Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 03:13 am BST

PDB ID	:	1KT8
Title	:	HUMAN BRANCHED CHAIN AMINO ACID AMINOTRANSFERASE (MITOCHONDRIAL): THREE DIMENSIONAL STRUCTURE OF ENZYME IN ITS KETIMINE FORM WITH THE SUBSTRATE L-ISOLEUCINE
Authors	:	Yennawar, N.H.; Conway, M.E.; Yennawar, H.P.; Farber, G.K.; Hutson, S.M.
Deposited on	:	2002-01-15
Resolution	:	1.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

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

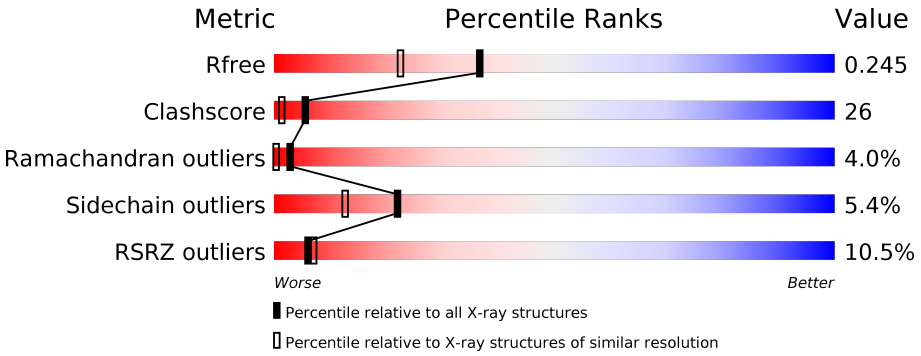
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	365	10% 64% 28% 5%
1	B	365	11% 65% 27% 6%

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
2	ILP	A	400	X	-	X	-

2 Entry composition

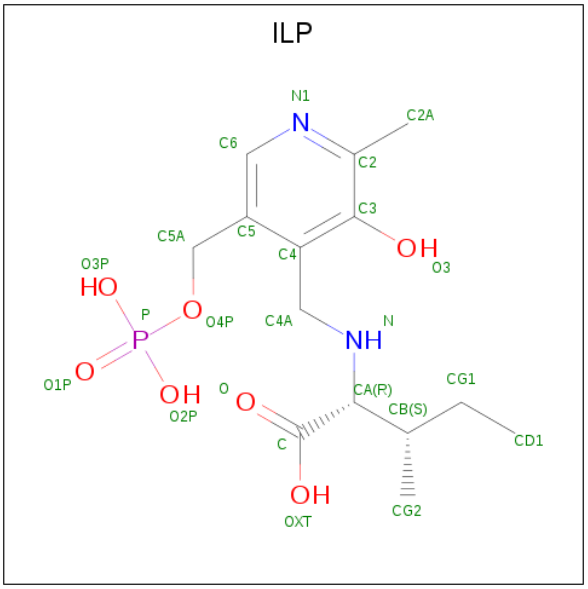
There are 5 unique types of molecules in this entry. The entry contains 6135 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 BRANCHED-CHAIN AMINO ACID AMINOTRANSFERASE, MITOCHONDRIAL.

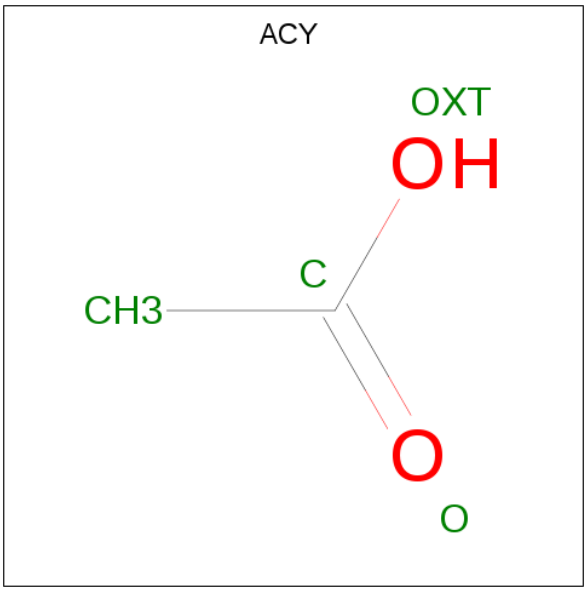
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	360	Total	C	N	O	S	0	3	0
			2904	1869	508	509	18			
1	B	365	Total	C	N	O	S	0	2	0
			2932	1887	515	512	18			

- Molecule 2 is N-[O-PHOSPHONO-PYRIDOXYL]-ISOLEUCINE (three-letter code: ILP) (formula: $C_{14}H_{23}N_2O_7P$).



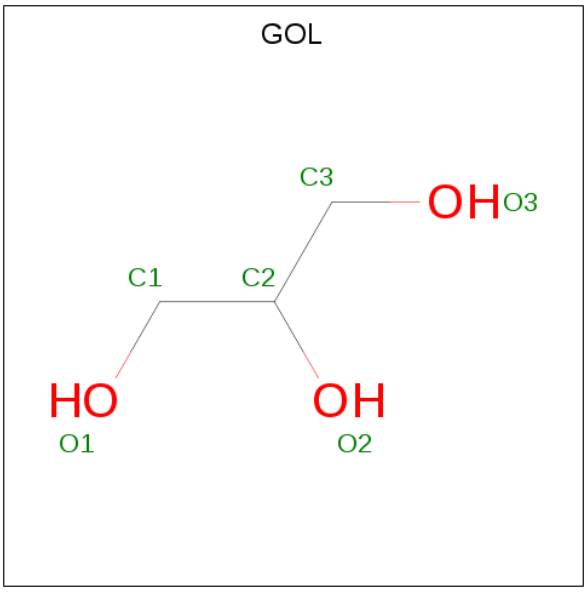
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			24	14	2	7	1		
2	B	1	Total	C	N	O	P	0	0
			24	14	2	7	1		

- Molecule 3 is ACETIC ACID (three-letter code: ACY) (formula: $C_2H_4O_2$).



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

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C O 6 3 3	0	0

- Molecule 5 is water.

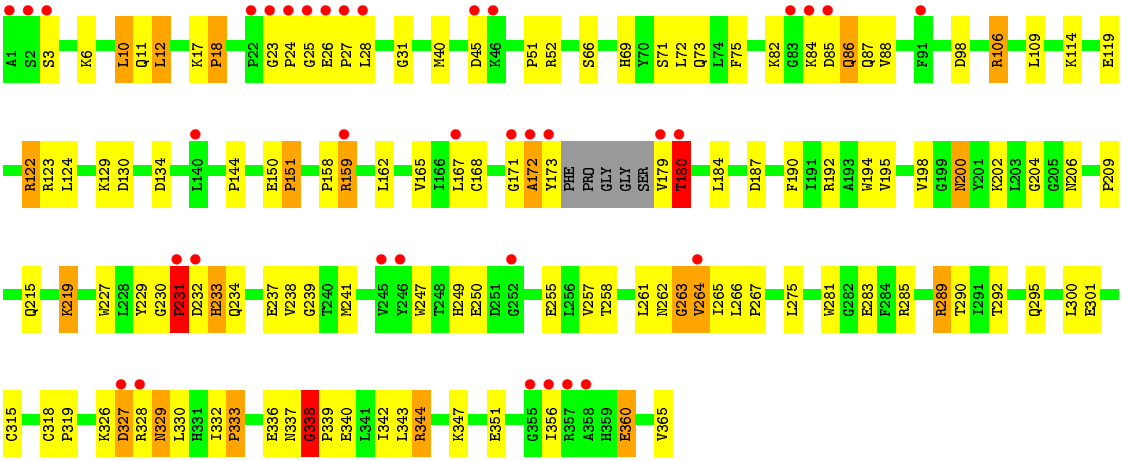
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	110	Total O 110 110	0	0
5	B	111	Total O 111 111	0	0

3 Residue-property plots (i)

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

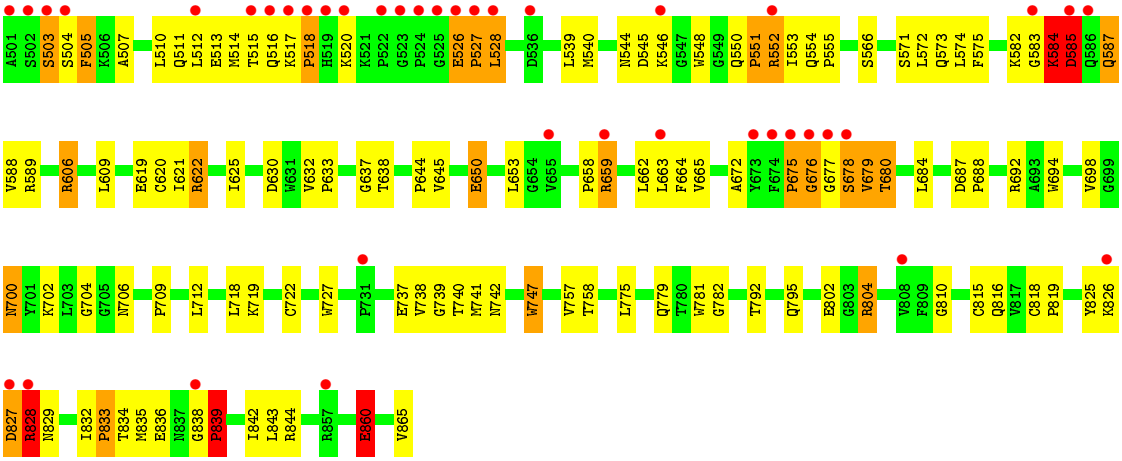
• Molecule 1: BRANCHED-CHAIN AMINO ACID AMINOTRANSFERASE, MITOCHONDRIAL

Chain A:



• Molecule 1: BRANCHED-CHAIN AMINO ACID AMINOTRANSFERASE, MITOCHONDRIAL

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	69.47Å 105.46Å 107.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.72 – 1.90 19.72 – 1.88	Depositor EDS
% Data completeness (in resolution range)	95.6 (19.72-1.90) 96.0 (19.72-1.88)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.14 (at 1.89Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.243 , 0.284 0.240 , 0.245	Depositor DCC
R_{free} test set	3128 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	30.9	Xtriage
Anisotropy	0.397	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 43.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.016 for -h,l,k	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	6135	wwPDB-VP
Average B, all atoms (Å^2)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, ACY, ILP

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	0.81	3/2978 (0.1%)	1.04	19/4039 (0.5%)
1	B	0.93	9/3009 (0.3%)	1.02	14/4081 (0.3%)
All	All	0.87	12/5987 (0.2%)	1.03	33/8120 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	3
All	All	0	6

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	828[A]	ARG	CB-CG	15.29	1.93	1.52
1	B	828[B]	ARG	CB-CG	15.29	1.93	1.52
1	B	828[A]	ARG	CG-CD	9.45	1.75	1.51
1	B	828[B]	ARG	CG-CD	9.45	1.75	1.51
1	A	168	CYS	CB-SG	-7.29	1.69	1.82
1	A	264[A]	VAL	CA-C	-5.95	1.37	1.52
1	A	264[B]	VAL	CA-C	-5.95	1.37	1.52
1	B	828[A]	ARG	CD-NE	5.47	1.55	1.46
1	B	828[B]	ARG	CD-NE	5.47	1.55	1.46
1	B	828[A]	ARG	N-CA	-5.23	1.35	1.46
1	B	828[B]	ARG	N-CA	-5.23	1.35	1.46
1	B	860	GLU	CB-CG	-5.10	1.42	1.52

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	231	PRO	CA-C-N	-11.24	92.46	117.20
1	B	606	ARG	NE-CZ-NH1	10.77	125.68	120.30
1	B	606	ARG	NE-CZ-NH2	-10.19	115.21	120.30
1	A	263	GLY	O-C-N	9.74	138.28	122.70
1	A	263	GLY	CA-C-N	-9.13	97.10	117.20
1	B	827	ASP	CA-C-N	-8.97	97.47	117.20
1	A	338	GLY	C-N-CD	8.92	147.13	128.40
1	A	327	ASP	CA-C-N	-8.89	97.64	117.20
1	B	828[A]	ARG	CB-CA-C	8.27	126.93	110.40
1	B	828[B]	ARG	CB-CA-C	8.27	126.93	110.40
1	B	827	ASP	O-C-N	8.24	135.88	122.70
1	B	828[A]	ARG	CG-CD-NE	8.19	129.00	111.80
1	B	828[B]	ARG	CG-CD-NE	8.19	129.00	111.80
1	B	828[A]	ARG	NE-CZ-NH1	8.12	124.36	120.30
1	B	828[B]	ARG	NE-CZ-NH1	8.12	124.36	120.30
1	A	106	ARG	NE-CZ-NH2	-8.05	116.28	120.30
1	A	106	ARG	NE-CZ-NH1	7.72	124.16	120.30
1	A	231	PRO	O-C-N	7.68	134.99	122.70
1	A	263	GLY	C-N-CA	7.39	140.17	121.70
1	A	231	PRO	C-N-CA	7.20	139.69	121.70
1	A	264[A]	VAL	N-CA-CB	7.20	127.33	111.50
1	A	264[B]	VAL	N-CA-CB	7.20	127.33	111.50
1	A	264[A]	VAL	CB-CA-C	-6.46	99.12	111.40
1	A	264[B]	VAL	CB-CA-C	-6.46	99.12	111.40
1	B	828[A]	ARG	CB-CG-CD	6.40	128.25	111.60
1	B	828[B]	ARG	CB-CG-CD	6.40	128.25	111.60
1	A	264[A]	VAL	CG1-CB-CG2	-6.28	100.86	110.90
1	A	264[B]	VAL	CG1-CB-CG2	-6.28	100.86	110.90
1	A	338	GLY	C-N-CA	-6.01	96.78	122.00
1	B	804	ARG	NE-CZ-NH2	-5.40	117.60	120.30
1	B	503	SER	N-CA-C	5.21	125.08	111.00
1	A	344	ARG	NE-CZ-NH1	5.04	122.82	120.30
1	A	289	ARG	NE-CZ-NH2	-5.00	117.80	120.30

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	231	PRO	Mainchain
1	A	263	GLY	Mainchain
1	A	327	ASP	Mainchain
1	B	551	PRO	Mainchain
1	B	827	ASP	Mainchain

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	2904	0	2935	147	0
1	B	2932	0	2961	169	0
2	A	24	0	19	9	0
2	B	24	0	19	6	0
3	A	12	0	9	1	0
3	B	12	0	9	1	0
4	B	6	0	8	1	0
5	A	110	0	0	8	0
5	B	111	0	0	4	0
All	All	6135	0	5960	307	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 26.

All (307) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:828[A]:ARG:CG	1:B:828[A]:ARG:CD	1.75	1.58
1:B:828[A]:ARG:CG	1:B:828[A]:ARG:CB	1.93	1.43
1:A:264[A]:VAL:O	1:A:265:ILE:HG12	1.37	1.20
1:A:262:ASN:OD1	1:A:264[A]:VAL:HB	1.43	1.17
1:A:338:GLY:O	1:A:340:GLU:N	1.79	1.13
1:A:264[A]:VAL:HG12	1:A:265:ILE:HG23	1.13	1.12
1:B:552[A]:ARG:NH2	1:B:554:GLN:NE2	2.07	1.02
1:B:552[A]:ARG:CZ	1:B:554:GLN:HE21	1.72	1.02
1:A:264[A]:VAL:O	1:A:265:ILE:CG1	2.15	0.94
1:B:511:GLN:H	1:B:552[A]:ARG:HA	1.32	0.93
1:B:515:THR:HG23	1:B:518:PRO:HG3	1.49	0.93
1:A:200:ASN:H	1:A:200:ASN:HD22	1.06	0.92
1:B:700:ASN:H	1:B:700:ASN:HD22	1.04	0.92
1:A:262:ASN:OD1	1:A:264[A]:VAL:CB	2.18	0.91
1:B:511:GLN:HB2	1:B:552[A]:ARG:HG3	1.53	0.91
1:B:584:LYS:HD2	1:B:587:GLN:HB2	1.51	0.91
1:B:511:GLN:O	1:B:552[A]:ARG:HG2	1.71	0.90
1:A:10:LEU:HD13	1:A:51:PRO:HB2	1.52	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:584:LYS:O	1:B:584:LYS:HG2	1.70	0.89
1:A:202:LYS:HZ1	2:A:400:ILP:HA	1.38	0.89
1:B:552[A]:ARG:CZ	1:B:554:GLN:NE2	2.34	0.89
1:A:249:HIS:CD2	1:A:250:GLU:H	1.92	0.87
1:B:552[A]:ARG:HD2	1:B:553:ILE:N	1.90	0.87
1:A:264[A]:VAL:CG1	1:A:265:ILE:HG23	2.01	0.86
1:A:232[A]:ASP:O	1:A:234:GLN:N	2.09	0.85
1:B:552[A]:ARG:HH22	1:B:554:GLN:NE2	1.75	0.85
1:B:702:LYS:NZ	2:B:900:ILP:HN2	1.74	0.84
1:B:702:LYS:HZ1	2:B:900:ILP:HN2	1.22	0.84
1:B:583:GLY:C	1:B:585:ASP:H	1.75	0.83
1:B:700:ASN:N	1:B:700:ASN:HD22	1.78	0.81
1:A:194:TRP:C	1:A:264[B]:VAL:HG12	2.00	0.81
1:B:520:LYS:HB3	1:B:520:LYS:NZ	1.96	0.81
1:A:123:ARG:NH1	5:A:1101:HOH:O	2.15	0.80
1:A:328[B]:ARG:O	1:A:329:ASN:HB2	1.81	0.80
1:B:552[A]:ARG:O	1:B:553:ILE:HB	1.83	0.79
1:B:511:GLN:HG2	1:B:552[A]:ARG:HB2	1.66	0.78
1:A:339:PRO:HB2	1:A:342:ILE:HB	1.65	0.78
1:B:825:TYR:O	1:B:828[A]:ARG:HB2	1.85	0.76
1:B:540:MET:O	1:B:552[B]:ARG:HB2	1.84	0.76
1:B:700:ASN:ND2	1:B:700:ASN:H	1.82	0.76
1:A:194:TRP:C	1:A:264[B]:VAL:CG1	2.55	0.75
1:B:583:GLY:O	1:B:585:ASP:N	2.19	0.75
1:A:71:SER:H	1:B:573:GLN:HE22	1.34	0.75
1:B:527:PRO:C	1:B:528:LEU:HD23	2.08	0.74
1:A:28:LEU:HB2	1:A:171:GLY:HA3	1.70	0.74
1:B:828[B]:ARG:O	1:B:828[B]:ARG:CD	2.36	0.74
1:B:839:PRO:HB3	1:B:842:ILE:HB	1.70	0.74
1:A:202:LYS:HZ1	2:A:400:ILP:HG23	1.50	0.73
1:B:637:GLY:O	1:B:672:ALA:HA	1.88	0.73
1:B:510:LEU:HD23	1:B:551:PRO:HB2	1.70	0.73
1:B:552[A]:ARG:NH1	1:B:554:GLN:NE2	2.37	0.73
1:B:676:GLY:C	1:B:678:SER:H	1.91	0.73
1:A:200:ASN:H	1:A:200:ASN:ND2	1.85	0.73
1:B:552[A]:ARG:O	1:B:553:ILE:CB	2.37	0.73
1:B:692:ARG:NH1	1:B:737:GLU:OE2	2.23	0.72
1:A:12:LEU:HD12	1:A:12:LEU:O	1.90	0.72
1:A:202:LYS:NZ	2:A:400:ILP:HA	2.03	0.72
1:B:552[A]:ARG:NH1	1:B:554:GLN:HB3	2.04	0.72
1:B:584:LYS:HZ2	1:B:587:GLN:HG3	1.53	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264[A]:VAL:HG12	1:A:265:ILE:CG2	2.07	0.72
1:B:511:GLN:H	1:B:552[A]:ARG:CA	2.03	0.71
1:B:835:MET:HA	1:B:839:PRO:HD3	1.72	0.71
1:B:511:GLN:HB2	1:B:552[A]:ARG:CG	2.20	0.70
1:B:828[B]:ARG:HD3	1:B:828[B]:ARG:O	1.92	0.69
1:A:281:TRP:CH2	1:A:344:ARG:HG2	2.28	0.69
1:A:10:LEU:CD1	1:A:51:PRO:HB2	2.21	0.68
1:A:202:LYS:NZ	2:A:400:ILP:HG23	2.07	0.68
1:A:264[A]:VAL:CG1	5:A:1013:HOH:O	2.42	0.68
1:A:82:LYS:HD3	1:A:134:ASP:HB3	1.76	0.68
1:B:552[A]:ARG:HH22	1:B:554:GLN:HE22	1.41	0.68
1:B:540:MET:SD	1:B:662:LEU:HD11	2.33	0.68
1:A:289:ARG:HG2	1:A:290:THR:N	2.08	0.67
1:B:740:THR:CG2	2:B:900:ILP:HG23	2.24	0.67
1:B:835:MET:HA	1:B:839:PRO:CD	2.24	0.67
1:B:584:LYS:NZ	1:B:587:GLN:HG3	2.09	0.67
1:A:73:GLN:HE22	1:B:571:SER:H	1.42	0.67
1:A:184:LEU:HD22	1:A:238:VAL:HG23	1.77	0.66
1:A:337:ASN:O	1:A:338:GLY:O	2.12	0.66
1:A:171:GLY:O	1:A:172:ALA:CB	2.43	0.66
1:A:200:ASN:HD22	1:A:200:ASN:N	1.80	0.66
1:A:106:ARG:NH2	1:B:709:PRO:O	2.28	0.66
1:A:202:LYS:HZ1	2:A:400:ILP:CG2	2.08	0.66
1:A:232[B]:ASP:O	1:A:233:HIS:HB2	1.95	0.66
1:B:520:LYS:HZ2	1:B:520:LYS:HB3	1.59	0.66
1:B:511:GLN:CG	1:B:552[A]:ARG:HB2	2.25	0.65
1:A:202:LYS:HZ1	2:A:400:ILP:CA	2.08	0.65
1:B:526:GLU:HB3	1:B:527:PRO:HD3	1.79	0.64
1:B:676:GLY:O	1:B:678:SER:N	2.31	0.64
1:B:552[A]:ARG:NH2	1:B:554:GLN:HE22	1.93	0.64
1:B:584:LYS:HE3	1:B:589:ARG:HH22	1.63	0.63
1:B:740:THR:HG21	2:B:900:ILP:CG2	2.28	0.63
1:B:552[A]:ARG:NH2	1:B:554:GLN:HE21	1.80	0.62
1:B:511:GLN:C	1:B:552[A]:ARG:HG2	2.20	0.62
1:A:229:TYR:O	1:A:232[A]:ASP:O	2.17	0.62
1:A:264[A]:VAL:O	1:A:265:ILE:CB	2.45	0.62
1:A:40:MET:SD	1:A:162:LEU:HD11	2.39	0.62
1:A:239:GLY:HA2	5:A:1044:HOH:O	2.00	0.62
1:A:338:GLY:C	1:A:340:GLU:N	2.53	0.61
1:A:26:GLU:N	1:A:27:PRO:HD3	2.15	0.61
1:B:572:LEU:HD13	1:B:658:PRO:HG3	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:LEU:HD13	1:A:158:PRO:HG3	1.82	0.61
1:B:511:GLN:N	1:B:552[A]:ARG:HA	2.09	0.61
1:A:184:LEU:HD22	1:A:238:VAL:CG2	2.30	0.61
1:A:264[A]:VAL:HG11	5:A:1013:HOH:O	2.01	0.61
1:A:12:LEU:C	1:A:12:LEU:HD12	2.20	0.61
1:A:262:ASN:HD21	1:A:264[A]:VAL:CG2	2.13	0.61
1:B:552[A]:ARG:NH1	1:B:554:GLN:HE21	1.96	0.61
1:B:678:SER:O	1:B:679:VAL:C	2.40	0.60
1:B:584:LYS:HE3	1:B:589:ARG:NH2	2.17	0.60
1:A:264[A]:VAL:CB	5:A:1013:HOH:O	2.49	0.59
1:B:815:CYS:SG	1:B:818:CYS:HB2	2.41	0.59
1:A:264[A]:VAL:HG12	1:A:264[A]:VAL:O	2.01	0.59
1:B:781:TRP:CH2	1:B:844:ARG:HG2	2.37	0.59
1:A:292:THR:OG1	1:A:295:GLN:HG3	2.03	0.59
1:B:512:LEU:HD22	1:B:514:MET:CE	2.33	0.58
1:A:11:GLN:NE2	1:A:52:ARG:HH11	2.03	0.57
1:A:159:ARG:HH11	1:A:159:ARG:CB	2.18	0.57
1:B:650:GLU:CG	1:B:658:PRO:HA	2.34	0.57
1:A:365:VAL:HG23	1:A:365:VAL:OXT	2.05	0.57
1:B:834:THR:O	1:B:839:PRO:HD3	2.05	0.56
1:A:66:SER:HB2	1:A:72:LEU:HD12	1.88	0.56
1:A:85:ASP:O	1:A:86:GLN:HB2	2.06	0.56
1:B:692:ARG:CZ	1:B:737:GLU:OE2	2.53	0.56
1:A:339:PRO:CB	1:A:342:ILE:HB	2.35	0.56
1:B:650:GLU:HG2	1:B:658:PRO:HA	1.88	0.55
1:B:684:LEU:HD22	1:B:738:VAL:HG23	1.88	0.55
1:B:552[A]:ARG:HD2	1:B:553:ILE:C	2.27	0.55
1:B:676:GLY:C	1:B:678:SER:N	2.59	0.55
1:A:262:ASN:OD1	1:A:264[B]:VAL:HG23	2.06	0.55
1:B:622:ARG:HG2	1:B:865:VAL:OXT	2.05	0.55
1:A:264[B]:VAL:HA	5:A:1070:HOH:O	2.06	0.55
1:B:573:GLN:HE21	1:B:704:GLY:HA3	1.70	0.55
1:A:179:VAL:HG22	1:A:180:THR:N	2.22	0.55
1:A:264[A]:VAL:HB	5:A:1013:HOH:O	2.07	0.55
1:B:539:LEU:HA	1:B:552[B]:ARG:O	2.06	0.55
1:A:28:LEU:HD12	1:A:171:GLY:HA2	1.88	0.54
1:A:300:LEU:O	1:A:301:GLU:C	2.46	0.54
1:B:511:GLN:NE2	1:B:550:GLN:OE1	2.41	0.54
1:B:552[A]:ARG:CZ	1:B:554:GLN:HB3	2.38	0.54
1:A:150:GLU:HG2	1:A:151:PRO:HD2	1.89	0.54
1:A:209:PRO:O	1:B:606:ARG:NH2	2.39	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:507:ALA:O	1:B:510:LEU:HD12	2.08	0.54
1:B:513:GLU:HG2	1:B:555:PRO:HD3	1.89	0.54
1:B:583:GLY:C	1:B:585:ASP:N	2.51	0.54
1:A:24:PRO:O	1:A:26:GLU:N	2.41	0.54
1:A:73:GLN:HE21	1:A:204:GLY:HA3	1.73	0.54
1:A:195:VAL:N	1:A:264[B]:VAL:CG1	2.71	0.53
1:B:510:LEU:CD2	1:B:551:PRO:HB2	2.38	0.53
1:A:122:ARG:CG	1:A:365:VAL:OXT	2.56	0.53
1:B:552[A]:ARG:NH1	1:B:554:GLN:CB	2.71	0.53
1:B:826:LYS:C	1:B:828[A]:ARG:H	2.08	0.53
1:A:171:GLY:O	1:A:172:ALA:HB3	2.08	0.53
1:B:566:SER:HB2	1:B:572:LEU:HD12	1.90	0.53
1:B:588:VAL:O	1:B:865:VAL:N	2.39	0.53
1:A:249:HIS:CD2	1:A:250:GLU:N	2.71	0.53
1:B:792:THR:OG1	1:B:795:GLN:HG3	2.08	0.53
1:A:144:PRO:HA	1:A:165:VAL:HG22	1.91	0.53
1:A:262:ASN:HD21	1:A:264[A]:VAL:HG23	1.74	0.53
1:B:511:GLN:CB	1:B:552[A]:ARG:HB2	2.39	0.53
1:A:261:LEU:HD13	3:A:1003:ACY:H3	1.91	0.53
1:A:71:SER:H	1:B:573:GLN:NE2	2.03	0.52
1:A:159:ARG:HH11	1:A:159:ARG:HB3	1.74	0.52
1:A:194:TRP:CA	1:A:264[B]:VAL:HG12	2.39	0.52
1:A:232[A]:ASP:CG	1:A:234:GLN:OE1	2.48	0.52
1:B:619:GLU:OE1	1:B:622:ARG:NH1	2.42	0.52
1:B:659:ARG:CB	1:B:659:ARG:HH11	2.23	0.52
1:B:515:THR:CG2	1:B:518:PRO:HG3	2.31	0.52
1:B:816:GLN:NE2	5:B:144:HOH:O	2.42	0.52
1:A:192:ARG:HB2	1:A:227:TRP:CE3	2.44	0.52
1:B:511:GLN:H	1:B:552[A]:ARG:CB	2.23	0.52
1:A:11:GLN:HE21	1:A:52:ARG:HH11	1.57	0.52
1:B:552[A]:ARG:NH1	1:B:554:GLN:CG	2.73	0.51
1:A:184:LEU:CD2	1:A:238:VAL:CG2	2.89	0.51
1:B:510:LEU:HD21	1:B:620:CYS:SG	2.51	0.51
1:B:835:MET:SD	1:B:839:PRO:HG2	2.51	0.51
1:A:119:GLU:OE1	1:A:122:ARG:NH1	2.43	0.51
1:A:202:LYS:CE	2:A:400:ILP:HA	2.40	0.50
1:A:122:ARG:HG2	1:A:365:VAL:OXT	2.10	0.50
1:B:622:ARG:CG	1:B:865:VAL:OXT	2.59	0.50
1:A:31:GLY:HA3	1:B:653:LEU:HB2	1.93	0.50
1:A:88:VAL:O	1:A:365:VAL:N	2.44	0.50
1:B:662:LEU:HD22	3:B:1001:ACY:H3	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:838:GLY:O	1:B:839:PRO:C	2.47	0.50
1:A:339:PRO:HB3	1:A:342:ILE:HD12	1.94	0.50
1:B:828[B]:ARG:CG	1:B:828[B]:ARG:O	2.58	0.50
1:A:98:ASP:OD2	1:A:114:LYS:HE3	2.11	0.50
1:A:283:GLU:CD	1:A:344:ARG:HH22	2.15	0.50
1:B:684:LEU:HD22	1:B:738:VAL:CG2	2.42	0.49
1:A:232[A]:ASP:OD1	1:A:234:GLN:OE1	2.30	0.49
1:B:659:ARG:HB3	1:B:659:ARG:HH11	1.77	0.49
1:A:82:LYS:HE2	1:A:86:GLN:CD	2.33	0.49
1:B:692:ARG:HB2	1:B:727:TRP:CE3	2.47	0.49
1:B:517:LYS:N	1:B:518:PRO:HD3	2.27	0.49
1:B:505:PHE:CE1	1:B:551:PRO:HG3	2.48	0.49
1:B:834:THR:HG22	1:B:839:PRO:HG3	1.95	0.49
1:B:679:VAL:HG12	1:B:819:PRO:HG2	1.95	0.49
1:B:860:GLU:H	1:B:860:GLU:HG3	1.46	0.49
1:B:865:VAL:OXT	1:B:865:VAL:HG23	2.13	0.49
1:B:544:ASN:OD1	1:B:546:LYS:N	2.42	0.48
1:A:179:VAL:O	1:A:180:THR:HG23	2.14	0.48
1:B:582:LYS:HE3	1:B:632:VAL:HB	1.94	0.48
1:A:124:LEU:HG	1:A:167:LEU:HD11	1.95	0.48
1:B:663:LEU:HD23	1:B:664:PHE:N	2.29	0.48
1:A:215:GLN:O	1:A:219:LYS:HG2	2.13	0.48
1:B:511:GLN:O	1:B:552[A]:ARG:O	2.31	0.48
1:A:200:ASN:ND2	1:A:200:ASN:N	2.53	0.48
1:B:520:LYS:HB3	1:B:520:LYS:HZ3	1.78	0.48
1:B:698:VAL:HG23	1:B:706:ASN:HD21	1.77	0.47
1:A:184:LEU:CD2	1:A:238:VAL:HG21	2.45	0.47
1:A:122:ARG:HE	1:A:365:VAL:HG23	1.79	0.47
1:A:315:CYS:SG	1:A:318:CYS:HB2	2.54	0.47
1:A:347:LYS:NZ	1:A:351:GLU:OE2	2.42	0.47
1:B:740:THR:HB	2:B:900:ILP:HG23	1.96	0.47
1:A:360:GLU:HG3	1:A:360:GLU:H	1.51	0.47
1:B:678:SER:O	1:B:680:THR:HG23	2.15	0.47
1:B:574:LEU:HD12	1:B:574:LEU:C	2.35	0.46
1:B:739:GLY:HA2	5:B:126:HOH:O	2.15	0.46
1:B:833:PRO:O	1:B:836:GLU:HG2	2.16	0.46
1:A:230:GLY:C	1:A:232[A]:ASP:N	2.55	0.46
1:B:552[A]:ARG:HH12	1:B:554:GLN:NE2	2.09	0.46
1:A:262:ASN:ND2	1:A:264[A]:VAL:HG23	2.31	0.46
1:A:328[A]:ARG:HG2	1:A:330:LEU:HD21	1.97	0.46
1:B:512:LEU:HD22	1:B:514:MET:HE3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:515:THR:HG23	1:B:518:PRO:CG	2.35	0.46
1:B:650:GLU:HG3	1:B:658:PRO:HA	1.98	0.46
1:B:572:LEU:HA	1:B:572:LEU:HD23	1.84	0.46
1:B:757:VAL:HG22	1:B:758:THR:N	2.31	0.45
1:B:644:PRO:HA	1:B:665:VAL:HG22	1.97	0.45
1:B:832:ILE:HA	1:B:833:PRO:HD3	1.77	0.45
1:A:257:VAL:HG22	1:A:258:THR:N	2.32	0.45
1:A:264[A]:VAL:C	1:A:265:ILE:HG23	2.38	0.45
1:A:109:LEU:HD23	1:A:109:LEU:HA	1.90	0.45
1:A:73:GLN:NE2	1:B:571:SER:H	2.10	0.45
1:B:737:GLU:HB2	1:B:741:MET:O	2.17	0.45
1:B:782:GLY:HA2	5:B:95:HOH:O	2.15	0.45
1:A:69:HIS:CE1	1:B:645:VAL:HG21	2.53	0.44
1:B:546:LYS:HB2	1:B:546:LYS:HE2	1.75	0.44
1:B:712:LEU:HG	4:B:2001:GOL:H12	2.00	0.44
1:B:747:TRP:C	1:B:747:TRP:CD1	2.89	0.44
1:A:11:GLN:HE21	1:A:52:ARG:NH1	2.15	0.44
1:A:28:LEU:CB	1:A:171:GLY:HA3	2.45	0.44
1:B:802:GLU:HB2	1:B:804:ARG:HD3	1.99	0.44
1:A:202:LYS:NZ	2:A:400:ILP:N	2.65	0.44
1:B:633:PRO:HB2	1:B:638:THR:OG1	2.18	0.44
1:B:825:TYR:O	1:B:828[A]:ARG:CB	2.62	0.44
1:A:75:PHE:CE2	1:A:202:LYS:HE3	2.53	0.44
1:B:621:ILE:O	1:B:625:ILE:HG13	2.18	0.44
1:A:98:ASP:OD2	1:A:114:LYS:CE	2.66	0.44
1:B:512:LEU:HD23	1:B:553:ILE:HB	2.00	0.44
1:B:511:GLN:H	1:B:552[A]:ARG:HB2	1.83	0.43
1:B:551:PRO:O	1:B:552[A]:ARG:CB	2.66	0.43
1:B:622:ARG:HE	1:B:865:VAL:HG23	1.83	0.43
1:A:85:ASP:O	1:A:86:GLN:CB	2.66	0.43
1:B:588:VAL:O	1:B:865:VAL:HA	2.18	0.43
1:B:552[A]:ARG:HH12	1:B:554:GLN:CD	2.22	0.43
1:B:838:GLY:O	1:B:843:LEU:HD12	2.19	0.43
1:A:129:LYS:HG3	1:A:130:ASP:N	2.34	0.43
1:B:684:LEU:CD2	1:B:738:VAL:CG2	2.96	0.43
1:A:187:ASP:HB3	1:A:190:PHE:CE1	2.54	0.43
1:A:198:VAL:HG23	1:A:206:ASN:HD21	1.82	0.43
1:A:264[A]:VAL:O	1:A:265:ILE:HG23	2.19	0.43
1:A:337:ASN:C	1:A:338:GLY:O	2.57	0.43
1:A:283:GLU:OE1	1:A:344:ARG:NH2	2.52	0.43
1:B:505:PHE:HB2	1:B:548:TRP:HB3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:742:ASN:O	1:B:810:GLY:HA2	2.19	0.42
1:A:194:TRP:HB2	1:B:694:TRP:CE3	2.54	0.42
1:B:816:GLN:HB2	5:B:10:HOH:O	2.19	0.42
1:B:516:GLN:C	1:B:518:PRO:HD3	2.40	0.42
1:A:333:PRO:O	1:A:336:GLU:HG2	2.18	0.42
1:B:692:ARG:HB2	1:B:727:TRP:CD2	2.54	0.42
1:A:339:PRO:O	1:A:343:LEU:HG	2.19	0.42
1:A:88:VAL:O	1:A:365:VAL:HA	2.20	0.42
1:A:179:VAL:HG22	1:A:180:THR:H	1.84	0.42
1:A:266:LEU:HA	1:A:267:PRO:HD3	1.89	0.42
1:B:505:PHE:CD1	1:B:551:PRO:HG3	2.54	0.42
1:B:687:ASP:HA	1:B:688:PRO:HD3	1.85	0.42
1:B:718:LEU:HD23	1:B:722:CYS:O	2.20	0.42
1:B:740:THR:CB	2:B:900:ILP:HG23	2.49	0.42
1:A:315:CYS:SG	1:A:318:CYS:SG	3.17	0.41
1:A:173:TYR:CE1	1:A:315:CYS:HB2	2.55	0.41
1:B:540:MET:HB3	1:B:552[B]:ARG:HB2	2.00	0.41
1:B:825:TYR:N	1:B:828[A]:ARG:O	2.43	0.41
1:A:24:PRO:C	1:A:26:GLU:N	2.73	0.41
1:A:264[B]:VAL:H	1:A:264[B]:VAL:HG23	1.18	0.41
1:A:237:GLU:HB2	1:A:241:MET:O	2.20	0.41
1:A:249:HIS:CG	1:A:250:GLU:H	2.35	0.41
1:A:255:GLU:HB3	1:A:285:ARG:HD2	2.02	0.41
1:B:609:LEU:HA	1:B:609:LEU:HD23	1.84	0.41
1:A:17:LYS:N	1:A:18:PRO:HD3	2.36	0.41
1:A:332:ILE:HA	1:A:333:PRO:HD3	1.75	0.41
1:B:552[A]:ARG:HD3	1:B:553:ILE:O	2.20	0.41
1:B:552[A]:ARG:O	1:B:553:ILE:CG1	2.68	0.41
1:A:84:LYS:HE3	1:A:356:ILE:HD12	2.03	0.41
1:B:552[A]:ARG:CD	1:B:553:ILE:O	2.69	0.41
1:A:318:CYS:HA	1:A:319:PRO:HD3	1.99	0.41
1:A:122:ARG:HG3	1:A:365:VAL:OXT	2.21	0.41
1:A:231:PRO:CD	1:A:232[A]:ASP:H	2.07	0.41
1:A:262:ASN:ND2	1:A:264[A]:VAL:CG2	2.82	0.41
1:A:12:LEU:C	1:A:12:LEU:CD1	2.89	0.40
1:A:202:LYS:HZ1	2:A:400:ILP:CB	2.33	0.40
1:A:262:ASN:HD21	1:A:264[A]:VAL:HG21	1.84	0.40
1:B:818:CYS:HA	1:B:819:PRO:HD3	1.92	0.40
1:B:575:PHE:CE2	1:B:702:LYS:HE3	2.55	0.40
1:A:69:HIS:HB2	5:A:1024:HOH:O	2.20	0.40
1:B:550:GLN:HA	1:B:551:PRO:HD3	1.93	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:LEU:HD21	1:A:238:VAL:HG21	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	359/365 (98%)	324 (90%)	22 (6%)	13 (4%)	3 0
1	B	365/365 (100%)	332 (91%)	17 (5%)	16 (4%)	2 0
All	All	724/730 (99%)	656 (91%)	39 (5%)	29 (4%)	3 0

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	172	ALA
1	A	180	THR
1	B	503	SER
1	B	504	SER
1	B	527	PRO
1	B	584	LYS
1	B	678	SER
1	B	679	VAL
1	B	829	ASN
1	B	839	PRO
1	A	3	SER
1	A	25	GLY
1	A	86	GLN
1	A	338	GLY
1	B	677	GLY
1	A	233	HIS
1	A	329	ASN

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Mol	Chain	Res	Type
1	B	526	GLU
1	A	326	LYS
1	B	585	ASP
1	B	528	LEU
1	B	675	PRO
1	A	23	GLY
1	B	518	PRO
1	B	833	PRO
1	A	18	PRO
1	A	231	PRO
1	A	333	PRO
1	B	676	GLY

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	316/316 (100%)	302 (96%)	14 (4%)	28 19
1	B	318/316 (101%)	296 (93%)	22 (7%)	15 7
All	All	634/632 (100%)	598 (94%)	36 (6%)	22 11

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LYS
1	A	10	LEU
1	A	12	LEU
1	A	45	ASP
1	A	87	GLN
1	A	122	ARG
1	A	151	PRO
1	A	159	ARG
1	A	180	THR
1	A	200	ASN
1	A	219	LYS

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Mol	Chain	Res	Type
1	A	247	TRP
1	A	275	LEU
1	A	360	GLU
1	B	505	PHE
1	B	545	ASP
1	B	552[A]	ARG
1	B	552[B]	ARG
1	B	584	LYS
1	B	585	ASP
1	B	587	GLN
1	B	622	ARG
1	B	630	ASP
1	B	650	GLU
1	B	659	ARG
1	B	675	PRO
1	B	680	THR
1	B	700	ASN
1	B	719	LYS
1	B	747	TRP
1	B	775	LEU
1	B	779	GLN
1	B	828[A]	ARG
1	B	828[B]	ARG
1	B	839	PRO
1	B	860	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	GLN
1	A	54	GLN
1	A	73	GLN
1	A	86	GLN
1	A	96	ASN
1	A	200	ASN
1	A	206	ASN
1	A	215	GLN
1	A	224	GLN
1	A	249	HIS
1	B	554	GLN
1	B	573	GLN
1	B	587	GLN

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Mol	Chain	Res	Type
1	B	700	ASN
1	B	706	ASN
1	B	715	GLN
1	B	724	GLN
1	B	742	ASN
1	B	779	GLN
1	B	853	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

9 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	ACY	B	1002	-	1,3,3	4.01	1 (100%)	0,3,3	0.00	-
4	GOL	B	2001	-	5,5,5	1.09	0	5,5,5	0.34	0
3	ACY	A	1003	-	1,3,3	1.99	0	0,3,3	0.00	-
3	ACY	B	1001	-	1,3,3	3.33	1 (100%)	0,3,3	0.00	-
2	ILP	A	400	-	21,24,24	2.02	6 (28%)	26,34,34	1.71	8 (30%)
3	ACY	B	1006	-	1,3,3	3.24	1 (100%)	0,3,3	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
3	ACY	A	1005	-	1,3,3	4.28	1 (100%)	0,3,3	0.00	-
2	ILP	B	900	-	21,24,24	3.18	6 (28%)	26,34,34	2.28	8 (30%)
3	ACY	A	1004	-	1,3,3	3.74	1 (100%)	0,3,3	0.00	-

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	B	2001	-	-	0/4/4/4	-
2	ILP	B	900	-	-	9/17/21/21	0/1/1/1
2	ILP	A	400	-	1/1/4/5	5/17/21/21	0/1/1/1

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	900	ILP	C4A-C4	11.22	1.66	1.51
2	A	400	ILP	C4A-C4	5.15	1.58	1.51
2	A	400	ILP	CA-N	-4.83	1.38	1.47
2	B	900	ILP	CA-N	-4.62	1.39	1.47
2	B	900	ILP	C3-C4	4.40	1.46	1.40
3	A	1005	ACY	CH3-C	4.28	1.54	1.48
2	B	900	ILP	C5-C4	4.16	1.46	1.40
3	B	1002	ACY	CH3-C	4.01	1.53	1.48
3	A	1004	ACY	CH3-C	3.74	1.53	1.48
3	B	1001	ACY	CH3-C	3.33	1.53	1.48
3	B	1006	ACY	CH3-C	3.24	1.52	1.48
2	B	900	ILP	C3-C2	2.94	1.43	1.40
2	A	400	ILP	C3-C4	2.76	1.44	1.40
2	A	400	ILP	C4A-N	-2.52	1.39	1.46
2	B	900	ILP	C6-N1	2.42	1.39	1.34
2	A	400	ILP	C5-C4	2.09	1.43	1.40
2	A	400	ILP	P-O3P	-2.05	1.47	1.54

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	900	ILP	C4A-N-CA	5.34	121.18	113.78
2	B	900	ILP	C3-C4-C5	-4.60	114.31	118.72
2	B	900	ILP	C4A-C4-C3	4.51	124.87	120.04

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	400	ILP	C4A-N-CA	4.13	119.50	113.78
2	B	900	ILP	C6-C5-C4	3.93	120.90	118.12
2	B	900	ILP	CG2-CB-CG1	-2.94	104.34	111.78
2	A	400	ILP	C6-C5-C4	2.67	120.01	118.12
2	A	400	ILP	O3P-P-O1P	2.46	120.33	110.68
2	B	900	ILP	O2P-P-O4P	-2.35	100.49	106.73
2	A	400	ILP	C4-C4A-N	2.32	118.18	111.78
2	A	400	ILP	C6-N1-C2	2.32	123.46	119.17
2	B	900	ILP	O4P-P-O1P	2.23	112.74	106.47
2	A	400	ILP	C5-C6-N1	-2.16	120.21	123.82
2	A	400	ILP	O4P-P-O1P	2.03	112.18	106.47
2	B	900	ILP	C6-N1-C2	2.02	122.91	119.17
2	A	400	ILP	O2P-P-O4P	-2.00	101.41	106.73

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	400	ILP	CA

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	400	ILP	CB-CA-N-C4A
2	A	400	ILP	C-CA-N-C4A
2	B	900	ILP	C-CA-CB-CG1
2	B	900	ILP	C5-C4-C4A-N
2	A	400	ILP	N-CA-CB-CG2
2	B	900	ILP	N-CA-CB-CG2
2	B	900	ILP	CG2-CB-CG1-CD1
2	A	400	ILP	N-CA-CB-CG1
2	B	900	ILP	C-CA-CB-CG2
2	B	900	ILP	C4-C4A-N-CA
2	A	400	ILP	CG2-CB-CG1-CD1
2	B	900	ILP	C3-C4-C4A-N
2	B	900	ILP	CA-CB-CG1-CD1
2	B	900	ILP	N-CA-CB-CG1

There are no ring outliers.

5 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	2001	GOL	1	0

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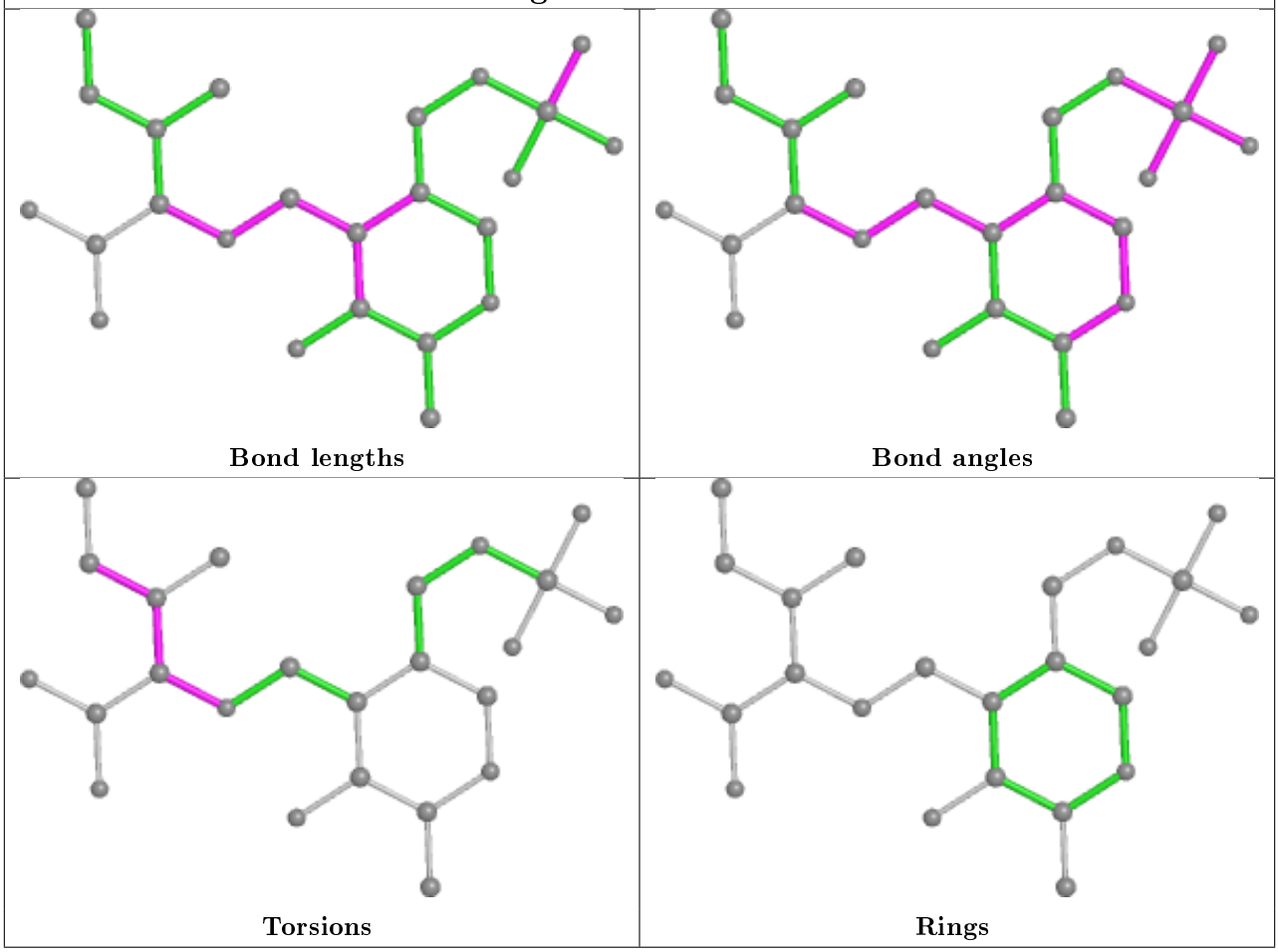
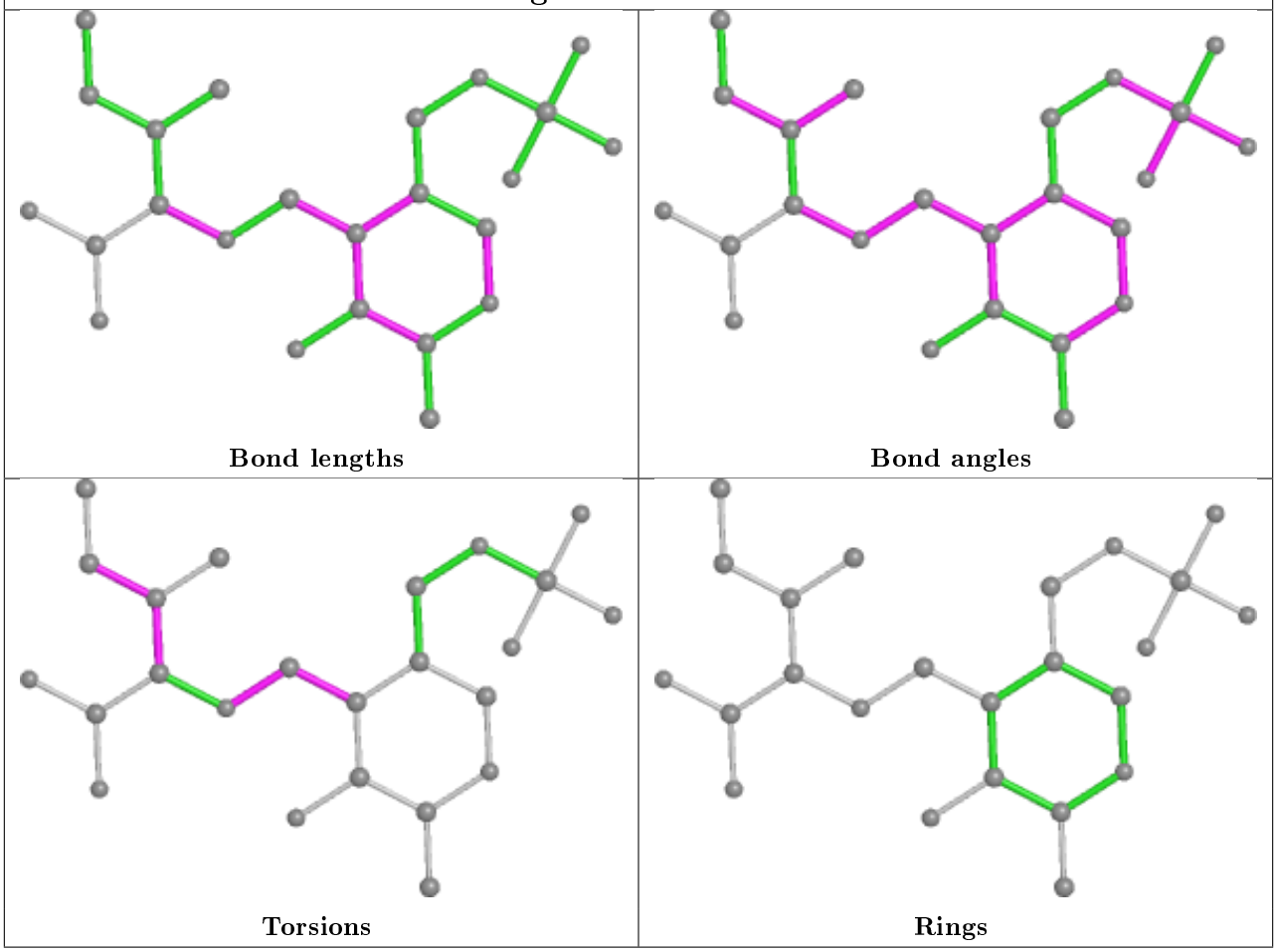
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1003	ACY	1	0
3	B	1001	ACY	1	0
2	A	400	ILP	9	0
2	B	900	ILP	6	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand ILP A 400



Ligand ILP B 900



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, 95^{th} 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	360/365 (98%)	0.67	36 (10%)	7	8	21, 35, 52, 65	2 (0%)
1	B	365/365 (100%)	0.77	40 (10%)	5	6	21, 37, 57, 69	0
All	All	725/730 (99%)	0.72	76 (10%)	6	7	21, 36, 55, 69	2 (0%)

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	264[A]	VAL	11.1
1	B	501	ALA	9.7
1	B	525	GLY	9.4
1	B	552[A]	ARG	8.6
1	B	502	SER	8.5
1	A	24	PRO	7.3
1	A	179	VAL	7.3
1	A	2	SER	7.3
1	B	503	SER	6.8
1	A	1	ALA	6.3
1	B	526	GLU	6.3
1	B	528	LEU	5.8
1	A	25	GLY	5.7
1	A	26	GLU	5.6
1	B	678	SER	5.4
1	B	527	PRO	5.2
1	B	828[A]	ARG	5.2
1	B	524	PRO	5.2
1	B	676	GLY	5.0
1	A	3	SER	5.0
1	A	357	ARG	4.7
1	A	173	TYR	4.7
1	B	523	GLY	4.7
1	A	172	ALA	4.6

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Mol	Chain	Res	Type	RSRZ
1	A	358	ALA	4.3
1	A	23	GLY	4.3
1	B	675	PRO	4.3
1	A	328[A]	ARG	4.0
1	A	28	LEU	4.0
1	A	159	ARG	4.0
1	A	84	LYS	3.8
1	B	583	GLY	3.8
1	B	585	ASP	3.5
1	B	674	PHE	3.5
1	B	673	TYR	3.5
1	A	252	GLY	3.4
1	A	46	LYS	3.4
1	B	520	LYS	3.3
1	A	45	ASP	3.2
1	B	518	PRO	3.2
1	A	27	PRO	3.0
1	A	171	GLY	2.9
1	B	517	LYS	2.9
1	B	546	LYS	2.9
1	A	231	PRO	2.8
1	B	586	GLN	2.8
1	B	519	HIS	2.8
1	A	356	ILE	2.8
1	A	232[A]	ASP	2.8
1	B	522	PRO	2.7
1	B	504	SER	2.7
1	B	516	GLN	2.6
1	B	731	PRO	2.6
1	B	857	ARG	2.6
1	B	827	ASP	2.6
1	B	515	THR	2.5
1	A	245	VAL	2.5
1	B	838	GLY	2.5
1	A	167	LEU	2.4
1	A	22	PRO	2.4
1	B	677	GLY	2.4
1	A	140	LEU	2.3
1	A	246	TYR	2.3
1	B	512	LEU	2.3
1	A	327	ASP	2.3
1	A	83	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	808	VAL	2.2
1	B	536	ASP	2.2
1	B	663	LEU	2.2
1	B	655	VAL	2.1
1	A	91	PHE	2.1
1	A	355	GLY	2.1
1	A	85	ASP	2.1
1	B	826	LYS	2.1
1	B	659	ARG	2.0
1	A	180	THR	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)

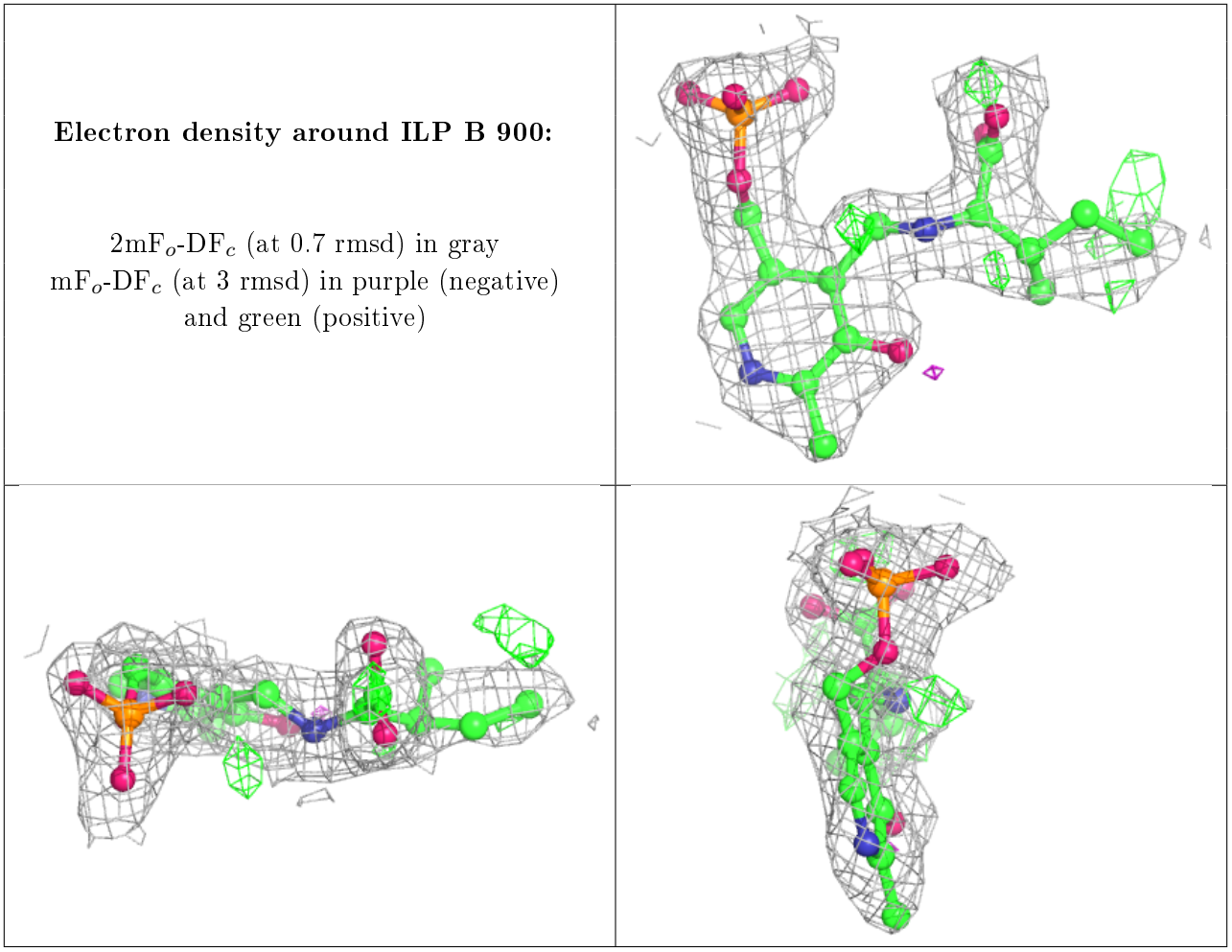
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	ACY	A	1003	4/4	0.78	0.19	40,40,41,43	0
3	ACY	B	1002	4/4	0.79	0.13	50,50,51,51	0
3	ACY	B	1006	4/4	0.82	0.17	44,45,46,47	0
4	GOL	B	2001	6/6	0.84	0.22	54,55,55,56	0
3	ACY	A	1005	4/4	0.84	0.14	42,43,44,46	0
3	ACY	A	1004	4/4	0.85	0.13	54,55,55,55	0
3	ACY	B	1001	4/4	0.90	0.11	49,49,49,50	0
2	ILP	B	900	24/24	0.91	0.15	32,43,46,48	8
2	ILP	A	400	24/24	0.93	0.14	24,28,38,40	8

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

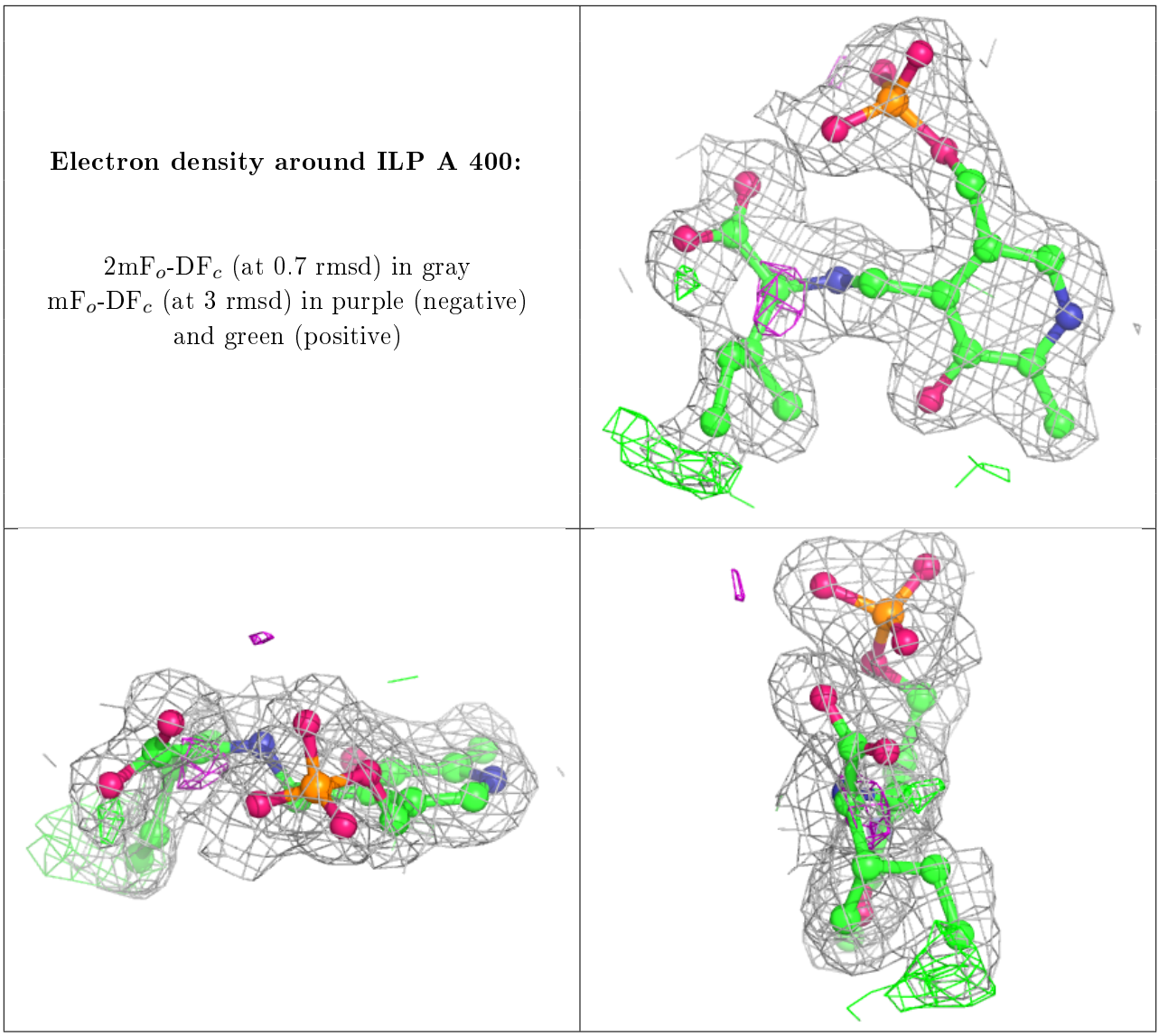
Electron density around ILP B 900:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



Electron density around ILP A 400:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



6.5 Other polymers

There are no such residues in this entry.