47
1:A:23:ASN:OD1	1:A:24:THR:N	2.42	0.47
1:B:208:ARG:HH11	1:B:211:ARG:HG3	1.79	0.47
1:A:608:ARG:CG	1:A:608:ARG:HH11	2.26	0.47
1:B:560:SER:CA	1:B:563:LEU:HD12	2.36	0.47
1:A:248:GLU:O	1:A:250:HIS:ND1	2.43	0.47
1:B:14:HIS:HB2	1:B:172:MET:HE3	1.95	0.47
1:A:323:TRP:CG	1:B:426:TYR:HD2	2.33	0.47
1:A:14:HIS:HB2	1:A:172:MET:HE3	1.96	0.47
1:B:614:PHE:O	1:B:640:ALA:HA	2.15	0.47
1:B:482:GLU:OE2	1:B:501:ARG:NH1	2.47	0.47
1:A:691:ALA:O	1:A:695:ILE:HG13	2.13	0.47
1:B:68:ARG:NH1	1:B:68:ARG:HB3	2.16	0.47
1:B:94:SER:HB3	1:B:116:LEU:HD11	1.97	0.47
1:A:140:ALA:O	1:A:142:ILE:N	2.43	0.47
1:B:335[B]:GLU:OE1	1:B:335[B]:GLU:CA	2.63	0.47
1:A:414:LEU:HD11	1:B:321:VAL:HG13	1.96	0.47
1:A:633:LEU:HB3	1:A:635:CYS:SG	2.54	0.47
1:A:73:LYS:HA	1:A:73:LYS:HD3	1.66	0.47
1:A:806:GLY:O	1:A:807:GLU:CG	2.63	0.47
1:B:628:THR:HG1	1:B:630:THR:HB	1.79	0.47
1:A:649:GLY:HA2	1:B:681:TYR:OH	2.15	0.47
1:A:555:TYR:O	1:A:559:LEU:HG	2.14	0.47
1:A:343:ARG:NH1	1:B:395:HIS:O	2.46	0.47
1:A:143:SER:HG	1:A:145:HIS:CE1	2.23	0.47
1:A:781:ILE:O	1:A:781:ILE:HG22	2.15	0.47
1:A:345:GLY:HA3	1:A:376:PRO:HD3	1.97	0.47
1:A:473:TYR:C	1:A:474:HIS:CD2	2.88	0.46
1:B:659:THR:HB	1:B:662:VAL:HG23	1.96	0.46
1:B:403:TYR:HD2	1:B:405:PRO:HD2	1.80	0.46
1:A:722:LEU:HD13	1:A:798:LEU:HB2	1.97	0.46
1:B:489:TYR:CE2	1:B:587:ALA:HB1	2.50	0.46
1:B:463:VAL:O	1:B:574:HIS:HB2	2.15	0.46
1:A:318:ALA:HB2	1:B:402:VAL:HB	1.98	0.46
1:B:208:ARG:O	1:B:208:ARG:HD2	2.16	0.46
1:B:775:ARG:HH21	3:B:1808:KAP:C	2.28	0.46
1:A:316:LYS:HB2	1:A:334:GLN:OE1	2.15	0.46
1:A:219:ASP:O	1:A:224:GLY:HA3	2.15	0.46
1:A:806:GLY:O	1:A:807:GLU:HG2	2.16	0.46
1:A:411:GLU:OE2	1:B:317:LEU:HD21	2.16	0.46
1:A:35:ILE:HD11	1:A:296:LEU:CD2	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:766:MET:HB3	1:A:804:ARG:CD	2.45	0.46
1:A:444:ARG:HH21	1:B:493:LEU:HB3	1.81	0.46
1:A:444:ARG:HG3	1:A:666:PHE:CZ	2.50	0.46
1:A:12:LEU:O	1:A:308:LEU:HD11	2.15	0.46
1:A:55:LYS:HE3	1:A:57:ILE:O	2.16	0.46
1:B:766:MET:HB3	1:B:804:ARG:CD	2.46	0.46
1:B:220:GLY:O	1:B:250:HIS:HB2	2.16	0.46
1:B:23:ASN:CG	1:B:24:THR:H	2.18	0.46
1:A:34:GLY:HA3	1:A:293:PHE:CE1	2.51	0.46
1:A:274:VAL:HG13	1:A:275:PRO:HD2	1.98	0.46
1:A:651:VAL:HG22	1:A:652:PRO:HD2	1.98	0.46
1:A:317:LEU:HG	1:B:407:LEU:HD11	1.97	0.46
1:B:766:MET:HB3	1:B:804:ARG:HD2	1.98	0.46
1:A:62:PRO:HG3	1:A:102:LEU:CD2	2.45	0.46
1:B:369:TRP:CE2	3:B:1808:KAP:HS2	2.50	0.46
1:B:764:LEU:HD22	1:B:774:THR:HG23	1.98	0.46
1:A:765:ILE:HG22	1:A:766:MET:N	2.30	0.45
1:A:298:GLU:O	1:A:302:LEU:HB2	2.16	0.45
1:A:208:ARG:HH12	1:A:211:ARG:CD	2.16	0.45
1:A:794:CYS:O	1:A:797:LEU:HB2	2.16	0.45
1:A:205:ASP:OD1	1:A:238:ARG:NE	2.39	0.45
1:A:446:PHE:CE1	1:A:450:HIS:ND1	2.84	0.45
1:A:766:MET:HE2	1:A:804:ARG:HD2	1.98	0.45
1:A:708:THR:HG23	1:A:709:SER:H	1.79	0.45
1:B:633:LEU:C	1:B:635:CYS:H	2.20	0.45
1:A:100:LYS:C	1:A:102:LEU:N	2.69	0.45
1:B:17:TYR:HD2	1:B:214:GLY:HA2	1.80	0.45
1:A:31:VAL:HG21	1:A:217:VAL:HG22	1.97	0.45
1:A:722:LEU:O	1:A:725:GLN:HB2	2.17	0.45
1:A:367:ALA:CB	1:A:372:GLN:HB2	2.46	0.45
1:B:446:PHE:CE1	1:B:450:HIS:ND1	2.85	0.45
1:B:539:ASP:N	1:B:539:ASP:OD1	2.49	0.45
1:B:539:ASP:O	1:B:541:ILE:N	2.50	0.45
1:A:323:TRP:HB3	1:B:426:TYR:CD2	2.51	0.45
1:B:10:LEU:HA	1:B:11:PRO:HD3	1.78	0.45
1:A:495:GLN:HG2	1:B:440:LYS:HE2	1.99	0.45
1:A:17:TYR:HD2	1:A:214:GLY:HA2	1.81	0.45
1:B:555:TYR:O	1:B:559:LEU:HG	2.17	0.45
1:B:521:LEU:CD1	1:B:521:LEU:N	2.72	0.45
1:B:100:LYS:O	1:B:102:LEU:N	2.50	0.45
1:B:62:PRO:HG3	1:B:102:LEU:CD2	2.45	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:473:TYR:C	1:A:474:HIS:CG	2.90	0.45
1:A:113:MET:HE2	1:A:136:TYR:HD1	1.82	0.45
1:A:614:PHE:O	1:A:640:ALA:HA	2.17	0.45
1:B:55:LYS:HE3	1:B:57:ILE:O	2.17	0.45
1:B:413:LEU:HD13	1:B:655:VAL:HG11	1.99	0.45
1:A:9:HIS:CE1	1:A:360:SER:HB3	2.52	0.45
1:B:274:VAL:HG13	1:B:275:PRO:HD2	1.99	0.45
1:B:521:LEU:H	1:B:521:LEU:CD1	2.20	0.44
1:A:466:VAL:HG12	1:A:502:GLY:HA3	1.99	0.44
1:A:426:TYR:HD2	1:B:323:TRP:HB3	1.81	0.44
3:A:1808:KAP:N	1:B:678:GLY:O	2.50	0.44
1:B:193:VAL:HG22	1:B:235:LEU:HD21	1.98	0.44
1:A:438:ALA:HB1	1:A:613:ILE:HD13	1.99	0.44
1:A:651:VAL:CG2	1:A:652:PRO:HD2	2.48	0.44
1:B:101:SER:C	1:B:102:LEU:HD23	2.38	0.44
1:B:473:TYR:C	1:B:474:HIS:CG	2.90	0.44
1:B:297:LYS:HB3	1:B:297:LYS:HE2	1.60	0.44
1:A:28:LYS:NZ	4:A:1810:TLA:H2	2.32	0.44
1:B:446:PHE:HD1	1:B:446:PHE:O	2.01	0.44
1:B:100:LYS:C	1:B:102:LEU:N	2.71	0.44
1:A:348:PHE:HE1	1:B:396:VAL:HG12	1.83	0.44
1:A:414:LEU:HD12	1:B:321:VAL:HG13	1.98	0.44
1:B:71:PHE:CE2	1:B:87:ILE:HG13	2.53	0.44
1:B:764:LEU:HD11	1:B:776:PRO:HB3	1.99	0.44
1:A:420:GLY:O	1:A:421:TRP:HB3	2.18	0.44
1:B:794:CYS:O	1:B:797:LEU:HB2	2.17	0.44
1:B:168:LEU:HA	1:B:168:LEU:HD23	1.73	0.44
1:A:446:PHE:HD1	1:A:446:PHE:O	1.99	0.44
1:A:626:VAL:CG2	1:A:628:THR:O	2.65	0.44
1:B:707:ILE:HG22	1:B:712:LYS:HA	2.00	0.44
1:B:356:ASN:O	1:B:357:SER:C	2.55	0.44
1:A:13:ASN:ND2	1:A:356:ASN:O	2.46	0.43
1:A:805:LEU:C	1:A:807:GLU:N	2.71	0.43
1:B:20:TRP:CE3	1:B:189:THR:HG21	2.53	0.43
1:A:402:VAL:HB	1:B:318:ALA:HB2	1.99	0.43
1:B:373:GLY:N	1:B:374:PRO:CD	2.77	0.43
1:B:370:TRP:CH2	3:B:1808:KAP:HE2	2.53	0.43
1:B:559:LEU:HD13	1:B:601:LEU:HA	2.00	0.43
1:B:417:VAL:HA	1:B:628:THR:CG2	2.49	0.43
1:A:94:SER:HB3	1:A:116:LEU:HD11	2.00	0.43
1:B:691:ALA:O	1:B:695:ILE:HG13	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:666:PHE:CE2	1:B:675:LEU:HB2	2.53	0.43
1:A:616:GLU:CD	1:A:628:THR:HG22	2.39	0.43
1:B:766:MET:HB3	1:B:804:ARG:NE	2.32	0.43
1:A:766:MET:HB3	1:A:804:ARG:HD2	2.00	0.43
1:A:744:LEU:C	1:A:744:LEU:HD12	2.38	0.43
1:A:446:PHE:CD1	1:A:450:HIS:ND1	2.87	0.43
1:B:486:PRO:HD3	1:B:498:TYR:OH	2.19	0.43
1:B:725:GLN:HB3	1:B:802:TYR:CZ	2.53	0.43
1:B:444:ARG:HG3	1:B:666:PHE:CZ	2.53	0.43
1:A:719:ASP:OD2	1:A:722:LEU:HD12	2.17	0.43
1:B:12:LEU:HB3	1:B:308:LEU:HD21	2.00	0.43
1:A:276:LYS:HD3	1:A:276:LYS:HA	1.68	0.43
1:B:608:ARG:CG	1:B:608:ARG:NH1	2.81	0.43
1:A:519:ILE:HG12	1:A:521:LEU:HD11	1.99	0.43
1:A:221:ARG:HE	1:A:221:ARG:HB2	1.65	0.43
1:A:324:TRP:HB2	1:A:327:THR:OG1	2.18	0.43
1:A:608:ARG:NH1	1:A:608:ARG:HG3	2.30	0.43
1:B:378:PHE:CE1	1:B:382:LEU:HG	2.54	0.43
1:A:764:LEU:HD13	1:A:776:PRO:HD3	2.00	0.43
1:A:297:LYS:HE2	1:A:297:LYS:HB3	1.54	0.43
1:A:446:PHE:O	1:A:448:VAL:N	2.50	0.43
1:A:766:MET:CE	1:A:804:ARG:HD2	2.49	0.43
1:A:261:TYR:CD1	1:A:261:TYR:C	2.91	0.43
1:A:50:LYS:HE3	1:A:131:LEU:HD12	2.00	0.43
1:A:628:THR:HB	1:A:630:THR:H	1.83	0.43
1:B:550:THR:O	1:B:554:ILE:HG13	2.19	0.43
1:A:8:PHE:C	1:A:8:PHE:HD1	2.21	0.42
1:A:766:MET:HE1	1:A:807:GLU:HB2	2.01	0.42
1:A:382:LEU:HA	1:A:382:LEU:HD23	1.92	0.42
1:B:706:ASN:HB3	1:B:714:LEU:CD2	2.48	0.42
1:A:403:TYR:HD2	1:A:405:PRO:HD2	1.84	0.42
1:A:378:PHE:CE1	1:A:382:LEU:HG	2.55	0.42
1:B:538:ARG:O	1:B:541:ILE:HG13	2.19	0.42
1:B:513:SER:O	1:B:515:GLY:N	2.52	0.42
1:B:373:GLY:HA2	1:B:786:GLY:HA3	2.00	0.42
1:A:369:TRP:CH2	1:B:399:PRO:HG3	2.55	0.42
1:B:219:ASP:O	1:B:224:GLY:HA3	2.19	0.42
1:B:486:PRO:HA	1:B:490:THR:OG1	2.19	0.42
1:A:103:GLY:O	1:A:104:LEU:CG	2.61	0.42
1:A:629:THR:C	1:A:631:GLU:N	2.71	0.42
1:A:659:THR:HB	1:A:662:VAL:HG23	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:SER:CB	1:B:145:HIS:ND1	2.80	0.42
1:B:697:TRP:CE2	1:B:703:THR:HB	2.55	0.42
1:A:770:ASP:CB	1:A:800:LYS:HD2	2.45	0.42
1:B:431:GLY:HA3	2:B:1644:PLP:O4P	2.20	0.42
1:B:625:GLY:HA2	1:B:714:LEU:CB	2.50	0.42
1:A:697:TRP:CE2	1:A:703:THR:HB	2.55	0.42
1:A:764:LEU:HD22	1:A:774:THR:HG23	2.02	0.42
1:A:782:TYR:CD2	1:A:784:MET:HE3	2.55	0.42
1:B:700:ASP:HA	1:B:701:PRO:HD2	1.79	0.42
1:A:219:ASP:HB3	1:A:224:GLY:HA2	2.02	0.42
1:A:707:ILE:HA	1:A:707:ILE:HD13	1.78	0.42
1:B:417:VAL:HA	1:B:628:THR:HG22	2.01	0.42
1:A:237:LEU:HD11	1:B:237:LEU:HD11	2.02	0.42
1:B:494:GLN:O	1:B:495:GLN:HB2	2.20	0.42
1:B:435:ILE:HG13	1:B:474:HIS:HB3	2.02	0.41
1:A:489:TYR:CE2	1:A:587:ALA:HB1	2.55	0.41
1:A:501:ARG:CD	1:B:497:TRP:CD2	3.03	0.41
1:A:94:SER:O	1:A:113:MET:HE3	2.20	0.41
1:B:65:SER:CB	1:B:68:ARG:HH12	2.25	0.41
1:A:225:ILE:HB	3:A:1809:KAP:OI2	2.20	0.41
1:B:446:PHE:CD1	1:B:450:HIS:ND1	2.88	0.41
1:B:369:TRP:CD2	3:B:1808:KAP:HS2	2.56	0.41
1:A:140:ALA:CB	1:A:142:ILE:HD13	2.48	0.41
1:A:281:ASP:O	1:A:282:LEU:HB2	2.20	0.41
1:B:643:ALA:O	1:B:645:LEU:N	2.52	0.41
1:A:594:ASP:OD1	1:A:596:LEU:HB3	2.20	0.41
1:A:435:ILE:HG13	1:A:474:HIS:HB3	2.03	0.41
1:B:219:ASP:HB3	1:B:224:GLY:HA2	2.02	0.41
1:A:627:GLU:OE2	1:A:699:LYS:HE2	2.19	0.41
1:A:321:VAL:O	1:A:321:VAL:HG12	2.19	0.41
1:B:183:LEU:C	1:B:183:LEU:HD23	2.40	0.41
1:A:484:GLN:NE2	1:A:489:TYR:HB2	2.34	0.41
1:A:316:LYS:O	1:A:320:GLU:HG3	2.20	0.41
1:B:92:LEU:HD12	1:B:135:LEU:CD2	2.51	0.41
1:A:208:ARG:HH11	1:A:211:ARG:HG3	1.84	0.41
1:B:431:GLY:O	1:B:434:ALA:HB3	2.20	0.41
1:B:256:VAL:HB	1:B:257:PRO:HD3	2.03	0.41
1:B:264:ASN:O	1:B:264:ASN:OD1	2.38	0.41
1:A:96:LEU:N	1:A:97:PRO:CD	2.84	0.41
1:B:744:LEU:C	1:B:744:LEU:HD12	2.41	0.41
1:A:501:ARG:CD	1:B:497:TRP:CE2	3.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:56:PRO:O	1:B:57:ILE:HD13	2.21	0.41
1:A:426:TYR:HD2	1:B:323:TRP:CG	2.39	0.41
1:A:676:LEU:HG	1:B:494:GLN:NE2	2.34	0.41
1:A:642:PHE:HB2	1:A:655:VAL:HG12	2.03	0.41
1:A:425:VAL:HG13	1:A:655:VAL:HG22	2.02	0.41
1:A:719:ASP:CB	1:A:722:LEU:HD12	2.49	0.41
1:B:594:ASP:HA	1:B:595:PRO:HD2	1.95	0.41
1:B:722:LEU:HD23	1:B:722:LEU:HA	1.73	0.41
1:A:446:PHE:HD2	1:A:576:GLY:CA	2.23	0.41
1:B:490:THR:HB	1:B:498:TYR:CE2	2.56	0.41
1:B:36:ALA:HA	1:B:51:LEU:CD1	2.42	0.41
1:B:333:HIS:HB3	1:B:335[A]:GLU:OE2	2.21	0.41
1:B:764:LEU:HD13	1:B:776:PRO:HD3	2.03	0.41
1:B:246:VAL:HA	1:B:269:LEU:O	2.21	0.41
1:B:624:LEU:HD22	1:B:738:ILE:HG23	2.03	0.41
1:B:73:LYS:HA	1:B:73:LYS:HD3	1.68	0.41
1:B:773:PHE:C	1:B:773:PHE:CD1	2.93	0.41
1:B:396:VAL:HG23	1:B:403:TYR:CE1	2.56	0.41
1:A:608:ARG:CG	1:A:608:ARG:NH1	2.83	0.41
1:B:584:ILE:HD13	1:B:592:MET:HA	2.02	0.41
1:A:395:HIS:HB3	1:B:649:GLY:O	2.21	0.41
1:A:624:LEU:C	1:A:739:GLY:HA3	2.41	0.41
1:B:205:ASP:OD1	1:B:238:ARG:NE	2.43	0.41
1:A:56:PRO:O	1:A:57:ILE:HD13	2.21	0.40
1:A:324:TRP:H	1:A:329:HIS:CE1	2.39	0.40
1:B:582:PRO:O	1:B:584:ILE:N	2.47	0.40
1:A:34:GLY:HA3	1:A:293:PHE:CZ	2.55	0.40
1:B:276:LYS:HA	1:B:276:LYS:HD3	1.81	0.40
1:A:348:PHE:CE1	1:B:396:VAL:HG12	2.56	0.40
1:A:168:LEU:HA	1:A:168:LEU:HD23	1.86	0.40
1:A:712:LYS:HB2	1:A:712:LYS:HE3	1.95	0.40
1:A:246:VAL:HA	1:A:269:LEU:O	2.21	0.40
1:B:484:GLN:NE2	1:B:489:TYR:HB2	2.32	0.40
1:A:714:LEU:HD12	1:A:714:LEU:HA	1.92	0.40
1:A:473:TYR:C	1:A:473:TYR:CD1	2.95	0.40
1:B:382:LEU:HD13	1:B:650:MET:HG3	2.04	0.40
1:B:614:PHE:CD2	1:B:637:PRO:HG3	2.56	0.40
1:A:559:LEU:HD12	1:A:600:VAL:HG12	2.02	0.40
1:B:146:LEU:HA	1:B:146:LEU:HD23	1.83	0.40
1:A:74:LEU:HD13	1:A:130:LEU:CD2	2.48	0.40
1:A:435:ILE:HD11	1:A:615:ASP:CG	2.42	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:281:ASP:O	1:B:282:LEU:HB2	2.22	0.40
1:A:236:LYS:HA	1:A:236:LYS:HD2	1.86	0.40
1:A:783:LEU:HD11	1:A:801:LEU:HD22	2.04	0.40
1:B:482:GLU:O	1:B:500:GLY:HA2	2.21	0.40
1:A:13:ASN:HD22	1:A:356:ASN:HB3	1.86	0.40
1:B:427:PHE:CD2	1:B:683:ALA:HB2	2.57	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:149:GLU:OE2	1:B:711:GLY:N[2_555]	2.14	0.06

## 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	735/831 (88%)	659 (90%)	62 (8%)	14 (2%)	10	32
1	B	723/831 (87%)	643 (89%)	59 (8%)	21 (3%)	6	19
All	All	1458/1662 (88%)	1302 (89%)	121 (8%)	35 (2%)	7	25

All (35) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	514	ASN
1	B	514	ASN
1	B	563	LEU
1	A	101	SER
1	A	103	GLY
1	A	141	ALA
1	A	421	TRP

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Mol	Chain	Res	Type
1	A	451	ASN
1	A	644	LYS
1	B	101	SER
1	B	118	PHE
1	B	355	ASP
1	B	357	SER
1	B	421	TRP
1	B	447	CYS
1	B	451	ASN
1	A	355	ASP
1	A	373	GLY
1	A	643	ALA
1	A	778	GLY
1	B	373	GLY
1	B	644	LYS
1	A	447	CYS
1	B	354	SER
1	B	495	GLN
1	B	540	GLU
1	B	643	ALA
1	B	778	GLY
1	B	730	SER
1	A	806	GLY
1	B	486	PRO
1	B	500	GLY
1	B	637	PRO
1	A	637	PRO
1	B	583	VAL

### 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	615/704 (87%)	571 (93%)	44 (7%)	18	45
1	B	598/704 (85%)	558 (93%)	40 (7%)	20	50
All	All	1213/1408 (86%)	1129 (93%)	84 (7%)	20	48

All (84) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	PHE
1	A	13	ASN
1	A	24	THR
1	A	42	GLN
1	A	68	ARG
1	A	72	SER
1	A	74	LEU
1	A	80	ARG
1	A	107	GLU
1	A	131	LEU
1	A	142	ILE
1	A	158	SER
1	A	193	VAL
1	A	206	LEU
1	A	208	ARG
1	A	252[A]	LEU
1	A	252[B]	LEU
1	A	263	ARG
1	A	279	SER
1	A	343	ARG
1	A	355	ASP
1	A	370	TRP
1	A	417	VAL
1	A	446	PHE
1	A	451	ASN
1	A	477	THR
1	A	521	LEU
1	A	524	SER
1	A	540	GLU
1	A	543	ASP
1	A	558	TYR
1	A	560	SER
1	A	628	THR
1	A	630	THR
1	A	635	CYS
1	A	655	VAL
1	A	708	THR
1	A	709	SER
1	A	713	THR
1	A	714	LEU
1	A	738	ILE
1	A	744	LEU

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Mol	Chain	Res	Type
1	A	765	ILE
1	A	775	ARG
1	B	8	PHE
1	B	9	HIS
1	B	13	ASN
1	B	24	THR
1	B	63	SER
1	B	68	ARG
1	B	72	SER
1	B	74	LEU
1	B	80	ARG
1	B	131	LEU
1	B	142	ILE
1	B	158	SER
1	B	193	VAL
1	B	206	LEU
1	B	208	ARG
1	B	279	SER
1	B	302	LEU
1	B	335[A]	GLU
1	B	335[B]	GLU
1	B	343	ARG
1	B	370	TRP
1	B	417	VAL
1	B	446	PHE
1	B	451	ASN
1	B	463	VAL
1	B	477	THR
1	B	499	THR
1	B	501	ARG
1	B	521	LEU
1	B	539	ASP
1	B	560	SER
1	B	628	THR
1	B	630	THR
1	B	635	CYS
1	B	655	VAL
1	B	709	SER
1	B	714	LEU
1	B	730	SER
1	B	738	ILE
1	B	775	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	82	GLN
1	A	162	GLN
1	A	312	ASN
1	A	329	HIS
1	A	484	GLN
1	A	514	ASN
1	A	518	ASN
1	A	724	GLN
1	A	733	GLN
1	B	162	GLN
1	B	312	ASN
1	B	484	GLN
1	B	518	ASN
1	B	591	HIS
1	B	725	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](#)

8 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
2	PLP	A	1644	1	15,15,16	1.83	1 (6%)	21,22,23	2.28	3 (14%)
3	KAP	A	1808	-	8,12,12	0.65	0	9,14,14	1.11	0
3	KAP	A	1809	-	8,12,12	1.07	1 (12%)	9,14,14	2.42	2 (22%)
4	TLA	A	1810	-	3,9,9	2.02	1 (33%)	6,12,12	3.03	2 (33%)
2	PLP	B	1644	1	15,15,16	1.67	2 (13%)	21,22,23	2.42	2 (9%)
3	KAP	B	1808	-	8,12,12	0.66	0	9,14,14	0.88	0
3	KAP	B	1809	-	8,12,12	0.72	0	9,14,14	1.63	2 (22%)
4	TLA	B	1810	-	3,9,9	2.21	2 (66%)	6,12,12	3.71	3 (50%)

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	PLP	A	1644	1	-	0/6/6/8	0/1/1/1
3	KAP	A	1808	-	-	0/8/12/12	0/0/0/0
3	KAP	A	1809	-	-	0/8/12/12	0/0/0/0
4	TLA	A	1810	-	-	0/4/12/12	0/0/0/0
2	PLP	B	1644	1	-	0/6/6/8	0/1/1/1
3	KAP	B	1808	-	-	0/8/12/12	0/0/0/0
3	KAP	B	1809	-	-	0/8/12/12	0/0/0/0
4	TLA	B	1810	-	-	0/4/12/12	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1644	PLP	O3-C3	-5.62	1.23	1.37
2	B	1644	PLP	O3-C3	-5.04	1.25	1.37
4	A	1810	TLA	O3-C3	-2.80	1.37	1.42
4	B	1810	TLA	C3-C2	-2.74	1.44	1.53
4	B	1810	TLA	O3-C3	-2.01	1.38	1.42
2	B	1644	PLP	C2-N1	2.28	1.38	1.34
3	A	1809	KAP	CE-CZ	2.79	1.55	1.51

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1810	TLA	C1-C2-C3	-7.41	98.17	113.35
4	A	1810	TLA	C1-C2-C3	-5.07	102.95	113.35
4	A	1810	TLA	C4-C3-C2	-4.91	103.28	113.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1810	TLA	C4-C3-C2	-4.32	104.49	113.35
3	A	1809	KAP	CZ-CH-N	-2.79	102.10	109.90
2	B	1644	PLP	O4P-P-O1P	-2.52	100.72	107.14
2	A	1644	PLP	O4P-P-O1P	-2.09	101.82	107.14
3	B	1809	KAP	O-CZ-CE	-2.02	117.60	121.73
2	A	1644	PLP	O3P-P-O1P	2.10	117.35	110.58
4	B	1810	TLA	O2-C2-C1	2.37	117.20	111.21
3	B	1809	KAP	CE-CZ-CH	2.85	122.79	116.79
3	A	1809	KAP	CD-CE-CZ	5.74	125.22	114.63
2	A	1644	PLP	O4P-C5A-C5	9.03	123.92	108.99
2	B	1644	PLP	O4P-C5A-C5	9.37	124.48	108.99

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1644	PLP	3	0
3	A	1808	KAP	3	0
3	A	1809	KAP	2	0
4	A	1810	TLA	5	0
2	B	1644	PLP	4	0
3	B	1808	KAP	6	0
3	B	1809	KAP	1	0
4	B	1810	TLA	4	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	748/831 (90%)	-0.09	20 (2%)	58 45	35, 62, 96, 134	0
1	B	738/831 (88%)	-0.09	25 (3%)	49 36	35, 64, 99, 127	0
All	All	1486/1662 (89%)	-0.09	45 (3%)	54 41	35, 63, 98, 134	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	428	SER	4.5
1	B	576	GLY	4.1
1	B	429	ASP	3.8
1	A	355	ASP	3.8
1	A	576	GLY	3.5
1	A	449	ASP	3.1
1	B	610	ILE	3.0
1	A	354	SER	2.9
1	B	781	ILE	2.8
1	A	432	SER	2.8
1	A	688	CYS	2.7
1	B	575	VAL	2.7
1	B	802	TYR	2.7
1	B	652	PRO	2.6
1	B	450	HIS	2.6
1	B	430	ASN	2.6
1	A	668	GLY	2.5
1	A	430	ASN	2.5
1	A	428	SER	2.5
1	B	373	GLY	2.5
1	B	449	ASP	2.4
1	A	448	VAL	2.4
1	B	727	SER	2.4
1	A	126	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	446	PHE	2.4
1	B	463	VAL	2.4
1	A	172	MET	2.4
1	A	805	LEU	2.3
1	A	687	GLY	2.3
1	B	433	THR	2.3
1	A	450	HIS	2.2
1	B	518	ASN	2.2
1	B	679	HIS	2.2
1	B	555	TYR	2.2
1	A	684	HIS	2.2
1	A	652	PRO	2.1
1	A	691	ALA	2.1
1	A	667	SER	2.1
1	A	654	ALA	2.1
1	B	432	SER	2.1
1	B	682	SER	2.1
1	B	641	CYS	2.1
1	B	559	LEU	2.0
1	B	535	PHE	2.0
1	B	654	ALA	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	KAP	B	1808	13/13	0.90	0.33	2.38	46,61,84,92	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	KAP	A	1809	13/13	0.92	0.21	2.14	30,52,71,84	0
3	KAP	A	1808	13/13	0.89	0.32	1.65	40,56,81,85	0
3	KAP	B	1809	13/13	0.96	0.18	1.59	28,46,73,78	0
4	TLA	A	1810	10/10	0.95	0.14	0.23	49,57,76,92	0
2	PLP	A	1644	15/16	0.94	0.27	-0.03	32,52,71,74	0
2	PLP	B	1644	15/16	0.97	0.24	-0.11	41,49,72,80	0
4	TLA	B	1810	10/10	0.96	0.14	-0.33	28,55,72,85	10

## 6.5 Other polymers [i](#)

There are no such residues in this entry.