



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:32 PM GMT

PDB ID : 1B7Y
Title : PHENYLALANYL TRNA SYNTHETASE COMPLEXED WITH PHENYL
ALANINYL-ADENYLATE
Authors : Reshetnikova, L.; Moor, N.; Lavrik, O.; Vassilyev, D.G.
Deposited on : 1999-01-26
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

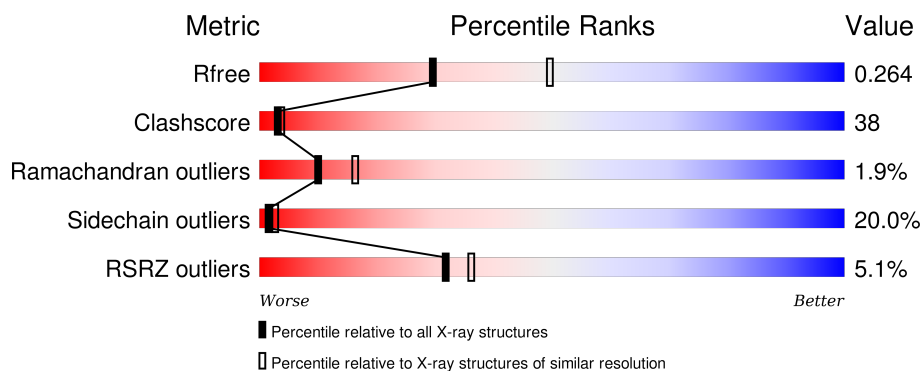
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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 ≥ 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	350	 2% 33% 34% 9% 24%
2	B	785	 6% 43% 43% 12% ..

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	A	1001	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8439 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 PROTEIN (PHENYLALANYL-TRNA SYNTHETASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	0	0
			2112	1382	359	364	7			

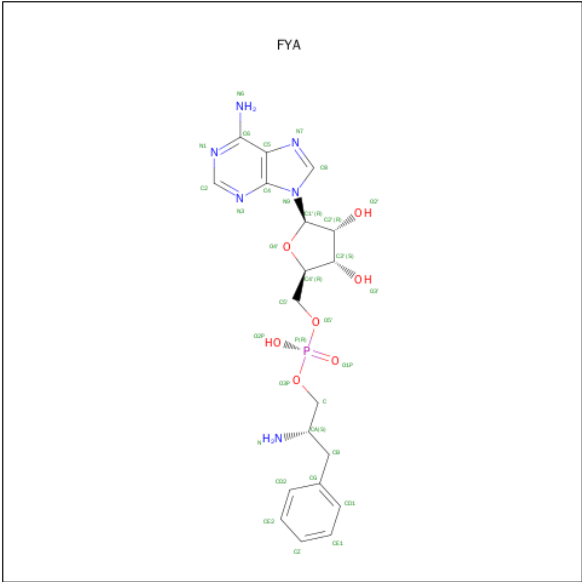
- Molecule 2 is a protein called PROTEIN (PHENYLALANYL-TRNA SYNTHETASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	775	Total	C	N	O	S	0	0	0
			6054	3879	1078	1087	10			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

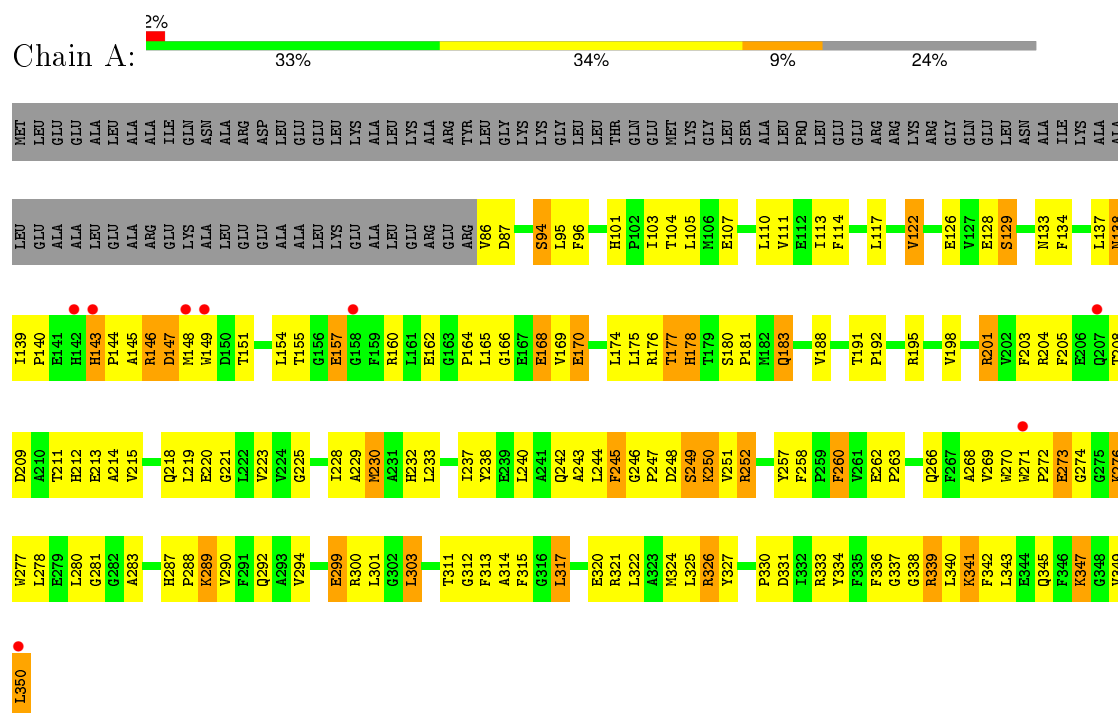
- Molecule 4 is ADENOSINE-5'-[PHENYLALANINOL-PHOSPHATE] (three-letter code: FYA) (formula: C₁₉H₂₅N₆O₇P).



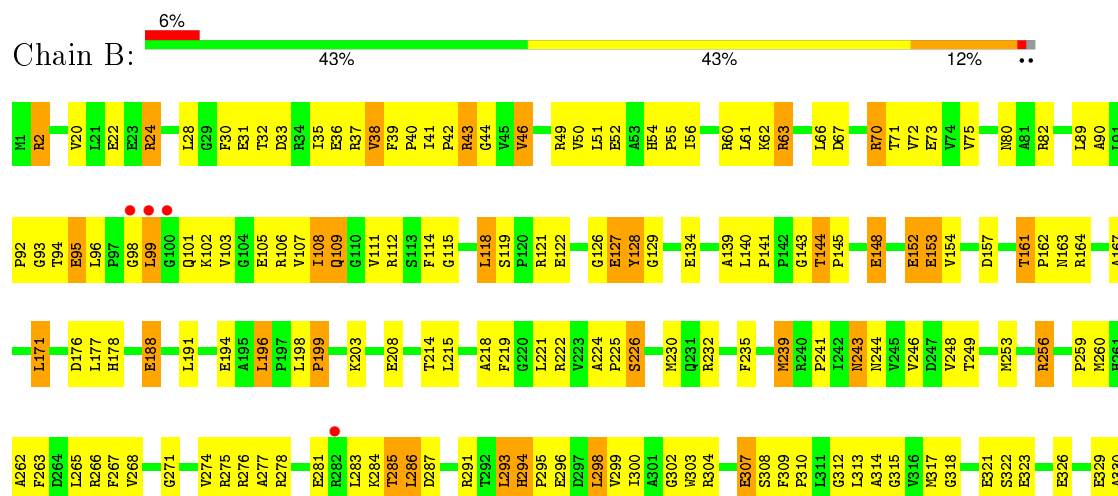
3 Residue-property plots

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: PROTEIN (PHENYLALANYL-TRNA SYNTHETASE)



• Molecule 2: PROTEIN (PHENYLALANYL-TRNA SYNTHETASE)





4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	174.50Å 174.50Å 140.20Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.50 47.41 – 2.50	Depositor EDS
% Data completeness (in resolution range)	95.7 (50.00-2.50) 95.7 (47.41-2.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	15112.19 (at 2.51Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.230 , 0.267 0.229 , 0.264	Depositor DCC
R_{free} test set	4107 reflections (5.03%)	DCC
Wilson B-factor (Å ²)	43.4	Xtriage
Anisotropy	0.241	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 82.6	EDS
Estimated twinning fraction	0.056 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Outliers	0 of 81641 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8439	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.41% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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, FYA

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	0.54	0/2180	0.76	0/2957
2	B	0.53	0/6205	0.79	4/8436 (0.0%)
All	All	0.53	0/8385	0.78	4/11393 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	115	GLY	N-CA-C	-5.83	98.54	113.10
2	B	601	GLU	CB-CA-C	-5.59	99.22	110.40
2	B	128	TYR	N-CA-C	5.35	125.44	111.00
2	B	570	LEU	CA-CB-CG	5.28	127.44	115.30

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	2112	0	2062	164	0
2	B	6054	0	6109	495	0
3	A	1	0	0	0	0
4	A	33	0	24	4	0
5	A	48	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	191	0	0	6	0
All	All	8439	0	8195	626	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 38.

All (626) 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:262:GLU:HG2	5:A:1015:HOH:O	1.45	1.16
2:B:600:LYS:HG2	2:B:601:GLU:H	1.04	1.14
2:B:75:VAL:HG11	2:B:108:ILE:HG21	1.31	1.12
2:B:285:THR:HG21	2:B:291:ARG:HE	1.20	1.02
2:B:294:HIS:CD2	2:B:296:GLU:H	1.79	1.00
1:A:213:GLU:HG3	1:A:214:ALA:H	1.28	0.99
2:B:601:GLU:O	2:B:602:ARG:HD2	1.64	0.98
2:B:600:LYS:HG2	2:B:601:GLU:N	1.78	0.96
2:B:614:GLU:HG2	2:B:624:PHE:HE1	1.29	0.95
2:B:602:ARG:HG2	2:B:602:ARG:HH11	1.28	0.95
2:B:596:LEU:HB2	2:B:599:ALA:HB3	1.46	0.95
1:A:263:PRO:HG3	2:B:461:GLU:HA	1.50	0.94
1:A:165:LEU:HD12	1:A:301:LEU:HD13	1.51	0.93
1:A:101:HIS:HB2	2:B:509:GLU:HG2	1.50	0.93
1:A:143:HIS:ND1	1:A:144:PRO:HD2	1.84	0.93
2:B:593:GLY:HA3	2:B:604:SER:HB3	1.50	0.92
2:B:707:TYR:HE1	2:B:711:GLU:HB2	1.32	0.91
2:B:764:ALA:HA	2:B:767:ARG:HG2	1.52	0.89
2:B:602:ARG:CG	2:B:602:ARG:HH11	1.84	0.89
2:B:294:HIS:HD2	2:B:296:GLU:H	0.89	0.89
1:A:213:GLU:HG2	1:A:215:VAL:H	1.38	0.88
2:B:673:LEU:HD22	2:B:673:LEU:N	1.90	0.86
2:B:710:VAL:O	2:B:714:VAL:HG23	1.75	0.86
2:B:707:TYR:CE1	2:B:711:GLU:HB2	2.11	0.85
1:A:138:ASN:HD21	1:A:289:LYS:HE2	1.39	0.85
1:A:257:TYR:HB3	2:B:161:THR:HG21	1.59	0.85
2:B:278:ARG:NH2	2:B:308:SER:HB3	1.92	0.84
1:A:209:ASP:O	1:A:333:ARG:HD2	1.77	0.84
2:B:729:ASP:N	2:B:744:ALA:HB3	1.92	0.84
2:B:596:LEU:CB	2:B:599:ALA:HB3	2.09	0.83
1:A:101:HIS:HD2	1:A:103:ILE:H	1.26	0.83
2:B:688:SER:HB3	2:B:752:PRO:HA	1.61	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:GLU:CG	1:A:214:ALA:H	1.92	0.82
2:B:497:ARG:O	2:B:501:VAL:HG23	1.78	0.82
1:A:287:HIS:HD2	1:A:289:LYS:H	1.24	0.82
2:B:278:ARG:HH21	2:B:308:SER:HB3	1.42	0.82
2:B:563:ASP:O	2:B:565:PRO:HD3	1.78	0.82
2:B:701:VAL:HG22	2:B:702:PRO:HD2	1.62	0.81
1:A:138:ASN:ND2	1:A:289:LYS:HE2	1.96	0.81
2:B:38:VAL:O	2:B:40:PRO:HD3	1.80	0.81
2:B:70:ARG:HG3	5:B:974:HOH:O	1.79	0.81
2:B:294:HIS:HD2	2:B:296:GLU:N	1.75	0.81
1:A:311:THR:HG22	1:A:312:GLY:N	1.94	0.81
1:A:140:PRO:O	1:A:146:ARG:HB2	1.81	0.81
2:B:549:PHE:O	2:B:553:VAL:HG23	1.81	0.80
1:A:140:PRO:HD2	1:A:145:ALA:HB3	1.63	0.80
2:B:121:ARG:HD3	2:B:127:GLU:O	1.80	0.80
2:B:751:HIS:HD2	2:B:752:PRO:HD2	1.47	0.80
2:B:198:LEU:HD12	2:B:393:LEU:HD13	1.64	0.79
2:B:99:LEU:HD13	2:B:101:GLN:HB2	1.63	0.79
2:B:757:ARG:HG3	2:B:759:GLU:HB2	1.63	0.79
1:A:213:GLU:CG	1:A:214:ALA:N	2.47	0.78
2:B:161:THR:HG22	2:B:162:PRO:HD2	1.64	0.78
2:B:438:GLU:HA	2:B:438:GLU:OE1	1.84	0.78
2:B:643:LEU:HA	2:B:647:GLU:O	1.84	0.77
2:B:519:GLU:HB3	2:B:523:ARG:HH12	1.47	0.77
1:A:250:LYS:H	1:A:270:TRP:HB3	1.48	0.77
1:A:218:GLN:NE2	4:A:1002:FYA:H5'1	1.98	0.77
2:B:427:ILE:HG12	2:B:466:ILE:HG21	1.67	0.76
2:B:614:GLU:HG2	2:B:624:PHE:CE1	2.19	0.76
2:B:656:HIS:HB3	2:B:659:ILE:HD12	1.68	0.75
1:A:311:THR:HG22	1:A:312:GLY:H	1.50	0.75
2:B:490:ALA:HB3	2:B:491:PRO:HD3	1.67	0.75
1:A:160:ARG:HG2	1:A:168:GLU:OE2	1.85	0.75
2:B:730:LEU:HD12	2:B:730:LEU:C	2.07	0.74
2:B:603:LEU:C	2:B:603:LEU:HD12	2.08	0.74
2:B:243:ASN:ND2	2:B:246:VAL:H	1.86	0.74
2:B:697:LEU:HD23	2:B:698:ALA:N	2.03	0.74
2:B:751:HIS:HB3	2:B:754:ARG:O	1.87	0.74
2:B:722:LEU:HD11	2:B:724:SER:O	1.88	0.74
2:B:688:SER:CB	2:B:752:PRO:HA	2.17	0.73
2:B:770:GLU:O	2:B:774:ALA:HB2	1.88	0.73
2:B:194:GLU:OE2	2:B:387:ARG:HG2	1.87	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:404:ILE:HD12	2:B:446:PRO:HD3	1.70	0.73
1:A:263:PRO:HG3	2:B:461:GLU:CA	2.18	0.73
2:B:700:VAL:HA	2:B:741:LYS:O	1.88	0.72
2:B:635:HIS:HB2	2:B:656:HIS:HA	1.70	0.72
2:B:730:LEU:HB2	2:B:742:SER:O	1.89	0.72
2:B:663:LEU:O	2:B:665:LEU:N	2.23	0.72
1:A:262:GLU:OE1	2:B:458:ASP:HA	1.90	0.71
2:B:666:PRO:HB2	2:B:667:PRO:CD	2.19	0.71
2:B:38:VAL:HG22	2:B:153:GLU:O	1.89	0.71
2:B:663:LEU:C	2:B:665:LEU:H	1.94	0.71
2:B:285:THR:HG21	2:B:291:ARG:NE	2.01	0.71
1:A:349:VAL:O	1:A:350:LEU:HD22	1.91	0.71
2:B:178:HIS:O	2:B:430:ARG:NH1	2.24	0.71
2:B:95:GLU:OE2	2:B:102:LYS:HE2	1.91	0.70
2:B:243:ASN:HD21	2:B:246:VAL:H	1.38	0.70
2:B:770:GLU:CG	2:B:771:ALA:N	2.52	0.70
1:A:101:HIS:CD2	1:A:103:ILE:H	2.08	0.70
1:A:103:ILE:HD11	1:A:320:GLU:HG3	1.73	0.70
2:B:736:LEU:HB2	2:B:737:PRO:HD2	1.73	0.70
2:B:761:VAL:HG23	2:B:762:GLU:N	2.06	0.70
2:B:672:GLU:C	2:B:673:LEU:HD22	2.12	0.70
2:B:294:HIS:NE2	2:B:296:GLU:HB2	2.07	0.70
2:B:761:VAL:HG23	2:B:762:GLU:H	1.56	0.70
1:A:271:TRP:HZ3	1:A:276:LYS:HE2	1.57	0.70
2:B:602:ARG:NH1	2:B:602:ARG:CG	2.49	0.70
2:B:589:LEU:HB3	2:B:609:LEU:HD12	1.73	0.70
2:B:757:ARG:HD3	2:B:759:GLU:H	1.56	0.70
2:B:303:TRP:HA	2:B:307:GLU:O	1.91	0.69
2:B:583:THR:HG22	2:B:675:LEU:HD12	1.75	0.69
2:B:728:PHE:HE1	2:B:745:PHE:C	1.96	0.69
2:B:287:ASP:H	2:B:317:MET:HE2	1.57	0.69
1:A:278:LEU:HD13	1:A:325:LEU:HD13	1.74	0.69
1:A:349:VAL:HG12	1:A:350:LEU:HD22	1.75	0.69
1:A:149:TRP:CD1	1:A:177:THR:HG23	2.27	0.69
2:B:604:SER:HA	2:B:608:LEU:HD22	1.74	0.68
1:A:331:ASP:HB3	1:A:334:TYR:CD2	2.28	0.68
2:B:673:LEU:CD2	2:B:673:LEU:N	2.57	0.68
1:A:209:ASP:OD2	1:A:212:HIS:HB2	1.93	0.68
1:A:271:TRP:CZ3	1:A:274:GLY:HA3	2.29	0.68
2:B:734:PRO:HB2	2:B:735:PRO:HD3	1.76	0.68
2:B:698:ALA:HA	2:B:743:LEU:O	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:724:SER:HB2	2:B:748:ARG:HB2	1.75	0.67
2:B:287:ASP:N	2:B:317:MET:HE2	2.09	0.67
1:A:155:THR:CG2	2:B:534:LEU:HD21	2.24	0.67
1:A:229:ALA:H	1:A:232:HIS:HD2	1.41	0.67
2:B:610:LYS:O	2:B:614:GLU:HG3	1.94	0.67
1:A:148:MET:HE2	2:B:162:PRO:HG2	1.75	0.67
2:B:751:HIS:HB2	2:B:756:LEU:CD2	2.25	0.67
1:A:349:VAL:O	1:A:350:LEU:HD13	1.94	0.67
2:B:652:LEU:HD12	2:B:670:LEU:O	1.94	0.67
1:A:278:LEU:HD13	1:A:325:LEU:CD1	2.25	0.67
2:B:520:ASP:OD2	2:B:524:PHE:HE1	1.79	0.66
2:B:239:MET:HE1	2:B:355:GLU:HG3	1.77	0.66
2:B:751:HIS:HB2	2:B:756:LEU:HD21	1.77	0.66
1:A:249:SER:HB2	1:A:270:TRP:O	1.96	0.66
2:B:693:ALA:HB3	2:B:749:PHE:HB2	1.77	0.65
1:A:331:ASP:O	1:A:334:TYR:HB2	1.96	0.65
1:A:246:GLY:HA2	1:A:248:ASP:N	2.12	0.65
2:B:767:ARG:N	2:B:767:ARG:HE	1.95	0.65
1:A:157:GLU:HG3	5:A:1035:HOH:O	1.96	0.65
1:A:205:PHE:CD1	1:A:205:PHE:O	2.50	0.64
1:A:290:VAL:O	1:A:294:VAL:HG23	1.96	0.64
2:B:108:ILE:HG22	2:B:109:GLN:N	2.11	0.64
2:B:656:HIS:NE2	2:B:658:GLU:HB2	2.12	0.64
2:B:557:LYS:HZ3	2:B:663:LEU:HD12	1.63	0.64
1:A:155:THR:HG21	2:B:534:LEU:HD21	1.80	0.64
2:B:645:GLU:OE2	2:B:680:LYS:HB2	1.98	0.64
2:B:194:GLU:HB2	2:B:196:LEU:CD2	2.28	0.64
2:B:588:LEU:HD23	2:B:588:LEU:C	2.18	0.64
2:B:108:ILE:O	2:B:109:GLN:HG2	1.98	0.64
2:B:139:ALA:O	2:B:140:LEU:HD12	1.98	0.64
2:B:757:ARG:HD2	2:B:759:GLU:HG2	1.81	0.63
1:A:248:ASP:HB2	5:A:1013:HOH:O	1.98	0.63
2:B:729:ASP:H	2:B:744:ALA:HB3	1.59	0.63
2:B:194:GLU:O	2:B:390:GLU:HG2	1.98	0.63
2:B:662:GLU:HG3	2:B:662:GLU:O	1.97	0.63
2:B:215:LEU:HB2	2:B:393:LEU:HB2	1.80	0.63
2:B:635:HIS:HB2	2:B:657:PRO:CD	2.28	0.63
2:B:758:ASP:O	2:B:761:VAL:HG22	1.99	0.62
2:B:770:GLU:HG2	2:B:771:ALA:H	1.64	0.62
2:B:671:PHE:HD1	2:B:673:LEU:HD21	1.64	0.62
2:B:767:ARG:CA	2:B:767:ARG:HE	2.12	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:299:VAL:HG23	2:B:312:GLY:O	1.98	0.62
2:B:553:VAL:O	2:B:556:LEU:HB3	1.99	0.62
2:B:219:PHE:HE2	2:B:387:ARG:HE	1.47	0.62
2:B:600:LYS:CG	2:B:601:GLU:H	1.93	0.62
2:B:751:HIS:CD2	2:B:752:PRO:HD2	2.32	0.62
2:B:767:ARG:NE	2:B:767:ARG:HA	2.14	0.62
1:A:213:GLU:HG2	1:A:215:VAL:N	2.12	0.61
2:B:214:THR:HG22	2:B:394:GLU:HG3	1.82	0.61
2:B:707:TYR:CD1	2:B:707:TYR:C	2.73	0.61
2:B:517:ASP:HB3	2:B:520:ASP:OD1	1.99	0.61
2:B:671:PHE:CD1	2:B:673:LEU:HD21	2.36	0.61
2:B:707:TYR:HD1	2:B:707:TYR:C	2.03	0.61
2:B:551:GLY:O	2:B:555:VAL:HG23	2.01	0.61
2:B:381:GLN:HE21	2:B:388:VAL:HG23	1.66	0.61
2:B:702:PRO:CB	2:B:704:PRO:HD2	2.31	0.61
1:A:311:THR:CG2	1:A:312:GLY:N	2.63	0.61
1:A:137:LEU:O	1:A:139:ILE:N	2.30	0.61
2:B:707:TYR:O	2:B:707:TYR:HD1	1.84	0.61
2:B:764:ALA:HA	2:B:767:ARG:CG	2.29	0.61
2:B:153:GLU:HG3	2:B:154:VAL:N	2.14	0.61
2:B:508:GLN:O	2:B:570:LEU:HB2	2.01	0.61
1:A:205:PHE:HE2	2:B:535:ASN:O	1.84	0.60
2:B:707:TYR:HE1	2:B:711:GLU:CB	2.10	0.60
1:A:341:LYS:HE3	2:B:563:ASP:OD2	2.01	0.60
2:B:589:LEU:CB	2:B:609:LEU:HD12	2.31	0.60
1:A:311:THR:CG2	1:A:312:GLY:H	2.15	0.60
2:B:243:ASN:HD22	2:B:243:ASN:C	2.05	0.60
2:B:757:ARG:HB3	2:B:760:GLU:OE2	2.01	0.60
2:B:762:GLU:O	2:B:765:VAL:HG12	2.02	0.60
2:B:635:HIS:ND1	2:B:636:PRO:O	2.35	0.59
2:B:603:LEU:O	2:B:603:LEU:HD12	2.02	0.59
2:B:285:THR:CG2	2:B:291:ARG:HE	2.06	0.59
2:B:657:PRO:HA	2:B:660:ALA:HB3	1.85	0.59
2:B:512:THR:HG22	2:B:545:ARG:HH21	1.68	0.59
2:B:771:ALA:HA	2:B:774:ALA:HB3	1.84	0.59
2:B:281:GLU:HG2	2:B:310:PRO:HG3	1.84	0.59
1:A:134:PHE:O	1:A:137:LEU:O	2.21	0.59
2:B:629:GLN:O	2:B:639:SER:HB3	2.03	0.59
2:B:39:PHE:HB2	2:B:152:GLU:HA	1.84	0.59
1:A:331:ASP:OD1	1:A:333:ARG:HD3	2.01	0.59
2:B:519:GLU:HB3	2:B:523:ARG:NH1	2.17	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:409:GLU:OE2	2:B:413:ARG:HD3	2.03	0.59
1:A:165:LEU:HD11	1:A:303:LEU:HD21	1.84	0.59
2:B:656:HIS:CD2	2:B:658:GLU:H	2.20	0.59
2:B:75:VAL:CG1	2:B:108:ILE:HG21	2.21	0.59
2:B:153:GLU:HB2	2:B:232:ARG:NH2	2.17	0.58
2:B:703:ALA:N	2:B:704:PRO:CD	2.66	0.58
2:B:710:VAL:HG11	2:B:743:LEU:CD1	2.33	0.58
1:A:341:LYS:H	1:A:341:LYS:HE2	1.68	0.58
1:A:299:GLU:HG3	1:A:300:ARG:N	2.18	0.58
2:B:649:VAL:HG23	2:B:674:ARG:H	1.69	0.58
2:B:119:SER:OG	2:B:122:GLU:HG3	2.04	0.58
2:B:259:PRO:HG2	2:B:360:PHE:CE2	2.38	0.58
2:B:161:THR:HB	2:B:163:ASN:OD1	2.03	0.58
2:B:671:PHE:HD1	2:B:673:LEU:CD2	2.15	0.58
2:B:294:HIS:CD2	2:B:296:GLU:HB2	2.39	0.58
2:B:706:PRO:O	2:B:710:VAL:HG23	2.04	0.58
2:B:265:LEU:HD23	2:B:268:VAL:HG21	1.86	0.58
2:B:108:ILE:CG2	2:B:109:GLN:N	2.67	0.57
2:B:728:PHE:HD1	2:B:728:PHE:H	1.49	0.57
2:B:99:LEU:HD22	2:B:101:GLN:NE2	2.19	0.57
1:A:148:MET:CE	2:B:162:PRO:HG2	2.34	0.57
2:B:757:ARG:CG	2:B:759:GLU:HB2	2.34	0.57
2:B:702:PRO:C	2:B:704:PRO:HD2	2.24	0.57
2:B:767:ARG:NE	2:B:767:ARG:CA	2.67	0.57
2:B:377:LEU:HD22	2:B:388:VAL:HG13	1.87	0.57
1:A:180:SER:O	1:A:183:GLN:HB2	2.05	0.57
2:B:635:HIS:HB2	2:B:657:PRO:HD3	1.85	0.57
2:B:293:LEU:HD13	2:B:293:LEU:N	2.18	0.57
2:B:657:PRO:HA	2:B:660:ALA:CB	2.35	0.57
2:B:286:LEU:HD11	2:B:323:GLU:CD	2.25	0.57
2:B:557:LYS:NZ	2:B:663:LEU:HD12	2.19	0.57
1:A:162:GLU:CG	1:A:166:GLY:HA2	2.35	0.57
1:A:126:GLU:HG3	2:B:575:ARG:HD2	1.87	0.56
2:B:707:TYR:HE2	2:B:727:LEU:HD22	1.70	0.56
2:B:38:VAL:O	2:B:40:PRO:CD	2.53	0.56
1:A:143:HIS:CG	1:A:144:PRO:HD2	2.40	0.56
2:B:702:PRO:HA	2:B:739:GLY:O	2.06	0.56
1:A:252:ARG:HE	1:A:268:ALA:HB3	1.71	0.56
2:B:758:ASP:HA	2:B:761:VAL:HG22	1.88	0.56
1:A:122:VAL:CG1	2:B:488:VAL:HG13	2.35	0.56
1:A:271:TRP:NE1	1:A:273:GLU:HB2	2.21	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:LYS:HG3	1:A:278:LEU:HD21	1.87	0.56
2:B:634:LEU:HB3	2:B:639:SER:OG	2.05	0.56
2:B:226:SER:OG	2:B:244:ASN:HA	2.05	0.56
2:B:71:THR:HG22	2:B:72:VAL:N	2.21	0.56
1:A:238:TYR:HA	1:A:251:VAL:HG11	1.88	0.55
2:B:366:PRO:C	2:B:367:LEU:HD22	2.27	0.55
1:A:287:HIS:CD2	1:A:288:PRO:HD2	2.41	0.55
2:B:514:SER:O	2:B:545:ARG:HG2	2.07	0.55
2:B:381:GLN:O	2:B:385:GLY:HA2	2.06	0.55
2:B:770:GLU:CG	2:B:771:ALA:H	2.19	0.55
2:B:55:PRO:HA	2:B:62:LYS:HA	1.89	0.55
2:B:221:LEU:CD2	2:B:386:ALA:HB2	2.36	0.55
2:B:527:ASP:HB3	2:B:528:PRO:HD2	1.89	0.55
2:B:38:VAL:C	2:B:40:PRO:HD3	2.27	0.55
1:A:96:PHE:CE1	2:B:567:ARG:HG3	2.42	0.54
2:B:458:ASP:O	2:B:462:GLU:HG2	2.08	0.54
2:B:731:TYR:HD1	2:B:742:SER:HG	1.55	0.54
1:A:94:SER:HB2	2:B:567:ARG:NH2	2.21	0.54
1:A:349:VAL:CG1	1:A:350:LEU:HD22	2.38	0.54
1:A:219:LEU:HB3	1:A:317:LEU:CD2	2.37	0.54
2:B:734:PRO:CB	2:B:735:PRO:HD3	2.37	0.54
2:B:370:VAL:N	2:B:371:PRO:HD2	2.23	0.54
2:B:530:ARG:HB2	2:B:530:ARG:HH11	1.72	0.54
2:B:509:GLU:HG3	2:B:510:VAL:N	2.22	0.54
1:A:343:LEU:HD13	2:B:509:GLU:O	2.07	0.54
2:B:757:ARG:CD	2:B:759:GLU:H	2.21	0.54
2:B:761:VAL:CG2	2:B:762:GLU:H	2.20	0.54
2:B:239:MET:CE	2:B:355:GLU:HG3	2.37	0.54
2:B:163:ASN:O	2:B:452:ASP:HB3	2.08	0.54
2:B:224:ALA:HB1	2:B:225:PRO:HD2	1.90	0.54
2:B:194:GLU:HB2	2:B:196:LEU:HD21	1.90	0.54
2:B:556:LEU:C	2:B:556:LEU:HD12	2.28	0.53
2:B:663:LEU:C	2:B:665:LEU:N	2.62	0.53
2:B:194:GLU:HB2	2:B:196:LEU:HD23	1.91	0.53
2:B:516:MET:HE3	2:B:545:ARG:HA	1.91	0.53
1:A:257:TYR:HB3	2:B:161:THR:CG2	2.36	0.53
2:B:702:PRO:HB2	2:B:704:PRO:HD2	1.90	0.53
2:B:707:TYR:O	2:B:707:TYR:CD1	2.62	0.53
2:B:773:ARG:C	2:B:775:ARG:N	2.62	0.53
2:B:650:GLY:HA3	2:B:673:LEU:HD13	1.90	0.53
2:B:669:HIS:C	2:B:670:LEU:HD13	2.29	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:GLN:OE1	1:A:247:PRO:HB3	2.09	0.53
2:B:275:ARG:O	2:B:299:VAL:HG12	2.08	0.53
2:B:575:ARG:HG2	2:B:581:GLU:HG2	1.90	0.53
2:B:523:ARG:NH1	2:B:523:ARG:HG3	2.23	0.52
2:B:498:LEU:O	2:B:498:LEU:HD12	2.08	0.52
2:B:617:PHE:HB3	2:B:622:LEU:O	2.09	0.52
2:B:729:ASP:CG	2:B:730:LEU:N	2.62	0.52
1:A:281:GLY:HA2	4:A:1002:FYA:O3'	2.09	0.52
1:A:148:MET:HE3	1:A:257:TYR:HE2	1.75	0.52
2:B:728:PHE:CD1	2:B:728:PHE:N	2.78	0.52
1:A:96:PHE:CZ	2:B:567:ARG:HG3	2.43	0.52
2:B:263:PHE:O	2:B:331:ILE:HB	2.10	0.52
2:B:459:LEU:O	2:B:462:GLU:HB2	2.10	0.52
2:B:285:THR:HG23	2:B:287:ASP:OD1	2.09	0.52
2:B:761:VAL:CG2	2:B:762:GLU:N	2.73	0.52
2:B:20:VAL:O	2:B:24:ARG:HG2	2.10	0.52
2:B:267:PHE:CE2	2:B:321:GLU:HG2	2.45	0.52
2:B:51:LEU:HD21	2:B:67:ASP:HB2	1.91	0.52
2:B:601:GLU:C	2:B:602:ARG:HD2	2.29	0.52
2:B:671:PHE:CD1	2:B:673:LEU:CD2	2.92	0.52
2:B:733:GLY:HA3	2:B:736:LEU:HD21	1.92	0.52
2:B:549:PHE:CG	2:B:550:PRO:HD3	2.45	0.52
2:B:699:VAL:HG12	2:B:773:ARG:NH2	2.25	0.52
2:B:447:PRO:HB2	2:B:449:HIS:CE1	2.45	0.52
1:A:165:LEU:HD12	1:A:301:LEU:CD1	2.33	0.51
2:B:621:GLY:O	2:B:680:LYS:HE2	2.10	0.51
2:B:666:PRO:HB2	2:B:667:PRO:HD2	1.92	0.51
2:B:63:ARG:HD2	2:B:73:GLU:OE2	2.10	0.51
2:B:697:LEU:C	2:B:697:LEU:HD23	2.29	0.51
2:B:359:ARG:HH11	2:B:359:ARG:HG3	1.75	0.51
2:B:191:LEU:HD13	2:B:378:SER:HA	1.91	0.51
2:B:563:ASP:C	2:B:565:PRO:HD3	2.30	0.51
1:A:246:GLY:HA2	1:A:248:ASP:H	1.74	0.51
2:B:635:HIS:HB2	2:B:657:PRO:HD2	1.91	0.51
2:B:224:ALA:HB1	2:B:225:PRO:CD	2.41	0.51
2:B:107:VAL:HA	2:B:112:ARG:HA	1.92	0.51
2:B:644:VAL:HG22	2:B:645:GLU:H	1.74	0.51
1:A:175:LEU:HB3	1:A:203:PHE:CD2	2.45	0.51
2:B:538:ALA:HB1	2:B:540:GLU:OE1	2.10	0.51
2:B:548:LEU:HD13	2:B:576:VAL:CG1	2.41	0.51
2:B:770:GLU:HG2	2:B:771:ALA:N	2.24	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:770:GLU:O	2:B:774:ALA:CB	2.58	0.51
2:B:337:CYS:HB2	2:B:369:GLN:NE2	2.25	0.51
2:B:285:THR:HG22	2:B:291:ARG:HG3	1.93	0.51
2:B:770:GLU:HG3	2:B:771:ALA:N	2.24	0.51
1:A:283:ALA:HB2	1:A:315:PHE:CB	2.40	0.51
2:B:723:GLU:OE1	2:B:750:ARG:HD2	2.10	0.51
2:B:643:LEU:C	2:B:644:VAL:O	2.47	0.51
1:A:94:SER:O	2:B:594:VAL:HB	2.11	0.51
2:B:762:GLU:C	2:B:765:VAL:HG12	2.31	0.51
1:A:271:TRP:CH2	1:A:274:GLY:HA3	2.45	0.51
2:B:409:GLU:OE1	2:B:413:ARG:NH1	2.43	0.51
2:B:635:HIS:CE1	2:B:636:PRO:O	2.64	0.50
2:B:530:ARG:NH1	2:B:530:ARG:HB2	2.25	0.50
1:A:101:HIS:CD2	1:A:103:ILE:HB	2.46	0.50
1:A:270:TRP:O	1:A:272:PRO:HD3	2.10	0.50
2:B:773:ARG:HB2	2:B:773:ARG:NH1	2.27	0.50
5:A:1006:HOH:O	2:B:31:GLU:HG3	2.11	0.50
1:A:110:LEU:HD11	1:A:322:LEU:HD23	1.93	0.50
2:B:286:LEU:HB3	2:B:317:MET:CE	2.42	0.50
2:B:105:GLU:HG3	2:B:114:PHE:CD1	2.47	0.50
2:B:36:GLU:O	2:B:154:VAL:HA	2.11	0.50
2:B:656:HIS:HB3	2:B:659:ILE:CD1	2.39	0.50
2:B:299:VAL:HG22	2:B:300:ILE:N	2.26	0.50
2:B:30:PHE:HA	5:B:976:HOH:O	2.10	0.50
2:B:262:ALA:HB1	2:B:331:ILE:HG13	1.94	0.50
2:B:274:VAL:HG12	2:B:298:LEU:HD21	1.92	0.50
2:B:2:ARG:HD2	5:B:828:HOH:O	2.12	0.50
2:B:557:LYS:HE3	2:B:664:GLU:OE2	2.11	0.50
1:A:105:LEU:HD22	1:A:349:VAL:CG1	2.41	0.50
2:B:253:MET:HG3	2:B:259:PRO:HA	1.93	0.50
2:B:649:VAL:CG2	2:B:674:ARG:H	2.24	0.50
2:B:72:VAL:HG21	2:B:114:PHE:CD2	2.47	0.50
2:B:188:GLU:HG3	2:B:188:GLU:O	2.12	0.49
1:A:101:HIS:CB	2:B:509:GLU:HG2	2.32	0.49
2:B:92:PRO:HG3	2:B:114:PHE:O	2.13	0.49
2:B:46:VAL:HB	2:B:143:GLY:O	2.12	0.49
1:A:198:VAL:N	1:A:220:GLU:O	2.38	0.49
2:B:598:TRP:CZ3	2:B:599:ALA:HB2	2.47	0.49
2:B:710:VAL:HG11	2:B:743:LEU:HD12	1.94	0.49
2:B:688:SER:OG	2:B:752:PRO:HA	2.13	0.49
2:B:520:ASP:HA	2:B:523:ARG:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:VAL:HG13	2:B:488:VAL:HG13	1.94	0.49
2:B:692:ALA:HA	2:B:749:PHE:O	2.12	0.49
2:B:50:VAL:HA	2:B:66:LEU:HD23	1.93	0.49
2:B:517:ASP:OD1	2:B:519:GLU:HB2	2.13	0.49
1:A:134:PHE:CZ	1:A:151:THR:HG21	2.48	0.49
2:B:765:VAL:HG13	2:B:766:SER:N	2.28	0.49
1:A:114:PHE:HA	1:A:117:LEU:HD13	1.95	0.49
2:B:286:LEU:HB2	2:B:318:GLY:O	2.13	0.48
2:B:609:LEU:HD21	2:B:671:PHE:HD2	1.77	0.48
2:B:613:LEU:O	2:B:617:PHE:HD1	1.96	0.48
2:B:538:ALA:HB3	2:B:541:LYS:HG2	1.95	0.48
1:A:268:ALA:HA	1:A:278:LEU:O	2.12	0.48
2:B:355:GLU:O	2:B:359:ARG:HG3	2.13	0.48
2:B:532:LEU:N	2:B:532:LEU:HD22	2.28	0.48
2:B:509:GLU:HA	2:B:571:PHE:CE1	2.47	0.48
2:B:587:GLY:HA3	2:B:671:PHE:CE2	2.47	0.48
2:B:575:ARG:HG2	2:B:581:GLU:CG	2.44	0.48
1:A:283:ALA:HB2	1:A:315:PHE:HB3	1.95	0.48
1:A:213:GLU:HG2	1:A:214:ALA:N	2.28	0.48
2:B:119:SER:OG	2:B:129:GLY:HA2	2.14	0.48
1:A:86:VAL:HG23	1:A:87:ASP:N	2.29	0.48
1:A:104:THR:HG1	2:B:511:TYR:HH	1.62	0.48
2:B:721:TYR:N	2:B:721:TYR:CD1	2.81	0.48
2:B:703:ALA:N	2:B:704:PRO:HD2	2.28	0.47
2:B:331:ILE:HG12	2:B:332:ALA:N	2.29	0.47
2:B:596:LEU:HD13	2:B:598:TRP:CH2	2.50	0.47
1:A:271:TRP:CZ3	1:A:276:LYS:HE2	2.43	0.47
2:B:642:VAL:HG23	2:B:651:PHE:HA	1.96	0.47
2:B:303:TRP:CA	2:B:307:GLU:O	2.61	0.47
1:A:205:PHE:HD1	1:A:205:PHE:O	1.97	0.47
1:A:342:PHE:O	1:A:345:GLN:HB2	2.15	0.47
1:A:337:GLY:O	1:A:339:ARG:N	2.41	0.47
1:A:339:ARG:NH1	2:B:562:LEU:HD22	2.30	0.47
1:A:349:VAL:CG1	1:A:349:VAL:O	2.63	0.47
2:B:650:GLY:HA3	2:B:672:GLU:O	2.15	0.47
2:B:244:ASN:O	2:B:248:VAL:HG23	2.15	0.47
2:B:486:ARG:HA	2:B:486:ARG:HD3	1.73	0.47
2:B:642:VAL:HG23	2:B:651:PHE:CA	2.45	0.47
1:A:208:THR:HG22	1:A:213:GLU:HA	1.96	0.47
2:B:609:LEU:HD22	2:B:652:LEU:CD1	2.45	0.47
2:B:253:MET:HE2	2:B:253:MET:HB3	1.67	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:293:LEU:CD1	2:B:293:LEU:N	2.78	0.47
2:B:671:PHE:C	2:B:671:PHE:CD1	2.88	0.47
2:B:702:PRO:O	2:B:741:LYS:HE2	2.15	0.47
2:B:751:HIS:HD2	2:B:752:PRO:CD	2.22	0.47
2:B:661:GLN:C	2:B:663:LEU:N	2.68	0.47
2:B:105:GLU:HG3	2:B:114:PHE:HD1	1.79	0.47
2:B:615:ALA:O	2:B:618:ALA:HB3	2.15	0.47
2:B:651:PHE:CE2	2:B:672:GLU:HB3	2.50	0.46
2:B:765:VAL:CG1	2:B:766:SER:N	2.78	0.46
2:B:523:ARG:HH11	2:B:523:ARG:HG3	1.79	0.46
2:B:533:LEU:HB2	2:B:536:PRO:HG3	1.97	0.46
1:A:209:ASP:C	1:A:333:ARG:HD2	2.35	0.46
2:B:243:ASN:ND2	2:B:243:ASN:C	2.69	0.46
2:B:536:PRO:HB3	2:B:542:ALA:HA	1.97	0.46
2:B:731:TYR:CE2	2:B:733:GLY:HA2	2.50	0.46
1:A:260:PHE:CD1	1:A:260:PHE:N	2.82	0.46
2:B:556:LEU:HD22	2:B:588:LEU:HD21	1.96	0.46
2:B:675:LEU:HB3	2:B:676:PRO:HA	1.96	0.46
1:A:221:GLY:O	1:A:314:ALA:HA	2.14	0.46
2:B:537:LEU:HA	2:B:537:LEU:HD23	1.60	0.46
2:B:374:ARG:HA	2:B:374:ARG:HD3	1.77	0.46
1:A:94:SER:HB2	2:B:567:ARG:CZ	2.46	0.46
1:A:147:ASP:N	1:A:147:ASP:OD1	2.49	0.46
1:A:215:VAL:HG21	2:B:513:TYR:CD2	2.51	0.46
2:B:266:ARG:NH1	2:B:267:PHE:CE1	2.83	0.46
1:A:213:GLU:HG3	1:A:214:ALA:N	2.01	0.46
2:B:93:GLY:O	2:B:102:LYS:HD3	2.16	0.46
1:A:134:PHE:HZ	1:A:151:THR:HG21	1.81	0.46
2:B:417:THR:HG22	2:B:473:PRO:HD2	1.97	0.46
1:A:169:VAL:HG22	1:A:170:GLU:N	2.31	0.46
2:B:294:HIS:CD2	2:B:296:GLU:N	2.63	0.46
1:A:164:PRO:HG2	1:A:188:VAL:HG11	1.97	0.46
2:B:695:ARG:HH22	2:B:761:VAL:HG21	1.80	0.46
2:B:516:MET:HE1	2:B:546:THR:H	1.81	0.46
2:B:631:PHE:HB2	2:B:634:LEU:HB2	1.97	0.46
2:B:730:LEU:HD12	2:B:730:LEU:O	2.16	0.45
1:A:219:LEU:HB3	1:A:317:LEU:HD22	1.97	0.45
2:B:549:PHE:CD2	2:B:550:PRO:HD3	2.51	0.45
2:B:660:ALA:O	2:B:663:LEU:O	2.34	0.45
2:B:219:PHE:CZ	2:B:387:ARG:HB3	2.51	0.45
1:A:238:TYR:HA	1:A:251:VAL:CG1	2.45	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:THR:HG22	1:A:192:PRO:HD2	1.98	0.45
2:B:695:ARG:HB2	2:B:747:LEU:HB2	1.99	0.45
1:A:270:TRP:CE3	1:A:270:TRP:HA	2.51	0.45
2:B:243:ASN:HD21	2:B:246:VAL:HG23	1.82	0.45
2:B:164:ARG:O	2:B:167:ALA:HB3	2.16	0.45
2:B:509:GLU:HB2	2:B:571:PHE:CE1	2.52	0.45
2:B:600:LYS:N	2:B:600:LYS:HD2	2.31	0.45
2:B:617:PHE:HD2	2:B:622:LEU:HB2	1.81	0.45
2:B:757:ARG:HD3	2:B:758:ASP:N	2.31	0.45
1:A:287:HIS:CD2	1:A:289:LYS:H	2.16	0.45
1:A:315:PHE:C	1:A:315:PHE:CD1	2.89	0.45
1:A:271:TRP:HB2	1:A:278:LEU:HD11	1.98	0.45
2:B:219:PHE:CE2	2:B:387:ARG:NE	2.82	0.45
2:B:300:ILE:HD13	2:B:314:ALA:HA	1.97	0.45
1:A:258:PHE:CZ	4:A:1002:FYA:HD1	2.52	0.45
2:B:427:ILE:HG22	2:B:428:LEU:N	2.32	0.45
2:B:44:GLY:HA3	2:B:94:THR:OG1	2.17	0.45
1:A:336:PHE:HB3	2:B:513:TYR:CE1	2.51	0.45
2:B:556:LEU:O	2:B:556:LEU:HD12	2.17	0.45
2:B:219:PHE:HE2	2:B:387:ARG:NE	2.12	0.45
2:B:399:LYS:HA	2:B:400:PRO:HD3	1.82	0.45
2:B:623:ALA:O	2:B:645:GLU:HA	2.17	0.44
1:A:126:GLU:HB3	1:A:203:PHE:HE2	1.83	0.44
1:A:169:VAL:CG2	1:A:170:GLU:N	2.80	0.44
2:B:28:LEU:HD13	2:B:176:ASP:HB3	1.98	0.44
2:B:218:ALA:O	2:B:330:ALA:HA	2.17	0.44
2:B:701:VAL:CG2	2:B:702:PRO:HD2	2.42	0.44
2:B:121:ARG:HA	2:B:126:GLY:O	2.17	0.44
1:A:218:GLN:NE2	4:A:1002:FYA:C5'	2.75	0.44
2:B:2:ARG:HH11	2:B:2:ARG:CG	2.30	0.44
2:B:35:ILE:O	2:B:36:GLU:HG2	2.16	0.44
2:B:657:PRO:O	2:B:660:ALA:HB3	2.18	0.44
1:A:105:LEU:HD22	1:A:349:VAL:HG12	1.98	0.44
2:B:759:GLU:O	2:B:763:GLU:HB2	2.18	0.44
1:A:180:SER:N	1:A:181:PRO:HD2	2.32	0.44
2:B:271:GLY:O	2:B:302:GLY:HA2	2.17	0.44
2:B:700:VAL:CG1	2:B:736:LEU:HD13	2.47	0.44
2:B:721:TYR:CD2	2:B:756:LEU:HD21	2.53	0.44
2:B:549:PHE:CD1	2:B:550:PRO:N	2.86	0.44
2:B:603:LEU:C	2:B:603:LEU:CD1	2.82	0.44
2:B:359:ARG:HH11	2:B:359:ARG:CG	2.30	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:PHE:HE2	1:A:269:VAL:HG11	1.82	0.44
2:B:548:LEU:HD22	2:B:584:HIS:CB	2.48	0.44
2:B:609:LEU:CD2	2:B:652:LEU:HD13	2.47	0.44
2:B:730:LEU:HD12	2:B:731:TYR:N	2.33	0.44
2:B:695:ARG:HH12	2:B:761:VAL:HG23	1.82	0.44
1:A:242:GLN:O	1:A:245:PHE:O	2.35	0.44
1:A:326:ARG:HG2	1:A:327:TYR:CD1	2.52	0.44
2:B:235:PHE:CZ	2:B:241:PRO:HD2	2.53	0.44
2:B:707:TYR:O	2:B:710:VAL:HB	2.18	0.44
2:B:643:LEU:HD13	2:B:648:GLU:HA	2.00	0.44
1:A:160:ARG:N	2:B:579:GLU:O	2.45	0.44
1:A:225:GLY:O	1:A:228:ILE:HG12	2.17	0.44
1:A:143:HIS:CB	1:A:144:PRO:HD2	2.48	0.44
2:B:728:PHE:O	2:B:729:ASP:HB2	2.18	0.44
1:A:339:ARG:NH1	2:B:563:ASP:OD1	2.51	0.44
2:B:198:LEU:HA	2:B:199:PRO:HD2	1.54	0.44
2:B:523:ARG:HH11	2:B:523:ARG:CG	2.30	0.43
2:B:557:LYS:HA	2:B:665:LEU:HD11	2.00	0.43
1:A:277:TRP:O	1:A:278:LEU:HD23	2.17	0.43
2:B:42:PRO:HA	2:B:43:ARG:NH1	2.33	0.43
2:B:732:GLN:HA	2:B:732:GLN:OE1	2.14	0.43
2:B:276:ARG:NH1	2:B:296:GLU:O	2.38	0.43
2:B:598:TRP:CE3	2:B:599:ALA:HB2	2.53	0.43
2:B:368:GLY:O	2:B:371:PRO:HG2	2.17	0.43
2:B:303:TRP:HB2	2:B:307:GLU:O	2.18	0.43
2:B:253:MET:O	2:B:256:ARG:O	2.35	0.43
2:B:705:THR:HG23	2:B:706:PRO:HD2	2.00	0.43
1:A:277:TRP:C	1:A:278:LEU:HD23	2.39	0.43
2:B:92:PRO:HD3	2:B:114:PHE:O	2.19	0.43
2:B:575:ARG:HD3	2:B:581:GLU:OE2	2.18	0.43
2:B:90:ALA:HB2	2:B:118:LEU:HD11	2.00	0.43
2:B:535:ASN:N	2:B:535:ASN:OD1	2.52	0.43
2:B:72:VAL:HG22	2:B:73:GLU:N	2.34	0.43
2:B:379:LEU:HA	2:B:379:LEU:HD23	1.92	0.43
1:A:237:ILE:O	1:A:240:LEU:HB3	2.18	0.43
2:B:600:LYS:HG2	2:B:601:GLU:HG3	2.00	0.43
1:A:149:TRP:CD1	1:A:177:THR:CG2	2.98	0.43
1:A:229:ALA:H	1:A:232:HIS:CD2	2.27	0.43
2:B:265:LEU:HD23	2:B:265:LEU:HA	1.81	0.43
1:A:230:MET:HB2	2:B:415:LEU:HA	2.01	0.43
2:B:549:PHE:CD1	2:B:549:PHE:C	2.92	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:768:VAL:O	2:B:772:LEU:HB2	2.19	0.43
2:B:520:ASP:OD2	2:B:554:ARG:NH1	2.49	0.43
2:B:309:PHE:N	2:B:309:PHE:CD1	2.87	0.43
1:A:201:ARG:HD2	1:A:215:VAL:CG1	2.48	0.43
2:B:589:LEU:HB3	2:B:609:LEU:CD1	2.43	0.43
1:A:341:LYS:HD3	1:A:341:LYS:HA	1.87	0.43
1:A:233:LEU:HB2	1:A:313:PHE:CG	2.54	0.43
2:B:650:GLY:CA	2:B:672:GLU:O	2.67	0.42
2:B:579:GLU:HG2	2:B:579:GLU:H	1.38	0.42
2:B:377:LEU:HD22	2:B:388:VAL:CG1	2.47	0.42
2:B:600:LYS:HB3	2:B:600:LYS:HE2	1.78	0.42
2:B:286:LEU:HB3	2:B:317:MET:HE2	2.00	0.42
2:B:520:ASP:OD2	2:B:524:PHE:CE1	2.67	0.42
1:A:122:VAL:HG13	2:B:488:VAL:CG1	2.49	0.42
2:B:498:LEU:HD12	2:B:498:LEU:C	2.39	0.42
2:B:80:ASN:HD22	2:B:134:GLU:HG3	1.84	0.42
2:B:602:ARG:NH1	2:B:602:ARG:HG3	2.31	0.42
2:B:276:ARG:HB3	2:B:295:PRO:O	2.19	0.42
2:B:703:ALA:HA	2:B:741:LYS:HG2	2.00	0.42
1:A:339:ARG:HH12	2:B:562:LEU:HD22	1.85	0.42
2:B:337:CYS:HB2	2:B:369:GLN:HE22	1.84	0.42
2:B:176:ASP:OD2	2:B:465:ARG:NH2	2.53	0.42
1:A:107:GLU:O	1:A:111:VAL:HG23	2.20	0.42
2:B:287:ASP:HB3	2:B:317:MET:HE1	2.01	0.42
2:B:717:ALA:C	2:B:768:VAL:HG22	2.40	0.42
1:A:177:THR:HG22	1:A:178:HIS:CD2	2.54	0.42
2:B:80:ASN:O	2:B:82:ARG:HD3	2.19	0.42
2:B:22:GLU:HG3	2:B:32:THR:HB	2.00	0.42
2:B:589:LEU:HD21	2:B:608:LEU:HD23	2.01	0.42
2:B:651:PHE:O	2:B:671:PHE:HA	2.19	0.42
2:B:307:GLU:OE1	2:B:307:GLU:HA	2.14	0.42
2:B:171:LEU:HD23	2:B:171:LEU:HA	1.74	0.42
2:B:700:VAL:HG12	2:B:736:LEU:CD1	2.49	0.42
2:B:430:ARG:NH2	5:B:786:HOH:O	2.53	0.42
2:B:249:THR:HB	2:B:260:MET:HG3	2.02	0.42
1:A:321:ARG:HA	1:A:324:MET:HE2	2.01	0.42
2:B:98:GLY:HA3	5:B:975:HOH:O	2.19	0.42
1:A:113:ILE:O	1:A:117:LEU:HD12	2.19	0.42
2:B:692:ALA:CA	2:B:749:PHE:O	2.68	0.42
2:B:538:ALA:HB3	2:B:541:LYS:CG	2.50	0.42
2:B:643:LEU:N	2:B:643:LEU:HD22	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:524:PHE:O	2:B:526:LEU:HG	2.20	0.42
2:B:509:GLU:CA	2:B:571:PHE:CE1	3.03	0.41
2:B:727:LEU:HA	2:B:744:ALA:O	2.20	0.41
2:B:519:GLU:O	2:B:522:ARG:HB2	2.20	0.41
2:B:535:ASN:N	2:B:536:PRO:HD3	2.35	0.41
1:A:128:GLU:HG3	1:A:129:SER:H	1.84	0.41
2:B:277:ALA:O	2:B:295:PRO:HA	2.19	0.41
2:B:467:GLN:HA	2:B:467:GLN:OE1	2.20	0.41
2:B:294:HIS:CD2	2:B:296:GLU:CB	3.03	0.41
1:A:271:TRP:CE3	1:A:274:GLY:HA3	2.55	0.41
2:B:265:LEU:HD23	2:B:268:VAL:CG2	2.50	0.41
2:B:636:PRO:C	2:B:638:VAL:H	2.24	0.41
2:B:572:GLU:HG3	2:B:573:VAL:N	2.36	0.41
2:B:177:LEU:HA	2:B:177:LEU:HD23	1.83	0.41
2:B:512:THR:HB	2:B:572:GLU:OE2	2.21	0.41
2:B:51:LEU:HD11	2:B:67:ASP:HB2	2.03	0.41
2:B:505:LEU:HD12	2:B:505:LEU:HA	1.84	0.41
2:B:600:LYS:H	2:B:600:LYS:HD2	1.85	0.41
2:B:304:ARG:O	2:B:307:GLU:HB2	2.20	0.41
2:B:529:PRO:CB	2:B:543:ALA:HB1	2.50	0.41
2:B:52:GLU:HB3	2:B:54:HIS:CE1	2.55	0.41
2:B:402:GLU:HG3	5:B:903:HOH:O	2.21	0.41
2:B:344:ARG:HG3	2:B:361:GLU:OE2	2.21	0.41
1:A:144:PRO:HB2	2:B:162:PRO:HB3	2.03	0.41
2:B:730:LEU:CD1	2:B:730:LEU:C	2.77	0.41
2:B:766:SER:C	2:B:768:VAL:N	2.72	0.41
2:B:633:PHE:HD1	2:B:634:LEU:HD13	1.85	0.41
2:B:609:LEU:HD23	2:B:613:LEU:CD2	2.51	0.41
2:B:669:HIS:C	2:B:670:LEU:CD1	2.89	0.41
2:B:520:ASP:HB3	2:B:524:PHE:CD1	2.55	0.41
1:A:133:ASN:HD21	1:A:177:THR:HB	1.85	0.41
1:A:133:ASN:OD1	1:A:176:ARG:HA	2.21	0.41
2:B:71:THR:CG2	2:B:72:VAL:N	2.83	0.41
1:A:347:LYS:HZ2	1:A:347:LYS:HG2	1.71	0.41
2:B:145:PRO:O	2:B:148:GLU:HB2	2.21	0.41
2:B:728:PHE:HE1	2:B:746:HIS:N	2.19	0.41
2:B:756:LEU:HA	2:B:756:LEU:HD23	1.82	0.41
1:A:126:GLU:CG	2:B:575:ARG:HD2	2.50	0.41
2:B:699:VAL:CG1	2:B:773:ARG:NH2	2.84	0.41
2:B:267:PHE:CD2	2:B:321:GLU:HG2	2.56	0.41
2:B:90:ALA:O	2:B:103:VAL:HG11	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:764:ALA:O	2:B:767:ARG:HB2	2.21	0.40
2:B:278:ARG:NH2	2:B:308:SER:O	2.55	0.40
2:B:549:PHE:N	2:B:550:PRO:CD	2.84	0.40
2:B:588:LEU:C	2:B:588:LEU:CD2	2.88	0.40
2:B:656:HIS:HD2	2:B:657:PRO:CD	2.34	0.40
2:B:566:GLU:CG	2:B:592:GLU:HG2	2.51	0.40
2:B:108:ILE:HG22	2:B:109:GLN:H	1.81	0.40
2:B:35:ILE:C	2:B:36:GLU:HG2	2.42	0.40
2:B:520:ASP:C	2:B:522:ARG:N	2.72	0.40
1:A:278:LEU:HD13	1:A:325:LEU:HD11	2.01	0.40
1:A:244:LEU:O	1:A:326:ARG:NH2	2.53	0.40
2:B:41:ILE:HA	2:B:42:PRO:HD2	1.92	0.40
2:B:507:PHE:N	2:B:507:PHE:CD1	2.89	0.40
2:B:773:ARG:HB2	2:B:773:ARG:HH11	1.86	0.40
2:B:531:LEU:C	2:B:532:LEU:HD22	2.42	0.40
2:B:141:PRO:HG2	2:B:144:THR:CG2	2.51	0.40
1:A:260:PHE:C	1:A:287:HIS:HB2	2.42	0.40
2:B:381:GLN:NE2	2:B:388:VAL:HG23	2.33	0.40
1:A:113:ILE:HD13	1:A:243:ALA:CB	2.52	0.40
1:A:195:ARG:HG2	1:A:223:VAL:HG22	2.03	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	263/350 (75%)	244 (93%)	14 (5%)	5 (2%)	10	16
2	B	773/785 (98%)	695 (90%)	63 (8%)	15 (2%)	10	16
All	All	1036/1135 (91%)	939 (91%)	77 (7%)	20 (2%)	10	16

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	138	ASN
2	B	488	VAL
2	B	664	GLU
1	A	338	GLY
2	B	708	GLY
1	A	94	SER
1	A	330	PRO
2	B	128	TYR
2	B	199	PRO
2	B	525	ARG
2	B	702	PRO
2	B	735	PRO
2	B	770	GLU
1	A	273	GLU
2	B	737	PRO
2	B	752	PRO
2	B	486	ARG
2	B	645	GLU
2	B	315	GLY
2	B	644	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	213/277 (77%)	176 (83%)	37 (17%)	2	4
2	B	623/630 (99%)	493 (79%)	130 (21%)	1	2
All	All	836/907 (92%)	669 (80%)	167 (20%)	1	3

All (167) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	95	LEU
1	A	122	VAL
1	A	129	SER
1	A	143	HIS

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Mol	Chain	Res	Type
1	A	146	ARG
1	A	147	ASP
1	A	154	LEU
1	A	157	GLU
1	A	168	GLU
1	A	170	GLU
1	A	174	LEU
1	A	177	THR
1	A	178	HIS
1	A	183	GLN
1	A	201	ARG
1	A	204	ARG
1	A	211	THR
1	A	230	MET
1	A	245	PHE
1	A	249	SER
1	A	250	LYS
1	A	252	ARG
1	A	260	PHE
1	A	266	GLN
1	A	276	LYS
1	A	280	LEU
1	A	289	LYS
1	A	292	GLN
1	A	299	GLU
1	A	303	LEU
1	A	317	LEU
1	A	326	ARG
1	A	339	ARG
1	A	340	LEU
1	A	341	LYS
1	A	347	LYS
1	A	350	LEU
2	B	2	ARG
2	B	24	ARG
2	B	33	ASP
2	B	37	ARG
2	B	38	VAL
2	B	43	ARG
2	B	46	VAL
2	B	49	ARG
2	B	56	ILE

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Mol	Chain	Res	Type
2	B	60	ARG
2	B	61	LEU
2	B	63	ARG
2	B	70	ARG
2	B	89	LEU
2	B	95	GLU
2	B	96	LEU
2	B	99	LEU
2	B	106	ARG
2	B	108	ILE
2	B	109	GLN
2	B	111	VAL
2	B	118	LEU
2	B	127	GLU
2	B	144	THR
2	B	148	GLU
2	B	152	GLU
2	B	153	GLU
2	B	157	ASP
2	B	161	THR
2	B	171	LEU
2	B	188	GLU
2	B	196	LEU
2	B	203	LYS
2	B	208	GLU
2	B	222	ARG
2	B	226	SER
2	B	230	MET
2	B	239	MET
2	B	243	ASN
2	B	256	ARG
2	B	283	LEU
2	B	284	LYS
2	B	285	THR
2	B	286	LEU
2	B	293	LEU
2	B	294	HIS
2	B	298	LEU
2	B	307	GLU
2	B	313	LEU
2	B	322	SER
2	B	326	GLU

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Mol	Chain	Res	Type
2	B	329	GLU
2	B	331	ILE
2	B	339	ASP
2	B	345	LYS
2	B	357	SER
2	B	361	GLU
2	B	364	VAL
2	B	397	SER
2	B	399	LYS
2	B	438	GLU
2	B	441	THR
2	B	451	LEU
2	B	459	LEU
2	B	461	GLU
2	B	484	ASP
2	B	486	ARG
2	B	494	LYS
2	B	497	ARG
2	B	499	ARG
2	B	512	THR
2	B	523	ARG
2	B	525	ARG
2	B	526	LEU
2	B	541	LYS
2	B	545	ARG
2	B	548	LEU
2	B	556	LEU
2	B	558	GLU
2	B	562	LEU
2	B	567	ARG
2	B	570	LEU
2	B	571	PHE
2	B	575	ARG
2	B	579	GLU
2	B	588	LEU
2	B	590	PHE
2	B	592	GLU
2	B	600	LYS
2	B	602	ARG
2	B	603	LEU
2	B	609	LEU
2	B	613	LEU

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Mol	Chain	Res	Type
2	B	624	PHE
2	B	633	PHE
2	B	634	LEU
2	B	635	HIS
2	B	638	VAL
2	B	639	SER
2	B	645	GLU
2	B	648	GLU
2	B	649	VAL
2	B	661	GLN
2	B	662	GLU
2	B	663	LEU
2	B	670	LEU
2	B	673	LEU
2	B	674	ARG
2	B	682	LEU
2	B	685	GLN
2	B	688	SER
2	B	689	ARG
2	B	700	VAL
2	B	701	VAL
2	B	707	TYR
2	B	713	LEU
2	B	715	ARG
2	B	716	GLU
2	B	727	LEU
2	B	728	PHE
2	B	730	LEU
2	B	731	TYR
2	B	732	GLN
2	B	738	GLU
2	B	741	LYS
2	B	746	HIS
2	B	748	ARG
2	B	767	ARG
2	B	770	GLU
2	B	775	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	101	HIS

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Mol	Chain	Res	Type
1	A	120	GLN
1	A	138	ASN
1	A	142	HIS
1	A	178	HIS
1	A	183	GLN
1	A	218	GLN
1	A	232	HIS
1	A	287	HIS
2	B	54	HIS
2	B	109	GLN
2	B	178	HIS
2	B	212	HIS
2	B	231	GLN
2	B	243	ASN
2	B	258	GLN
2	B	261	HIS
2	B	294	HIS
2	B	350	HIS
2	B	449	HIS
2	B	656	HIS
2	B	751	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	FYA	A	1002	-	31,36,36	1.23	2 (6%)	30,52,52	0.69	1 (3%)

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	FYA	A	1002	-	-	0/16/36/36	0/4/4/4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1002	FYA	O3P-C	-4.19	1.27	1.44
4	A	1002	FYA	O4'-C1'	2.81	1.44	1.41

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1002	FYA	C2'-C1'-N9	-2.16	111.00	114.29

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1002	FYA	4	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	265/350 (75%)	0.12	8 (3%) 54 59	27, 54, 103, 129	0
2	B	775/785 (98%)	0.14	45 (5%) 26 30	21, 59, 108, 128	0
All	All	1040/1135 (91%)	0.13	53 (5%) 32 36	21, 58, 107, 129	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	149	TRP	5.8
2	B	718	ALA	5.3
1	A	350	LEU	5.1
2	B	739	GLY	5.0
2	B	753	LYS	5.0
2	B	736	LEU	4.7
1	A	142	HIS	4.5
2	B	732	GLN	4.5
2	B	700	VAL	4.4
2	B	99	LEU	4.3
2	B	731	TYR	4.3
2	B	743	LEU	4.3
2	B	697	LEU	4.1
2	B	768	VAL	4.0
1	A	148	MET	4.0
2	B	759	GLU	3.9
2	B	756	LEU	3.8
2	B	696	ASP	3.7
2	B	713	LEU	3.6
2	B	689	ARG	3.5
1	A	158	GLY	3.4
2	B	100	GLY	3.3
2	B	638	VAL	3.3
2	B	693	ALA	3.3

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Mol	Chain	Res	Type	RSRZ
2	B	699	VAL	3.2
2	B	733	GLY	3.2
2	B	772	LEU	3.1
2	B	701	VAL	3.1
2	B	633	PHE	2.9
2	B	771	ALA	2.9
2	B	705	THR	2.8
1	A	207	GLN	2.8
2	B	752	PRO	2.8
2	B	692	ALA	2.7
2	B	734	PRO	2.7
2	B	282	ARG	2.7
2	B	694	PHE	2.7
2	B	761	VAL	2.5
2	B	745	PHE	2.5
1	A	143	HIS	2.5
1	A	271	TRP	2.4
2	B	704	PRO	2.4
2	B	738	GLU	2.3
2	B	751	HIS	2.3
2	B	749	PHE	2.2
2	B	775	ARG	2.2
2	B	757	ARG	2.1
2	B	98	GLY	2.1
2	B	637	GLY	2.1
2	B	740	HIS	2.1
2	B	659	ILE	2.0
2	B	762	GLU	2.0
2	B	525	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	MG	A	1001	1/1	0.84	0.26	9.05	42,42,42,42	0
4	FYA	A	1002	33/33	0.97	0.15	-0.01	40,65,73,76	0

6.5 Other polymers [i](#)

There are no such residues in this entry.