



# Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 02:54 pm BST

PDB ID : 2H23  
Title : Structure of Rubisco LSMT bound to Trimethyllysine and AdoHcy  
Authors : Couture, J.F.; Hauk, G.; Trievel, R.C.  
Deposited on : 2006-05-17  
Resolution : 2.45 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (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

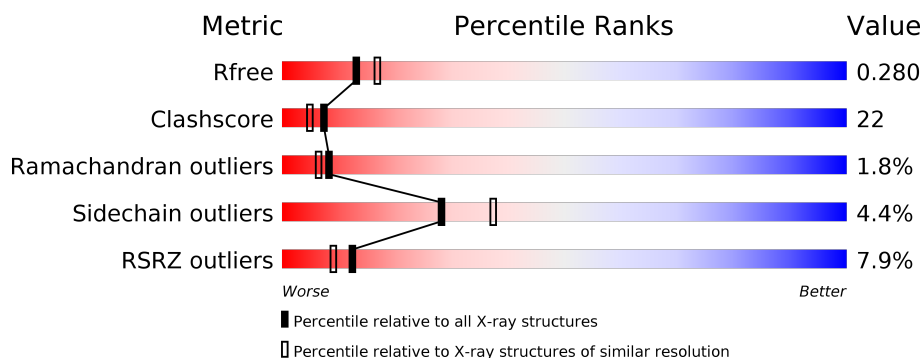
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.45 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	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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 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  $\leq 5\%$ . 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	440	<div> <div>14%</div> <div> <div></div> <div>57%</div> <div>36%</div> <div></div> </div> <div></div> </div>
1	B	440	<div> <div>5%</div> <div> <div></div> <div>65%</div> <div>31%</div> <div></div> </div> <div></div> </div>
1	C	440	<div> <div>4%</div> <div> <div></div> <div>63%</div> <div>33%</div> <div></div> </div> <div></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	M3L	A	501	-	-	-	X
3	M3L	C	500	-	-	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10988 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.

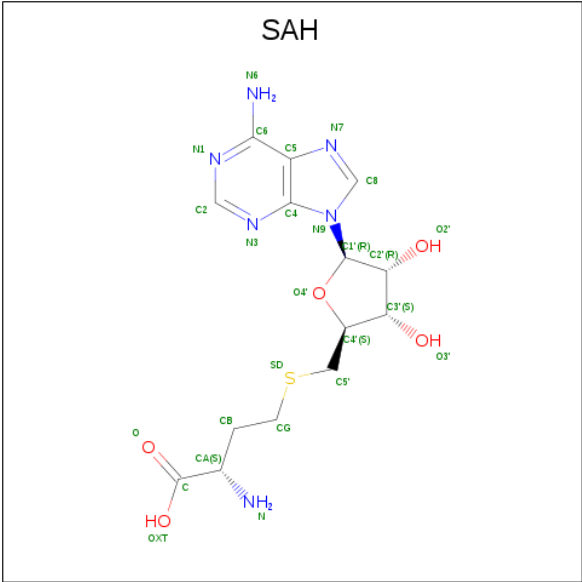
- Molecule 1 is a protein called Ribulose-1,5 bisphosphate carboxylase/oxygenase large subunit N-methyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	424	Total	C	N	O	S	0	0	0
			3413	2189	562	655	7			
1	B	440	Total	C	N	O	S	0	0	0
			3542	2270	585	680	7			
1	C	438	Total	C	N	O	S	0	0	0
			3526	2262	582	675	7			

There are 18 discrepancies between the modelled and reference sequences:

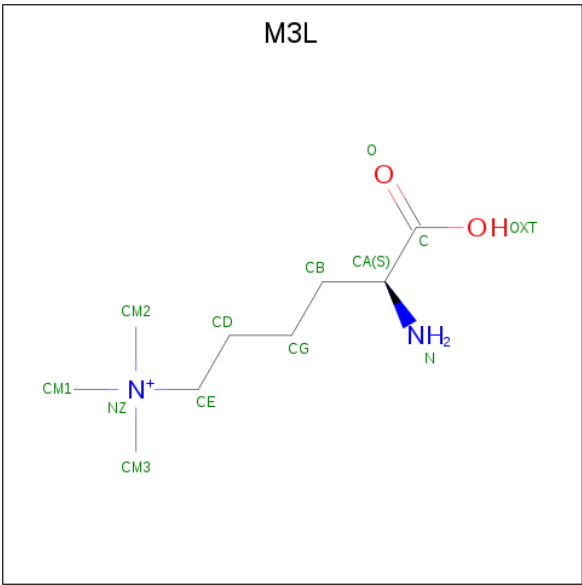
Chain	Residue	Modelled	Actual	Comment	Reference
A	483	GLU	-	CLONING ARTIFACT	UNP Q43088
A	484	ASN	-	CLONING ARTIFACT	UNP Q43088
A	485	LEU	-	CLONING ARTIFACT	UNP Q43088
A	486	TYR	-	CLONING ARTIFACT	UNP Q43088
A	487	PHE	-	CLONING ARTIFACT	UNP Q43088
A	488	GLN	-	CLONING ARTIFACT	UNP Q43088
B	483	GLU	-	CLONING ARTIFACT	UNP Q43088
B	484	ASN	-	CLONING ARTIFACT	UNP Q43088
B	485	LEU	-	CLONING ARTIFACT	UNP Q43088
B	486	TYR	-	CLONING ARTIFACT	UNP Q43088
B	487	PHE	-	CLONING ARTIFACT	UNP Q43088
B	488	GLN	-	CLONING ARTIFACT	UNP Q43088
C	483	GLU	-	CLONING ARTIFACT	UNP Q43088
C	484	ASN	-	CLONING ARTIFACT	UNP Q43088
C	485	LEU	-	CLONING ARTIFACT	UNP Q43088
C	486	TYR	-	CLONING ARTIFACT	UNP Q43088
C	487	PHE	-	CLONING ARTIFACT	UNP Q43088
C	488	GLN	-	CLONING ARTIFACT	UNP Q43088

- Molecule 2 is S-ADENOSYL-L-HOMOCYSTEINE (three-letter code: SAH) (formula: C<sub>14</sub>H<sub>20</sub>N<sub>6</sub>O<sub>5</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
2	B	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
2	C	1	Total	C	N	O	S	0	0
			26	14	6	5	1		

- Molecule 3 is N-TRIMETHYLLYSINE (three-letter code: M3L) (formula: C<sub>9</sub>H<sub>21</sub>N<sub>2</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			13	9	2	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			13	9	2	2		
3	C	1	Total	C	N	O	0	0
			13	9	2	2		

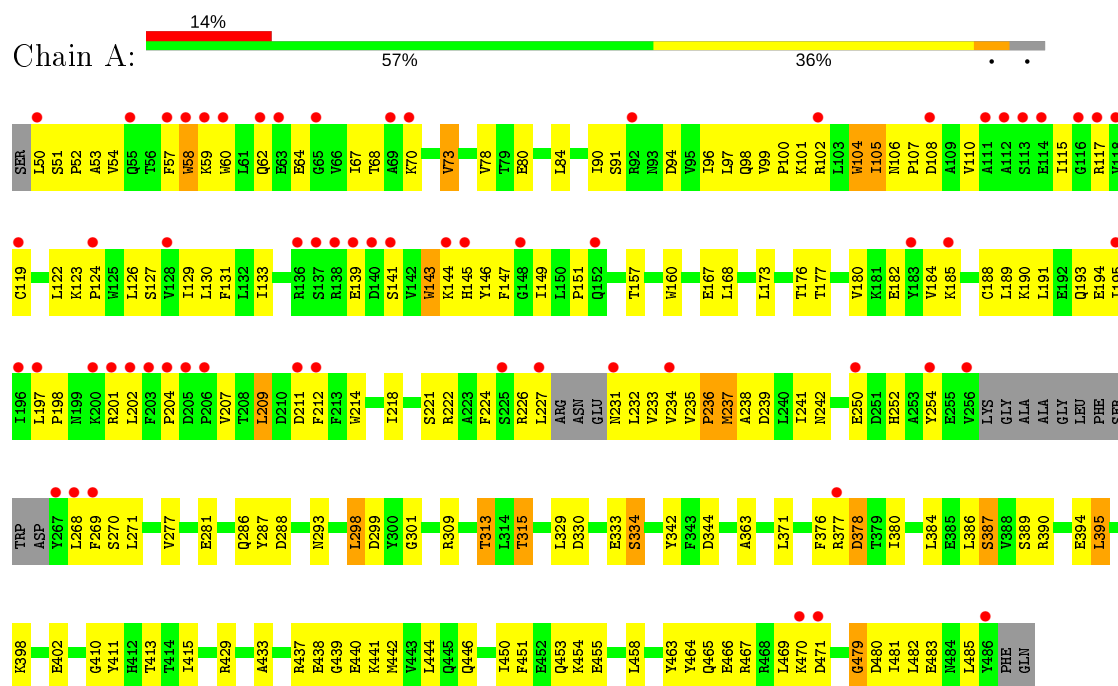
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	111	Total	O	0	0
			111	111		
4	B	136	Total	O	0	0
			136	136		
4	C	143	Total	O	0	0
			143	143		

### 3 Residue-property plots

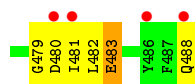
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

- Molecule 1: Ribulose-1,5 biphosphate carboxylase/oxygenase large subunit N-methyltransferase

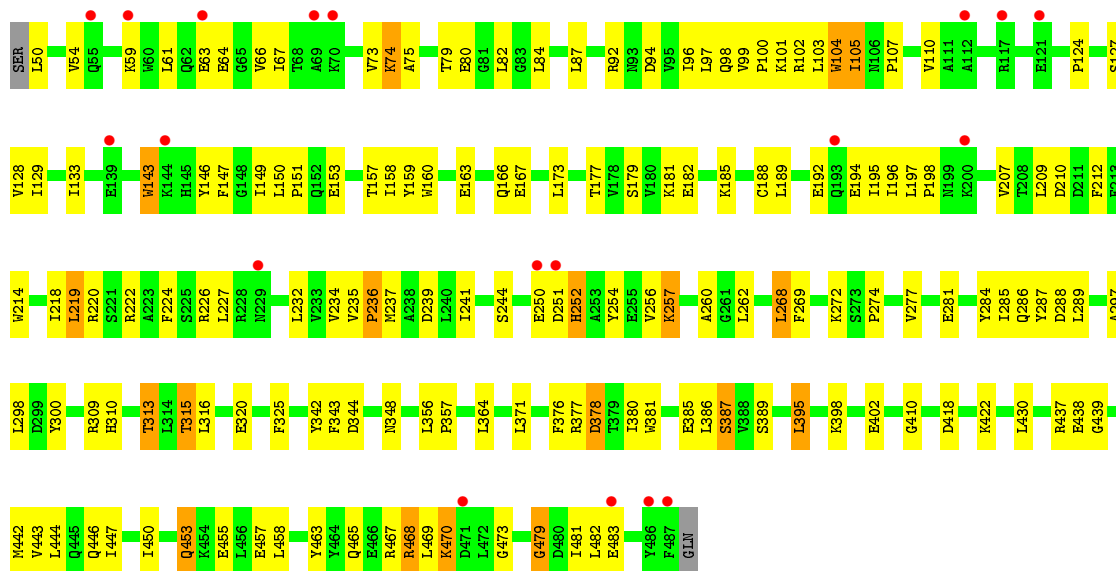


- Molecule 1: Ribulose-1,5 biphosphate carboxylase/oxygenase large subunit N-methyltransferase





- Molecule 1: Ribulose-1,5 biphosphate carboxylase/oxygenase large subunit N-methyltransferase





## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	131.86Å 157.55Å 267.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.98 – 2.45 14.98 – 2.44	Depositor EDS
% Data completeness (in resolution range)	96.5 (14.98-2.45) 96.6 (14.98-2.44)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.52 (at 2.45Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.253 , 0.288 0.245 , 0.280	Depositor DCC
$R_{free}$ test set	10008 reflections (9.74%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	55.0	Xtriage
Anisotropy	0.454	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 50.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	10988	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.03% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SAH, M3L

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	0.41	0/3482	0.58	0/4723
1	B	0.41	0/3617	0.62	1/4906 (0.0%)
1	C	0.39	0/3601	0.60	0/4886
All	All	0.40	0/10700	0.60	1/14515 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	236	PRO	N-CA-C	5.03	125.17	112.10

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	3413	0	3384	194	0
1	B	3542	0	3501	138	0
1	C	3526	0	3488	157	0
2	A	26	0	19	0	0
2	B	26	0	19	0	0
2	C	26	0	19	3	0
3	A	13	0	20	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	13	0	20	5	0
3	C	13	0	20	8	0
4	A	111	0	0	21	0
4	B	136	0	0	9	0
4	C	143	0	0	11	0
All	All	10988	0	10490	471	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 22.

All (471) 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)
3:C:500:M3L:CG	3:C:500:M3L:CD	1.79	1.59
1:B:395:LEU:HD22	1:C:469:LEU:HD12	1.44	0.99
1:A:482:LEU:HD11	1:C:179:SER:HB3	1.43	0.98
1:A:395:LEU:HD22	1:B:469:LEU:HD12	1.45	0.98
1:C:177:THR:HB	4:C:820:HOH:O	1.64	0.96
1:B:438:GLU:HG2	1:B:442:MET:HE2	1.46	0.95
1:B:286:GLN:HE21	1:B:288:ASP:H	1.14	0.95
1:C:470:LYS:HA	1:C:470:LYS:HE2	1.50	0.94
1:C:50:LEU:HD22	1:C:54:VAL:HG11	1.51	0.91
1:A:363:ALA:HB1	1:A:395:LEU:HD13	1.55	0.86
3:C:500:M3L:CB	3:C:500:M3L:CD	2.52	0.86
1:A:254:TYR:HB3	1:A:270:SER:HB3	1.58	0.85
1:A:439:GLY:HA2	1:A:442:MET:HE3	1.57	0.85
1:A:469:LEU:HD12	1:C:395:LEU:HD22	1.59	0.83
1:A:194:GLU:C	1:A:195:ILE:HD12	1.99	0.83
1:A:438:GLU:HG2	1:A:442:MET:HE2	1.63	0.81
1:C:226:ARG:HG2	1:C:252:HIS:CE1	2.15	0.81
1:B:179:SER:HB3	1:C:482:LEU:HD11	1.62	0.80
1:B:439:GLY:HA2	1:B:442:MET:HE3	1.63	0.79
1:B:97:LEU:HD13	1:B:237:MET:HE3	1.64	0.78
1:C:73:VAL:HG21	1:C:84:LEU:HB3	1.65	0.78
1:A:130:LEU:HD11	1:A:191:LEU:HD22	1.64	0.78
1:B:286:GLN:HE21	1:B:288:ASP:N	1.82	0.77
1:A:106:ASN:HD21	1:A:108:ASP:HB2	1.49	0.77
1:B:286:GLN:NE2	1:B:288:ASP:H	1.83	0.77
1:A:102:ARG:HA	4:A:904:HOH:O	1.84	0.77
1:A:126:LEU:HD11	4:A:832:HOH:O	1.86	0.76
1:C:97:LEU:HB2	1:C:237:MET:HE1	1.66	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:96:ILE:HD11	1:B:273:SER:HB2	1.69	0.75
1:B:173:LEU:O	1:B:177:THR:HG23	1.88	0.74
1:B:78:VAL:HG11	1:B:282:GLN:HE22	1.53	0.73
1:C:348:ASN:H	1:C:446:GLN:HE22	1.37	0.73
1:A:100:PRO:HA	1:A:268:LEU:HD12	1.71	0.73
1:C:185:LYS:HD3	4:C:852:HOH:O	1.86	0.73
1:C:371:LEU:HB3	1:C:380:ILE:HD13	1.68	0.73
1:B:315:THR:HG22	4:C:808:HOH:O	1.87	0.73
1:C:104:TRP:HH2	1:C:269:PHE:H	1.37	0.73
1:A:395:LEU:CD2	1:B:469:LEU:HD12	2.18	0.72
1:A:139:GLU:HG3	4:A:836:HOH:O	1.89	0.72
1:A:214:TRP:O	1:A:218:ILE:HG12	1.89	0.72
1:B:371:LEU:HB3	1:B:380:ILE:HD13	1.70	0.72
1:A:115:ILE:HG22	1:A:202:LEU:HD13	1.71	0.72
1:C:177:THR:HG22	1:C:298:LEU:HD13	1.71	0.71
1:A:479:GLY:HA3	1:C:342:TYR:CE2	2.25	0.71
1:A:119:CYS:HA	1:A:122:LEU:HD12	1.73	0.71
1:B:99:VAL:O	1:B:104:TRP:CH2	2.45	0.70
1:A:207:VAL:HB	4:A:806:HOH:O	1.92	0.70
1:A:286:GLN:HE21	1:A:288:ASP:C	1.95	0.69
1:A:147:PHE:HA	4:A:845:HOH:O	1.91	0.69
1:A:99:VAL:O	1:A:104:TRP:HH2	1.74	0.69
1:C:73:VAL:CG2	1:C:84:LEU:HB3	2.23	0.68
1:B:99:VAL:O	1:B:104:TRP:HH2	1.76	0.68
1:A:226:ARG:HH12	1:A:250:GLU:HG2	1.59	0.68
1:C:194:GLU:C	1:C:195:ILE:HD12	2.13	0.68
1:C:74:LYS:HG2	1:C:87:LEU:HD21	1.75	0.68
1:A:173:LEU:O	1:A:177:THR:HG23	1.93	0.68
1:B:107:PRO:O	1:B:110:VAL:HG22	1.93	0.68
1:A:129:ILE:O	1:A:133:ILE:HG13	1.93	0.67
1:A:101:LYS:HA	1:A:104:TRP:CD2	2.29	0.67
1:A:226:ARG:C	1:A:227:LEU:HD22	2.15	0.67
1:B:395:LEU:CD2	1:C:469:LEU:HD12	2.24	0.67
1:A:465:GLN:HB3	1:C:395:LEU:HD11	1.77	0.67
1:B:287:TYR:CZ	3:B:502:M3L:HM13	2.30	0.66
1:B:142:VAL:HG23	4:B:819:HOH:O	1.94	0.66
1:B:467:ARG:HA	1:B:470:LYS:HG2	1.77	0.66
1:A:188:CYS:HB3	1:A:212:PHE:CD1	2.31	0.65
1:B:197:LEU:HB2	1:B:198:PRO:HD3	1.78	0.65
1:B:67:ILE:HD11	1:B:237:MET:SD	2.37	0.65
1:B:73:VAL:HG21	1:B:84:LEU:HB3	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:TRP:HB2	1:A:147:PHE:CE1	2.32	0.65
1:A:198:PRO:HG3	4:A:805:HOH:O	1.97	0.65
1:A:98:GLN:HG2	1:A:270:SER:HA	1.78	0.64
1:C:438:GLU:CD	1:C:442:MET:HE2	2.18	0.64
1:A:330:ASP:O	1:A:334:SER:HB2	1.96	0.64
1:A:53:ALA:C	1:A:149:ILE:HD11	2.18	0.64
1:B:401:ARG:O	1:B:405:LYS:HG3	1.97	0.64
1:C:222:ARG:NH2	2:C:802:SAH:O	2.29	0.64
1:A:101:LYS:HA	1:A:104:TRP:CE3	2.33	0.64
1:A:167:GLU:HG3	1:A:437:ARG:NH1	2.12	0.64
1:C:468:ARG:HH11	1:C:468:ARG:HB3	1.63	0.63
1:A:313:THR:HB	1:A:344:ASP:OD1	1.98	0.63
1:B:128:VAL:HG11	1:B:219:LEU:HD11	1.81	0.63
1:C:250:GLU:OE2	1:C:288:ASP:HA	1.99	0.63
1:C:300:TYR:OH	3:C:500:M3L:HB3	1.99	0.63
1:A:98:GLN:HE21	1:A:270:SER:HB2	1.64	0.62
1:B:222:ARG:HA	3:B:502:M3L:HM23	1.80	0.62
1:C:214:TRP:O	1:C:218:ILE:HG12	1.99	0.62
1:A:160:TRP:CD1	1:A:429:ARG:HD3	2.35	0.62
1:A:105:ILE:O	1:A:105:ILE:HG22	1.99	0.62
1:B:101:LYS:HA	1:B:104:TRP:CD2	2.33	0.62
1:C:67:ILE:HD11	1:C:237:MET:SD	2.39	0.62
1:A:222:ARG:HD3	1:A:239:ASP:OD2	2.00	0.62
1:A:481:ILE:HG23	1:C:220:ARG:HH22	1.65	0.62
1:A:99:VAL:O	1:A:104:TRP:CH2	2.52	0.62
1:A:238:ALA:HB1	1:A:271:LEU:HD22	1.81	0.62
1:A:222:ARG:HA	3:A:501:M3L:HM23	1.81	0.61
1:C:107:PRO:O	1:C:110:VAL:HG22	2.00	0.61
1:A:78:VAL:HG23	1:A:80:GLU:HG2	1.81	0.61
1:B:438:GLU:HG2	1:B:442:MET:CE	2.25	0.61
1:B:78:VAL:HG12	1:B:80:GLU:H	1.65	0.61
1:B:427:ASP:HA	4:B:893:HOH:O	1.99	0.61
1:C:227:LEU:HD11	1:C:256:VAL:HG23	1.82	0.61
1:C:222:ARG:O	3:C:500:M3L:HM13	2.01	0.61
1:C:286:GLN:HE21	1:C:288:ASP:C	2.03	0.61
1:C:479:GLY:O	1:C:483:GLU:HG2	2.00	0.61
1:B:176:THR:HG21	1:C:481:ILE:HD12	1.82	0.60
1:A:99:VAL:HG21	1:A:237:MET:HB3	1.82	0.60
1:C:157:THR:HA	1:C:160:TRP:CD1	2.36	0.60
1:C:97:LEU:HD22	1:C:237:MET:HE3	1.84	0.60
1:A:130:LEU:HD21	1:A:191:LEU:HD13	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:463:TYR:O	1:A:467:ARG:HG3	2.02	0.60
1:A:437:ARG:O	1:A:441:LYS:HG3	2.02	0.60
1:A:58:TRP:CE3	1:A:58:TRP:HA	2.36	0.59
1:A:293:ASN:ND2	1:A:309:ARG:HB2	2.18	0.59
1:C:188:CYS:SG	1:C:212:PHE:CD1	2.95	0.59
1:B:381:TRP:O	1:B:385:GLU:HG3	2.02	0.59
1:A:218:ILE:HG23	1:A:222:ARG:HD2	1.83	0.59
1:C:453:GLN:O	1:C:457:GLU:HG3	2.02	0.59
1:A:105:ILE:HD11	1:A:143:TRP:CD2	2.37	0.59
1:A:413:THR:O	1:A:441:LYS:HE2	2.03	0.58
1:A:106:ASN:ND2	1:A:108:ASP:HB2	2.17	0.58
1:A:78:VAL:CG2	1:A:80:GLU:HG2	2.33	0.58
1:A:313:THR:HG22	4:A:909:HOH:O	2.04	0.58
1:A:107:PRO:O	1:A:110:VAL:HG22	2.03	0.58
1:A:466:GLU:O	1:A:470:LYS:HG2	2.04	0.58
1:A:189:LEU:O	1:A:193:GLN:HG2	2.03	0.58
1:B:49:SER:O	1:B:52:PRO:HG2	2.03	0.58
1:B:59:LYS:O	1:B:63:GLU:HG3	2.04	0.58
1:B:61:LEU:HB3	1:B:67:ILE:HG12	1.84	0.58
1:C:97:LEU:HD22	1:C:237:MET:CE	2.34	0.58
1:A:99:VAL:CG2	1:A:237:MET:HB3	2.33	0.58
1:C:226:ARG:HD2	1:C:252:HIS:NE2	2.19	0.58
1:C:92:ARG:HG3	1:C:274:PRO:O	2.03	0.57
1:C:103:LEU:O	1:C:143:TRP:CZ3	2.57	0.57
1:A:64:GLU:HA	1:A:102:ARG:HH12	1.69	0.57
1:A:54:VAL:N	1:A:149:ILE:HD11	2.19	0.57
1:B:188:CYS:HB3	1:B:212:PHE:CD1	2.39	0.57
1:A:110:VAL:HG12	1:A:131:PHE:CG	2.39	0.57
1:C:398:LYS:O	1:C:402:GLU:HG2	2.05	0.57
1:A:479:GLY:HA3	1:C:342:TYR:CZ	2.40	0.57
1:C:157:THR:OG1	1:C:177:THR:HG21	2.04	0.56
1:B:227:LEU:HD22	1:B:255:GLU:CD	2.26	0.56
1:C:104:TRP:CH2	1:C:269:PHE:HB2	2.40	0.56
1:C:286:GLN:HE22	1:C:309:ARG:HH22	1.53	0.56
1:A:101:LYS:HG3	1:A:104:TRP:CD1	2.41	0.56
1:A:221:SER:O	3:A:501:M3L:HM12	2.06	0.56
1:C:222:ARG:HD3	1:C:239:ASP:OD2	2.06	0.56
1:A:455:GLU:O	1:A:458:LEU:HB2	2.05	0.56
1:C:104:TRP:HH2	1:C:269:PHE:N	2.02	0.56
1:A:105:ILE:HD11	1:A:143:TRP:CE2	2.40	0.56
1:B:78:VAL:HG12	1:B:80:GLU:HG2	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:470:LYS:HA	1:C:470:LYS:CE	2.31	0.56
1:A:167:GLU:HG3	1:A:437:ARG:HH12	1.71	0.56
1:A:376:PHE:O	1:A:378:ASP:N	2.39	0.56
1:C:463:TYR:O	1:C:467:ARG:HG3	2.06	0.56
1:C:73:VAL:HG11	1:C:96:ILE:HG22	1.88	0.55
1:C:222:ARG:HH11	1:C:239:ASP:CG	2.10	0.55
1:A:286:GLN:HE22	1:A:309:ARG:HH22	1.54	0.55
1:C:74:LYS:HD3	1:C:75:ALA:O	2.06	0.55
1:A:226:ARG:O	1:A:227:LEU:HD13	2.07	0.55
1:B:110:VAL:HG21	1:B:127:SER:HB3	1.87	0.55
1:B:78:VAL:CG1	1:B:79:THR:N	2.69	0.55
1:A:51:SER:HB3	1:A:54:VAL:HB	1.88	0.54
1:B:194:GLU:C	1:B:195:ILE:HD12	2.28	0.54
1:C:105:ILE:HB	1:C:234:VAL:HB	1.90	0.54
1:B:103:LEU:O	1:B:143:TRP:HZ3	1.91	0.54
1:A:315:THR:HG23	1:B:473:GLY:O	2.06	0.54
1:A:277:VAL:HG13	1:A:281:GLU:HB2	1.90	0.54
1:A:376:PHE:C	1:A:378:ASP:H	2.10	0.54
1:B:73:VAL:CG2	1:B:84:LEU:HB3	2.37	0.54
1:B:73:VAL:HG23	1:B:85:VAL:O	2.08	0.54
1:A:97:LEU:HD12	1:A:97:LEU:O	2.09	0.53
1:C:173:LEU:O	1:C:177:THR:HG23	2.08	0.53
1:B:241:ILE:CD1	1:B:285:ILE:HG23	2.39	0.53
1:B:80:GLU:HG3	1:B:242:ASN:ND2	2.24	0.53
1:B:51:SER:N	1:B:52:PRO:HD2	2.23	0.53
1:C:250:GLU:HB3	1:C:286:GLN:CG	2.38	0.53
1:C:316:LEU:HA	4:C:942:HOH:O	2.07	0.53
1:C:455:GLU:O	1:C:458:LEU:HB2	2.09	0.53
1:C:54:VAL:HA	1:C:149:ILE:HD11	1.91	0.53
1:A:398:LYS:O	1:A:402:GLU:HG2	2.09	0.52
1:A:58:TRP:NE1	4:A:852:HOH:O	2.32	0.52
1:C:386:LEU:O	1:C:387:SER:CB	2.57	0.52
1:A:315:THR:HG22	4:B:807:HOH:O	2.09	0.52
1:A:104:TRP:CZ2	1:A:269:PHE:HB2	2.45	0.52
1:A:91:SER:O	1:A:94:ASP:HB2	2.10	0.52
1:A:182:GLU:O	1:A:185:LYS:HB3	2.09	0.52
1:B:78:VAL:HG12	1:B:79:THR:N	2.24	0.52
1:A:141:SER:HB3	4:A:808:HOH:O	2.09	0.52
1:A:123:LYS:CB	1:A:124:PRO:HD2	2.40	0.52
1:B:177:THR:HG22	1:B:298:LEU:HD13	1.92	0.52
1:B:222:ARG:HH11	1:B:239:ASP:CG	2.13	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:LEU:O	1:B:143:TRP:CZ3	2.63	0.52
1:B:227:LEU:HB2	1:B:230:GLU:HB2	1.91	0.52
1:A:180:VAL:O	1:A:184:VAL:HG23	2.10	0.51
1:A:123:LYS:HB3	1:A:124:PRO:HD2	1.91	0.51
1:A:197:LEU:HB2	1:A:198:PRO:HD3	1.91	0.51
1:A:157:THR:OG1	1:A:177:THR:HG21	2.11	0.51
1:A:54:VAL:HG22	1:A:149:ILE:HD12	1.92	0.51
1:B:174:LEU:O	1:B:178:VAL:HG23	2.10	0.51
1:B:428:SER:O	1:B:432:ILE:HG13	2.09	0.51
1:A:50:LEU:HD22	1:A:54:VAL:HG11	1.93	0.51
1:B:479:GLY:O	1:B:483:GLU:HB2	2.10	0.51
1:C:438:GLU:HG2	1:C:442:MET:HE3	1.92	0.51
1:A:119:CYS:CA	1:A:122:LEU:HD12	2.41	0.51
1:A:157:THR:OG1	1:A:177:THR:CG2	2.59	0.51
1:A:211:ASP:O	1:A:214:TRP:HB3	2.11	0.51
4:A:809:HOH:O	1:C:315:THR:HG22	2.11	0.51
1:B:78:VAL:CG1	1:B:80:GLU:HG2	2.42	0.50
1:C:103:LEU:O	1:C:143:TRP:HZ3	1.94	0.50
1:C:80:GLU:N	1:C:80:GLU:OE2	2.39	0.50
1:A:410:GLY:HA3	4:A:824:HOH:O	2.11	0.50
1:B:106:ASN:HD21	1:B:108:ASP:HB2	1.76	0.50
1:C:224:PHE:HD2	1:C:254:TYR:CE2	2.28	0.50
1:A:277:VAL:CG1	1:A:281:GLU:HB2	2.41	0.50
1:B:101:LYS:HA	1:B:104:TRP:CE2	2.47	0.50
1:B:386:LEU:O	1:B:387:SER:CB	2.59	0.50
1:B:92:ARG:O	1:B:93:ASN:HB2	2.10	0.50
1:A:371:LEU:HB3	1:A:380:ILE:HD13	1.92	0.50
1:A:68:THR:C	1:A:70:LYS:H	2.14	0.50
1:A:104:TRP:CH2	1:A:269:PHE:HB2	2.46	0.50
1:C:226:ARG:HG2	1:C:252:HIS:NE2	2.25	0.50
1:C:438:GLU:HG2	1:C:442:MET:CE	2.42	0.50
3:C:500:M3L:CE	3:C:500:M3L:CG	2.80	0.50
1:A:301:GLY:HA3	4:A:804:HOH:O	2.11	0.50
1:B:222:ARG:CA	3:B:502:M3L:HM23	2.42	0.50
1:A:151:PRO:HD2	1:A:222:ARG:HH21	1.77	0.49
1:A:286:GLN:NE2	1:A:288:ASP:H	2.10	0.49
1:C:250:GLU:C	1:C:252:HIS:H	2.16	0.49
1:A:60:TRP:HB2	1:A:145:HIS:CD2	2.47	0.49
1:C:66:VAL:HA	4:C:849:HOH:O	2.13	0.49
1:A:100:PRO:HA	1:A:268:LEU:CD1	2.41	0.49
1:B:128:VAL:CG1	1:B:219:LEU:HD11	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:316:LEU:HD12	1:C:343:PHE:CE1	2.47	0.49
1:A:380:ILE:O	1:A:384:LEU:HG	2.13	0.49
1:B:226:ARG:HE	1:B:226:ARG:HA	1.77	0.49
1:A:390:ARG:HA	1:A:463:TYR:CZ	2.48	0.49
1:A:59:LYS:HA	1:A:62:GLN:OE1	2.13	0.49
1:B:446:GLN:O	1:B:450:ILE:HG13	2.13	0.49
1:C:446:GLN:O	1:C:450:ILE:HG13	2.13	0.49
1:A:224:PHE:CZ	3:A:501:M3L:HM22	2.48	0.49
1:A:57:PHE:CE1	1:A:146:TYR:HA	2.48	0.49
1:B:453:GLN:O	1:B:457:GLU:HG3	2.13	0.49
1:A:96:ILE:HG13	1:A:97:LEU:HG	1.94	0.49
1:C:61:LEU:HB3	1:C:67:ILE:HG12	1.94	0.49
1:C:74:LYS:HG2	1:C:87:LEU:CD2	2.42	0.49
1:A:231:ASN:O	1:A:232:LEU:HG	2.13	0.48
1:A:329:LEU:O	1:A:333:GLU:HG3	2.13	0.48
1:A:51:SER:HB3	1:A:54:VAL:CG2	2.44	0.48
1:B:92:ARG:HB2	1:B:92:ARG:CZ	2.42	0.48
1:C:260:ALA:O	1:C:262:LEU:HD22	2.13	0.48
1:A:58:TRP:HE3	1:A:58:TRP:HA	1.78	0.48
1:B:376:PHE:O	1:B:378:ASP:N	2.43	0.48
1:C:124:PRO:O	1:C:128:VAL:HG23	2.14	0.48
1:B:110:VAL:HG12	1:B:131:PHE:CG	2.48	0.48
1:B:167:GLU:HG2	1:B:430:LEU:HD12	1.95	0.48
1:C:235:VAL:HG21	1:C:269:PHE:CD1	2.48	0.48
1:C:192:GLU:HA	1:C:196:ILE:HB	1.94	0.48
1:A:233:VAL:HG12	1:A:234:VAL:N	2.28	0.48
1:B:167:GLU:HG3	1:B:437:ARG:NH1	2.28	0.48
1:B:480:ASP:HA	4:B:860:HOH:O	2.12	0.48
1:C:241:ILE:CD1	1:C:285:ILE:HG23	2.44	0.48
1:C:167:GLU:HG2	1:C:430:LEU:CD1	2.43	0.48
1:A:64:GLU:CA	1:A:102:ARG:HH12	2.27	0.48
1:A:144:LYS:HA	4:A:808:HOH:O	2.13	0.48
1:A:58:TRP:O	1:A:62:GLN:HG3	2.13	0.48
1:B:177:THR:O	1:B:181:LYS:HG3	2.14	0.48
1:A:209:LEU:O	1:A:212:PHE:HB2	2.13	0.48
1:B:225:SER:C	1:B:227:LEU:H	2.17	0.48
1:A:143:TRP:O	1:A:147:PHE:HD1	1.96	0.48
1:C:226:ARG:NH1	4:C:809:HOH:O	2.47	0.47
1:A:107:PRO:HA	1:A:110:VAL:HG22	1.96	0.47
1:C:182:GLU:O	1:C:185:LYS:HB3	2.14	0.47
1:A:415:ILE:HG13	1:A:441:LYS:HD2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:241:ILE:HD13	1:B:285:ILE:HG23	1.94	0.47
1:B:264:SER:HA	1:B:267:TYR:CZ	2.49	0.47
1:B:228:ARG:O	1:B:229:ASN:HB2	2.15	0.47
1:C:128:VAL:HG11	1:C:219:LEU:HD11	1.95	0.47
1:C:251:ASP:HB2	4:C:912:HOH:O	2.13	0.47
1:C:348:ASN:H	1:C:446:GLN:NE2	2.09	0.47
1:A:386:LEU:O	1:A:387:SER:CB	2.63	0.47
1:B:227:LEU:CB	1:B:230:GLU:HB2	2.44	0.47
1:C:250:GLU:O	1:C:250:GLU:HG3	2.13	0.47
1:C:224:PHE:CZ	3:C:500:M3L:HM12	2.49	0.47
1:A:464:TYR:HB3	4:A:838:HOH:O	2.15	0.47
1:C:104:TRP:CH2	1:C:269:PHE:N	2.81	0.47
1:A:57:PHE:CD1	1:A:146:TYR:HA	2.50	0.47
1:B:277:VAL:HG13	1:B:281:GLU:HB2	1.97	0.47
1:C:226:ARG:CD	1:C:252:HIS:NE2	2.79	0.47
1:A:193:GLN:HA	1:A:197:LEU:HD12	1.97	0.46
1:B:251:ASP:OD2	1:B:272:LYS:NZ	2.47	0.46
1:B:264:SER:HA	1:B:267:TYR:CE2	2.50	0.46
1:A:194:GLU:O	1:A:195:ILE:HD12	2.14	0.46
1:B:398:LYS:O	1:B:402:GLU:HG2	2.15	0.46
1:C:226:ARG:HH12	1:C:250:GLU:HB2	1.80	0.46
1:C:250:GLU:HG2	1:C:289:LEU:HD12	1.97	0.46
1:C:250:GLU:HB3	1:C:286:GLN:HG3	1.97	0.46
1:C:226:ARG:NE	1:C:252:HIS:CD2	2.83	0.46
1:A:101:LYS:HA	1:A:104:TRP:CE2	2.49	0.46
1:B:97:LEU:HB2	1:B:237:MET:CE	2.45	0.46
1:B:220:ARG:HH22	1:C:481:ILE:HG23	1.79	0.46
1:B:143:TRP:HB2	1:B:147:PHE:CE1	2.51	0.46
1:B:363:ALA:O	1:B:364:LEU:C	2.54	0.46
1:C:277:VAL:HG13	1:C:281:GLU:HB2	1.97	0.46
1:A:286:GLN:NE2	1:A:288:ASP:C	2.67	0.46
1:A:96:ILE:HD11	1:A:271:LEU:HD23	1.98	0.46
1:C:376:PHE:HA	4:C:859:HOH:O	2.16	0.46
1:B:301:GLY:HA3	4:B:804:HOH:O	2.15	0.46
1:C:99:VAL:O	1:C:104:TRP:HH2	1.99	0.46
1:A:106:ASN:HB2	1:A:107:PRO:CD	2.46	0.45
1:A:395:LEU:HD22	1:B:469:LEU:CD1	2.31	0.45
1:A:73:VAL:HG21	1:A:84:LEU:HB3	1.99	0.45
1:C:157:THR:OG1	1:C:177:THR:CG2	2.63	0.45
1:C:59:LYS:O	1:C:63:GLU:HG3	2.16	0.45
1:B:146:TYR:CE1	1:B:236:PRO:HA	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:227:LEU:HD22	1:A:227:LEU:N	2.31	0.45
1:A:482:LEU:HD11	1:C:179:SER:CB	2.31	0.45
1:B:257:LYS:HA	1:B:267:TYR:CD1	2.52	0.45
1:C:310:HIS:CD2	1:C:439:GLY:HA3	2.51	0.45
1:A:168:LEU:HD21	1:A:433:ALA:HA	1.99	0.45
1:A:54:VAL:HA	1:A:149:ILE:HD11	1.99	0.45
1:C:64:GLU:OE1	1:C:102:ARG:NH1	2.50	0.45
1:C:320:GLU:HA	1:C:325:PHE:CD1	2.51	0.45
1:B:313:THR:HB	1:B:344:ASP:OD1	2.17	0.45
1:C:158:ILE:HA	1:C:181:LYS:HD2	1.99	0.45
1:C:92:ARG:NH1	1:C:92:ARG:HB2	2.31	0.45
1:A:157:THR:HA	1:A:160:TRP:CD1	2.52	0.45
1:A:177:THR:HG22	1:A:298:LEU:HG	1.98	0.45
1:B:192:GLU:HA	1:B:196:ILE:HB	1.98	0.45
1:B:50:LEU:C	1:B:52:PRO:HD2	2.36	0.45
1:A:212:PHE:HB3	4:A:843:HOH:O	2.17	0.45
1:B:97:LEU:HB2	1:B:237:MET:HE1	1.98	0.45
1:B:310:HIS:CD2	1:B:439:GLY:HA3	2.51	0.45
1:B:376:PHE:C	1:B:378:ASP:H	2.20	0.45
1:C:94:ASP:O	1:C:272:LYS:HB2	2.17	0.45
1:C:92:ARG:HH11	1:C:92:ARG:HB2	1.82	0.45
1:A:146:TYR:CD2	4:A:845:HOH:O	2.57	0.44
1:A:67:ILE:HD11	4:A:852:HOH:O	2.17	0.44
1:C:97:LEU:HD12	1:C:97:LEU:C	2.37	0.44
1:A:73:VAL:CG2	1:A:84:LEU:HB3	2.47	0.44
1:C:241:ILE:HD13	1:C:285:ILE:HG23	1.99	0.44
1:C:381:TRP:O	1:C:385:GLU:HG3	2.18	0.44
1:B:329:LEU:O	1:B:333:GLU:HG3	2.17	0.44
3:B:502:M3L:HB2	3:B:502:M3L:HE2	1.75	0.44
1:B:49:SER:C	1:B:52:PRO:HD2	2.37	0.44
1:A:147:PHE:CA	4:A:845:HOH:O	2.59	0.44
1:B:167:GLU:HG2	1:B:430:LEU:CD1	2.47	0.44
1:B:255:GLU:HA	1:B:268:LEU:O	2.17	0.44
1:C:110:VAL:HG21	1:C:127:SER:HB3	1.98	0.44
1:C:97:LEU:HD13	1:C:237:MET:HE3	1.99	0.44
1:A:389:SER:HA	1:A:463:TYR:CG	2.52	0.44
1:B:439:GLY:HA2	1:B:442:MET:CE	2.41	0.44
1:A:363:ALA:HB2	1:B:469:LEU:HD11	2.00	0.44
1:C:143:TRP:HB2	1:C:147:PHE:CE1	2.53	0.44
1:B:64:GLU:OE2	1:B:64:GLU:HA	2.18	0.44
1:C:153:GLU:HG2	1:C:159:TYR:CD1	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:LEU:C	1:A:204:PRO:HD3	2.38	0.43
1:A:446:GLN:O	1:A:450:ILE:HG13	2.18	0.43
1:B:185:LYS:O	1:B:189:LEU:HG	2.18	0.43
1:C:244:SER:HB2	1:C:284:TYR:CG	2.54	0.43
1:A:242:ASN:HA	1:A:287:TYR:OH	2.19	0.43
1:C:99:VAL:O	1:C:104:TRP:CH2	2.71	0.43
1:C:252:HIS:N	1:C:252:HIS:CD2	2.85	0.43
1:C:348:ASN:ND2	4:C:816:HOH:O	2.34	0.43
1:C:146:TYR:CE1	1:C:236:PRO:HA	2.52	0.43
1:C:98:GLN:HA	1:C:269:PHE:O	2.17	0.43
1:C:313:THR:HB	1:C:344:ASP:OD1	2.18	0.43
1:C:167:GLU:HG2	1:C:430:LEU:HD12	2.00	0.43
1:A:54:VAL:CA	1:A:149:ILE:HD11	2.47	0.43
1:B:189:LEU:O	1:B:193:GLN:HG2	2.19	0.43
1:C:348:ASN:N	1:C:446:GLN:HE22	2.11	0.43
1:A:104:TRP:O	1:A:105:ILE:HG13	2.18	0.43
1:A:190:LYS:O	1:A:194:GLU:HB2	2.18	0.43
1:A:222:ARG:NH1	1:A:239:ASP:OD2	2.46	0.43
1:A:98:GLN:HA	1:A:269:PHE:O	2.18	0.43
1:A:176:THR:HG21	1:B:481:ILE:HD12	2.00	0.43
1:A:224:PHE:CE1	3:A:501:M3L:HM22	2.53	0.43
1:C:410:GLY:HA3	4:C:834:HOH:O	2.18	0.43
1:A:451:PHE:O	1:A:455:GLU:HG3	2.19	0.43
1:A:479:GLY:O	1:A:483:GLU:HG3	2.18	0.43
1:A:97:LEU:HD13	1:A:237:MET:HG3	2.01	0.43
1:C:54:VAL:HA	1:C:149:ILE:CD1	2.48	0.43
1:A:106:ASN:HB2	1:A:107:PRO:HD2	1.99	0.43
1:A:119:CYS:CB	1:A:122:LEU:HD12	2.48	0.43
1:B:426:LEU:HD23	1:B:431:ALA:HA	2.01	0.43
1:C:418:ASP:O	1:C:422:LYS:HG3	2.19	0.43
1:A:411:TYR:OH	1:A:440:GLU:OE1	2.25	0.43
1:C:371:LEU:HB3	1:C:380:ILE:CD1	2.43	0.43
1:B:167:GLU:HG3	1:B:437:ARG:HH12	1.84	0.43
1:B:104:TRP:CH2	1:B:269:PHE:HB2	2.54	0.43
1:B:488:GLN:HB3	4:B:864:HOH:O	2.19	0.43
1:C:100:PRO:HA	1:C:268:LEU:HB3	2.01	0.43
1:C:101:LYS:HA	1:C:104:TRP:CD2	2.53	0.43
1:C:438:GLU:OE2	1:C:442:MET:HE2	2.19	0.43
1:B:348:ASN:N	1:B:446:GLN:HE22	2.18	0.42
1:C:443:VAL:O	1:C:447:ILE:HG13	2.19	0.42
1:C:185:LYS:O	1:C:189:LEU:HG	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:TYR:CZ	1:B:479:GLY:HA3	2.54	0.42
1:C:286:GLN:NE2	1:C:288:ASP:C	2.70	0.42
1:A:176:THR:O	1:A:180:VAL:HG23	2.19	0.42
1:B:78:VAL:HG11	1:B:282:GLN:NE2	2.28	0.42
1:C:257:LYS:CB	1:C:257:LYS:NZ	2.83	0.42
1:B:157:THR:HA	1:B:160:TRP:CD1	2.55	0.42
1:A:68:THR:C	1:A:70:LYS:N	2.73	0.42
1:C:129:ILE:O	1:C:133:ILE:HG13	2.19	0.42
1:A:100:PRO:C	1:A:102:ARG:H	2.23	0.42
1:B:348:ASN:H	1:B:446:GLN:HE22	1.68	0.42
1:A:146:TYR:CE1	1:A:236:PRO:HA	2.54	0.42
1:B:395:LEU:HD13	1:C:469:LEU:HD11	2.01	0.42
1:C:287:TYR:CZ	3:C:500:M3L:HM33	2.55	0.42
1:B:427:ASP:CB	4:B:893:HOH:O	2.68	0.42
1:B:444:LEU:HA	1:B:444:LEU:HD12	1.77	0.42
1:B:257:LYS:HB3	1:B:258:GLY:H	1.61	0.42
1:A:483:GLU:C	1:A:485:LEU:H	2.23	0.41
1:B:130:LEU:HD11	1:B:191:LEU:HD22	2.01	0.41
1:B:441:LYS:NZ	4:B:806:HOH:O	2.51	0.41
1:C:386:LEU:O	1:C:387:SER:HB3	2.19	0.41
1:A:390:ARG:HB2	1:A:463:TYR:CE1	2.55	0.41
1:C:167:GLU:HG3	1:C:437:ARG:NH1	2.35	0.41
1:A:117:ARG:HB2	4:A:869:HOH:O	2.20	0.41
1:C:195:ILE:HD12	1:C:195:ILE:N	2.34	0.41
1:C:197:LEU:N	1:C:198:PRO:HD2	2.35	0.41
1:A:241:ILE:HG23	1:A:241:ILE:O	2.20	0.41
1:B:142:VAL:HG12	1:B:142:VAL:O	2.19	0.41
1:A:454:LYS:HE2	4:A:853:HOH:O	2.19	0.41
1:B:180:VAL:O	1:B:184:VAL:HG23	2.20	0.41
1:A:398:LYS:NZ	1:B:466:GLU:OE1	2.44	0.41
1:B:463:TYR:O	1:B:467:ARG:HG3	2.21	0.41
1:A:102:ARG:HG2	1:A:102:ARG:O	2.20	0.41
1:A:376:PHE:C	1:A:378:ASP:N	2.74	0.41
1:B:287:TYR:CE2	3:B:502:M3L:HD3	2.56	0.41
1:B:315:THR:HG23	1:C:473:GLY:O	2.20	0.41
1:A:151:PRO:HD2	1:A:222:ARG:NH2	2.35	0.41
1:A:394:GLU:HB2	1:A:458:LEU:HD21	2.03	0.41
1:B:395:LEU:HD11	1:C:465:GLN:HB3	2.03	0.41
1:B:61:LEU:CB	1:B:67:ILE:HG12	2.48	0.41
1:B:58:TRP:CZ2	1:B:74:LYS:HA	2.56	0.41
1:C:251:ASP:O	1:C:252:HIS:O	2.39	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:356:LEU:N	1:C:357:PRO:HD2	2.36	0.41
1:C:376:PHE:C	1:C:378:ASP:H	2.24	0.41
1:A:96:ILE:HD11	1:A:271:LEU:CD2	2.50	0.41
2:C:802:SAH:SD	3:C:500:M3L:HM31	2.60	0.41
1:A:62:GLN:HG2	1:A:67:ILE:HG23	2.03	0.41
1:B:51:SER:N	1:B:52:PRO:CD	2.83	0.41
1:C:150:LEU:HA	1:C:151:PRO:HD3	1.93	0.41
1:A:470:LYS:HE2	1:A:470:LYS:HA	2.02	0.41
1:C:82:LEU:O	2:C:802:SAH:N	2.54	0.41
1:A:102:ARG:HB2	4:A:829:HOH:O	2.21	0.41
1:B:427:ASP:OD2	1:B:428:SER:N	2.53	0.41
1:C:163:GLU:O	1:C:166:GLN:HB2	2.21	0.41
1:C:97:LEU:HD22	1:C:237:MET:SD	2.61	0.41
1:C:250:GLU:HB3	1:C:286:GLN:HG2	2.02	0.41
1:C:286:GLN:HG2	1:C:289:LEU:HG	2.03	0.40
1:A:218:ILE:O	1:A:222:ARG:HB2	2.22	0.40
1:A:53:ALA:O	1:A:57:PHE:HB2	2.21	0.40
1:A:90:ILE:HD13	1:A:96:ILE:HG22	2.03	0.40
1:B:188:CYS:HB3	1:B:212:PHE:CE1	2.56	0.40
1:B:390:ARG:HB2	1:B:463:TYR:CE1	2.57	0.40
1:C:389:SER:HB2	4:C:858:HOH:O	2.21	0.40
1:A:438:GLU:CG	1:A:442:MET:HE2	2.43	0.40
1:C:297:ALA:O	1:C:300:TYR:O	2.38	0.40
1:A:201:ARG:HH11	1:A:201:ARG:HG2	1.85	0.40
1:A:99:VAL:HG21	1:A:235:VAL:HG12	2.04	0.40
1:A:91:SER:HB2	1:A:94:ASP:OD1	2.22	0.40
1:A:100:PRO:C	1:A:102:ARG:N	2.75	0.40
1:A:101:LYS:HA	1:A:104:TRP:CZ3	2.56	0.40
1:A:60:TRP:O	1:A:64:GLU:HG2	2.21	0.40
1:B:367:THR:OG1	4:B:816:HOH:O	2.22	0.40
1:B:364:LEU:HD23	1:B:372:LEU:HD11	2.03	0.40
1:C:196:ILE:HD13	1:C:207:VAL:HG21	2.03	0.40
1:C:226:ARG:CD	1:C:252:HIS:CD2	3.05	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	418/440 (95%)	366 (88%)	44 (10%)	8 (2%)	8	6
1	B	438/440 (100%)	412 (94%)	19 (4%)	7 (2%)	9	8
1	C	436/440 (99%)	402 (92%)	26 (6%)	8 (2%)	8	6
All	All	1292/1320 (98%)	1180 (91%)	89 (7%)	23 (2%)	8	6

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	364	LEU
1	C	387	SER
1	A	73	VAL
1	A	105	ILE
1	A	377	ARG
1	A	387	SER
1	A	479	GLY
1	B	364	LEU
1	B	387	SER
1	C	252	HIS
1	C	470	LYS
1	A	52	PRO
1	A	236	PRO
1	A	471	ASP
1	B	377	ARG
1	C	232	LEU
1	B	226	ARG
1	B	288	ASP
1	B	366	GLY
1	C	377	ARG
1	C	105	ILE
1	C	479	GLY
1	B	236	PRO

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	374/386 (97%)	357 (96%)	17 (4%)	27	36
1	B	386/386 (100%)	370 (96%)	16 (4%)	30	40
1	C	384/386 (100%)	367 (96%)	17 (4%)	28	37
All	All	1144/1158 (99%)	1094 (96%)	50 (4%)	28	37

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	TRP
1	A	104	TRP
1	A	127	SER
1	A	143	TRP
1	A	209	LEU
1	A	237	MET
1	A	252	HIS
1	A	298	LEU
1	A	299	ASP
1	A	313	THR
1	A	315	THR
1	A	334	SER
1	A	378	ASP
1	A	395	LEU
1	A	444	LEU
1	A	453	GLN
1	A	480	ASP
1	B	80	GLU
1	B	94	ASP
1	B	104	TRP
1	B	143	TRP
1	B	209	LEU
1	B	210	ASP
1	B	226	ARG
1	B	236	PRO
1	B	313	THR

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Mol	Chain	Res	Type
1	B	315	THR
1	B	378	ASP
1	B	395	LEU
1	B	444	LEU
1	B	453	GLN
1	B	482	LEU
1	B	483	GLU
1	C	74	LYS
1	C	79	THR
1	C	104	TRP
1	C	143	TRP
1	C	209	LEU
1	C	210	ASP
1	C	219	LEU
1	C	236	PRO
1	C	257	LYS
1	C	268	LEU
1	C	313	THR
1	C	315	THR
1	C	378	ASP
1	C	395	LEU
1	C	444	LEU
1	C	453	GLN
1	C	468	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	GLN
1	A	93	ASN
1	A	98	GLN
1	A	106	ASN
1	A	145	HIS
1	A	152	GLN
1	A	169	GLN
1	A	286	GLN
1	B	55	GLN
1	B	106	ASN
1	B	152	GLN
1	B	169	GLN
1	B	286	GLN
1	B	310	HIS

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Mol	Chain	Res	Type
1	B	453	GLN
1	B	488	GLN
1	C	55	GLN
1	C	98	GLN
1	C	145	HIS
1	C	152	GLN
1	C	166	GLN
1	C	169	GLN
1	C	286	GLN
1	C	310	HIS
1	C	392	ASN
1	C	446	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](#)

6 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	M3L	C	500	-	8,12,12	2.40	2 (25%)	10,16,16	1.19	1 (10%)
3	M3L	B	502	-	8,12,12	0.73	0	10,16,16	0.38	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	SAH	B	801	-	21,28,28	1.18	3 (14%)	20,40,40	0.89	0
2	SAH	C	802	-	21,28,28	0.95	2 (9%)	20,40,40	0.90	0
3	M3L	A	501	-	8,12,12	0.71	0	10,16,16	0.46	0
2	SAH	A	800	-	21,28,28	0.95	2 (9%)	20,40,40	1.02	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	M3L	C	500	-	-	0/8/12/12	-
3	M3L	B	502	-	-	3/8/12/12	-
2	SAH	B	801	-	-	2/7/31/31	0/3/3/3
2	SAH	C	802	-	-	2/7/31/31	0/3/3/3
3	M3L	A	501	-	-	0/8/12/12	-
2	SAH	A	800	-	-	1/7/31/31	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	500	M3L	CD-CG	4.91	1.79	1.51
3	C	500	M3L	CG-CB	-4.22	1.34	1.52
2	B	801	SAH	C2-N3	3.10	1.37	1.32
2	B	801	SAH	C4-N3	3.06	1.39	1.35
2	A	800	SAH	C2-N3	2.70	1.36	1.32
2	C	802	SAH	C2-N3	2.55	1.36	1.32
2	C	802	SAH	C8-N7	-2.19	1.30	1.34
2	B	801	SAH	C8-N7	-2.06	1.31	1.34
2	A	800	SAH	C8-N7	-2.02	1.31	1.34

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	500	M3L	CG-CB-CA	-2.88	104.29	113.35

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	502	M3L	N-CA-CB-CG
3	B	502	M3L	C-CA-CB-CG
3	B	502	M3L	CE-CD-CG-CB
2	B	801	SAH	CB-CG-SD-C5'
2	B	801	SAH	C3'-C4'-C5'-SD
2	A	800	SAH	CB-CG-SD-C5'
2	C	802	SAH	CB-CG-SD-C5'
2	C	802	SAH	CA-CB-CG-SD

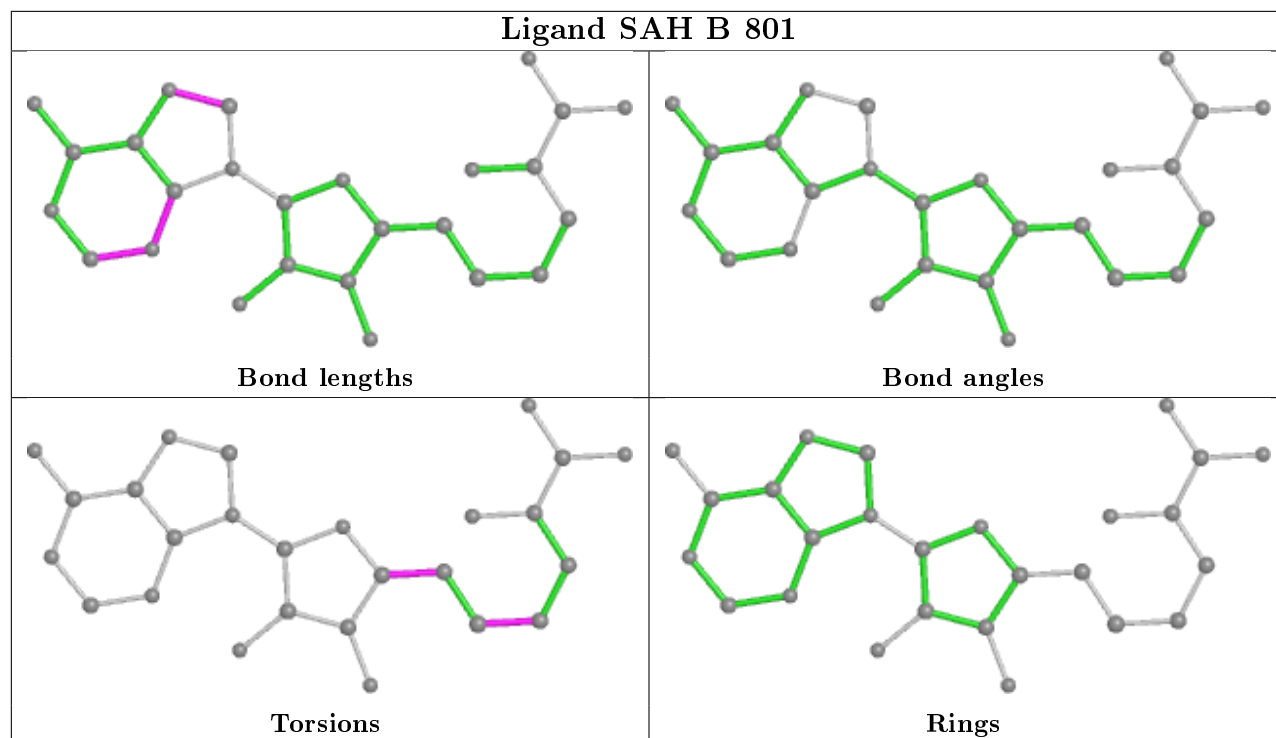
There are no ring outliers.

4 monomers are involved in 19 short contacts:

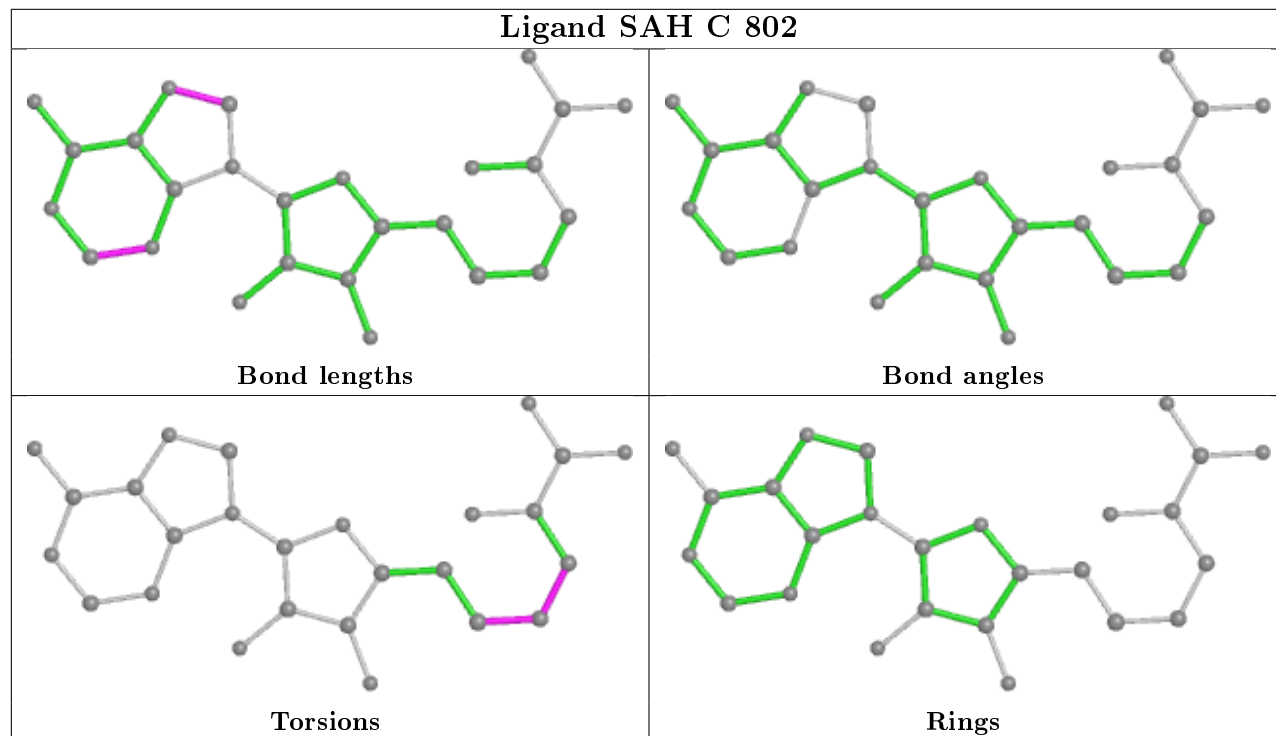
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	500	M3L	8	0
3	B	502	M3L	5	0
2	C	802	SAH	3	0
3	A	501	M3L	4	0

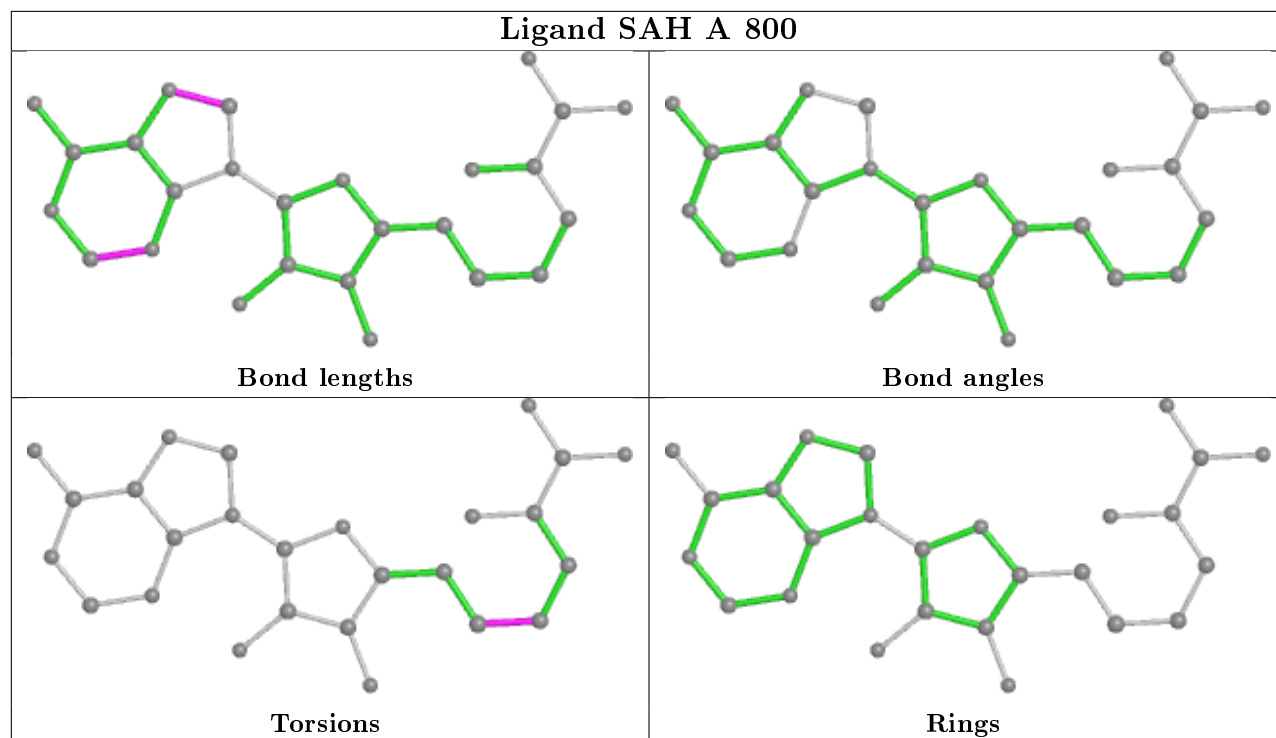
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 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 SAH B 801



## Ligand SAH C 802





## 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	424/440 (96%)	0.60	62 (14%) 2 1	45, 66, 123, 137	0
1	B	440/440 (100%)	0.09	22 (5%) 28 26	37, 60, 97, 127	0
1	C	438/440 (99%)	0.09	19 (4%) 35 32	40, 62, 95, 124	0
All	All	1302/1320 (98%)	0.25	103 (7%) 12 9	37, 62, 105, 137	0

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	487	PHE	8.2
1	A	267	TYR	7.0
1	A	140	ASP	6.9
1	A	112	ALA	6.5
1	C	250	GLU	6.4
1	A	145	HIS	6.3
1	A	139	GLU	5.1
1	C	471	ASP	5.1
1	B	486	TYR	5.0
1	A	59	LYS	4.9
1	A	117	ARG	4.9
1	B	480	ASP	4.5
1	A	206	PRO	4.2
1	A	144	LYS	4.2
1	A	148	GLY	4.1
1	A	254	TYR	4.0
1	B	471	ASP	4.0
1	C	112	ALA	4.0
1	A	231	ASN	4.0
1	A	60	TRP	3.9
1	A	486	TYR	3.9
1	A	201	ARG	3.9
1	C	486	TYR	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	57	PHE	3.8
1	A	471	ASP	3.8
1	A	63	GLU	3.8
1	B	226	ARG	3.6
1	A	124	PRO	3.6
1	A	50	LEU	3.4
1	C	139	GLU	3.4
1	A	55	GLN	3.3
1	A	114	GLU	3.3
1	A	116	GLY	3.3
1	A	137	SER	3.3
1	A	118	VAL	3.2
1	C	251	ASP	3.2
1	C	70	LYS	3.2
1	A	205	ASP	3.1
1	A	211	ASP	3.1
1	C	117	ARG	3.1
1	A	197	LEU	3.1
1	A	256	VAL	3.0
1	C	193	GLN	2.9
1	A	108	ASP	2.9
1	A	128	VAL	2.9
1	B	488	GLN	2.9
1	B	193	GLN	2.8
1	C	121	GLU	2.8
1	A	65	GLY	2.8
1	B	375	LEU	2.8
1	B	143	TRP	2.8
1	A	183	TYR	2.8
1	A	141	SER	2.8
1	A	268	LEU	2.7
1	C	69	ALA	2.7
1	A	92	ARG	2.7
1	B	55	GLN	2.7
1	A	70	LYS	2.6
1	A	138	ARG	2.6
1	A	269	PHE	2.5
1	A	204	PRO	2.5
1	A	136	ARG	2.5
1	A	62	GLN	2.5
1	A	152	GLN	2.5
1	B	117	ARG	2.5

*Continued on next page...*



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Mol	Chain	Res	Type	RSRZ
1	B	229	ASN	2.5
1	A	200	LYS	2.5
1	B	228	ARG	2.5
1	A	196	ILE	2.4
1	B	378	ASP	2.4
1	A	119	CYS	2.4
1	B	336	GLY	2.4
1	A	203	PHE	2.4
1	C	200	LYS	2.3
1	A	227	LEU	2.3
1	A	234	VAL	2.3
1	A	377	ARG	2.3
1	B	104	TRP	2.3
1	C	144	LYS	2.3
1	A	225	SER	2.3
1	A	69	ALA	2.3
1	A	470	LYS	2.2
1	B	201	ARG	2.2
1	B	121	GLU	2.2
1	A	250	GLU	2.2
1	B	405	LYS	2.2
1	B	409	ALA	2.2
1	C	59	LYS	2.1
1	A	212	PHE	2.1
1	C	483	GLU	2.1
1	A	195	ILE	2.1
1	B	460	GLN	2.1
1	B	163	GLU	2.1
1	A	102	ARG	2.1
1	C	55	GLN	2.1
1	A	185	LYS	2.1
1	A	111	ALA	2.0
1	A	113	SER	2.0
1	C	229	ASN	2.0
1	B	481	ILE	2.0
1	A	58	TRP	2.0
1	A	202	LEU	2.0
1	C	63	GLU	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 ⓘ

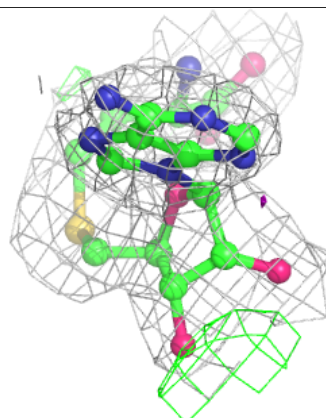
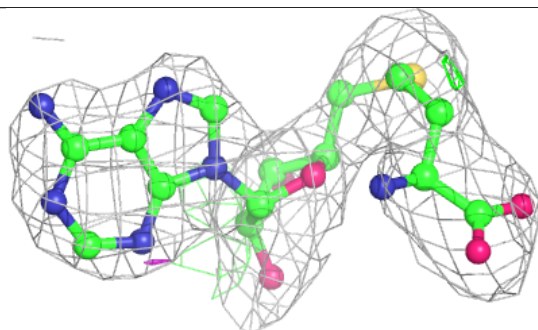
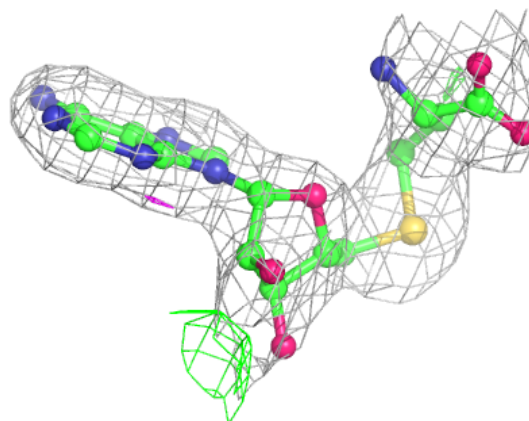
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	M3L	A	501	13/13	0.73	0.78	148,161,189,191	0
3	M3L	B	502	13/13	0.82	0.66	111,128,164,165	0
3	M3L	C	500	13/13	0.84	0.63	86,110,156,158	0
2	SAH	A	800	26/26	0.93	0.14	57,59,70,72	0
2	SAH	C	802	26/26	0.96	0.13	43,48,49,50	0
2	SAH	B	801	26/26	0.97	0.10	36,40,43,46	0

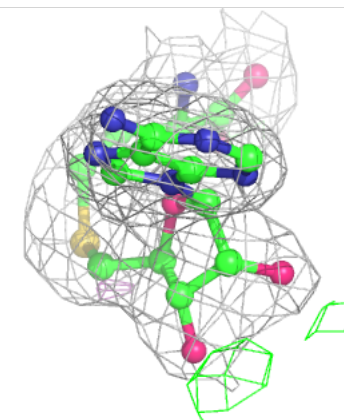
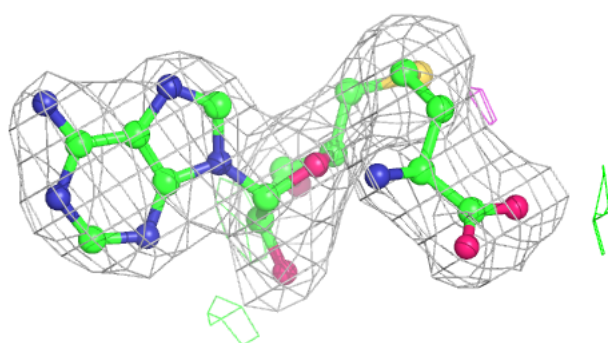
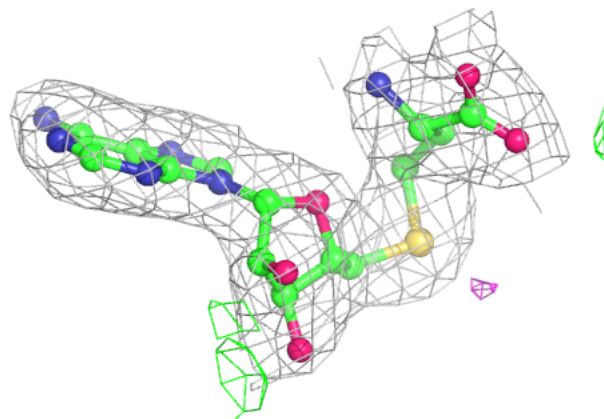
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

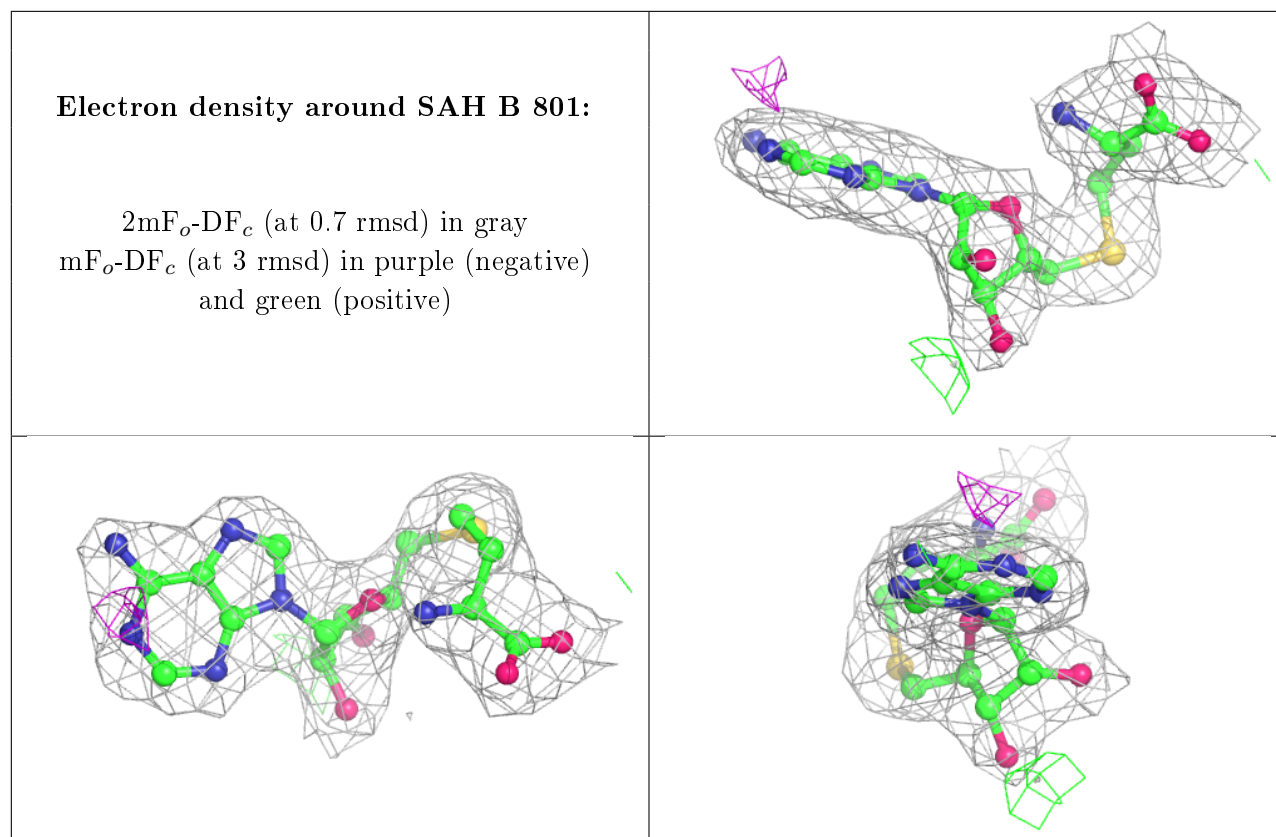
**Electron density around SAH A 800:**

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

**Electron density around SAH C 802:**

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





## 6.5 Other polymers [i](#)

There are no such residues in this entry.