



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 09:56 PM BST

PDB ID : 5C71  
Title : The structure of *Aspergillus oryzae*  $\alpha$ -glucuronidase complexed with glycyrrhetic acid monoglucuronide  
Authors : Sun, H.L.; Lv, B.; Huang, S.; Li, C.; Jiang, T.  
Deposited on : 2015-06-24  
Resolution : 2.62 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 : 2.13.1  
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) : 2.13.1

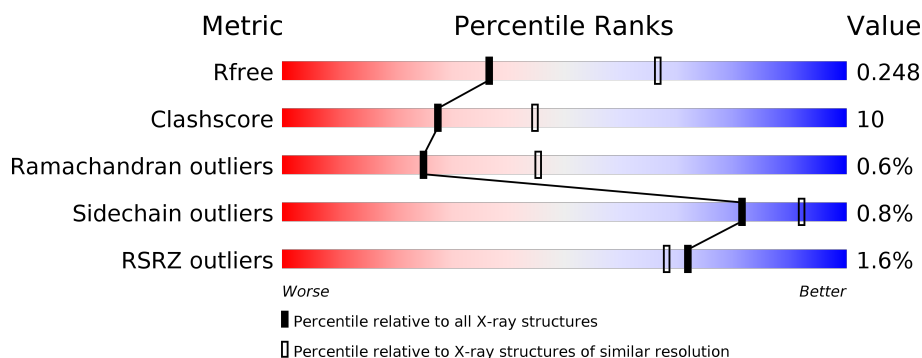
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.62 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	3797 (2.64-2.60)
Clashscore	141614	4168 (2.64-2.60)
Ramachandran outliers	138981	4093 (2.64-2.60)
Sidechain outliers	138945	4093 (2.64-2.60)
RSRZ outliers	127900	3731 (2.64-2.60)

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	637	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>72%</span> <span>18%</span> <span>• 8%</span> </div> </div>
1	B	637	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>2%</span> <span>78%</span> <span>12%</span> <span>• 9%</span> </div> </div>
1	C	637	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>75%</span> <span>15%</span> <span>• 8%</span> </div> </div>
1	D	637	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>77%</span> <span>12%</span> <span>• 9%</span> </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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	CBW	C	701	-	-	X	X
3	GCU	A	702	-	-	X	X
3	GCU	C	702	-	-	X	-

## 2 Entry composition [i](#)

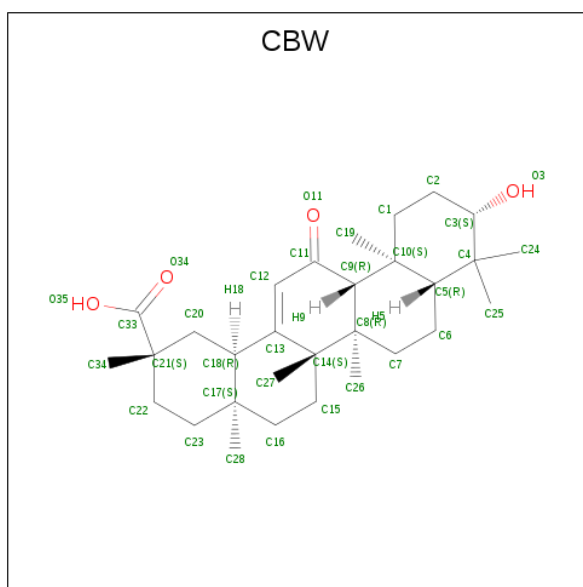
There are 4 unique types of molecules in this entry. The entry contains 18844 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glucuronidase.

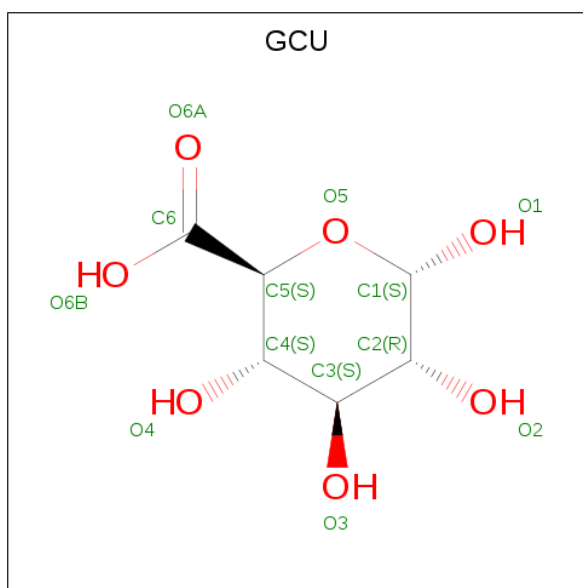
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	583	Total	C	N	O	S	0	0	0
			4648	2949	802	884	13			
1	B	582	Total	C	N	O	S	0	0	0
			4638	2943	799	883	13			
1	C	583	Total	C	N	O	S	0	0	0
			4648	2949	802	884	13			
1	D	582	Total	C	N	O	S	0	0	0
			4638	2943	799	883	13			

- Molecule 2 is (3BETA,5BETA,14BETA)-3-HYDROXY-11-OXOOLEAN-12-EN-29-OIC ACID (three-letter code: CBW) (formula:  $C_{30}H_{46}O_4$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			34	30	4		
2	C	1	Total	C	O	0	0
			34	30	4		

- Molecule 3 is alpha-D-glucopyranuronic acid (three-letter code: GCU) (formula: C<sub>6</sub>H<sub>10</sub>O<sub>7</sub>).

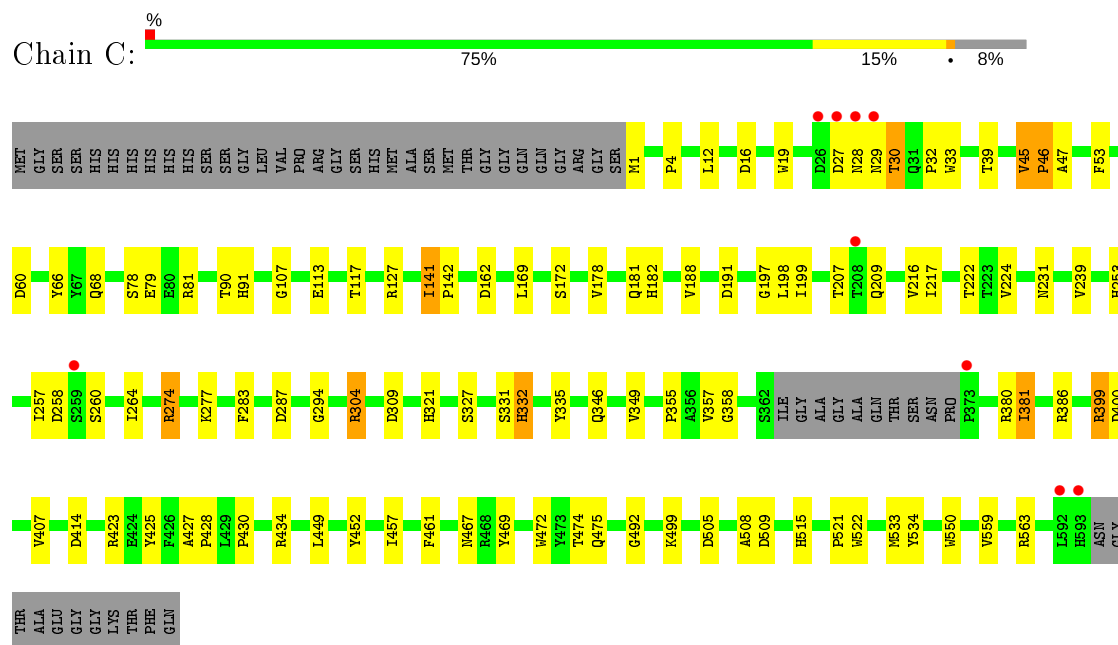


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	6	6		
3	C	1	Total	C	O	0	0
			12	6	6		

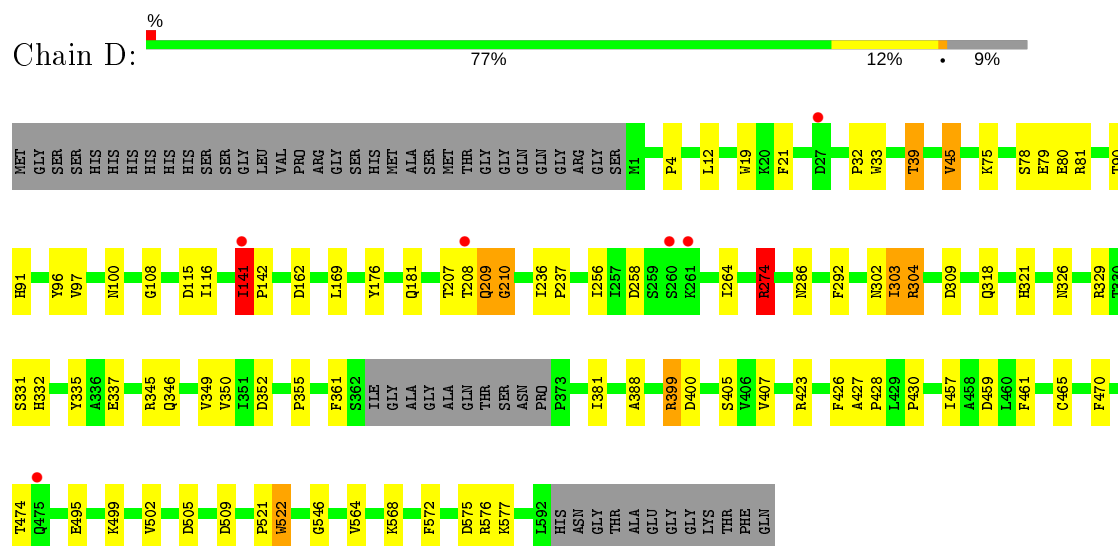
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	49	Total	O	0	0
			49	49		
4	B	44	Total	O	0	0
			44	44		
4	C	41	Total	O	0	0
			41	41		
4	D	46	Total	O	0	0
			46	46		





• Molecule 1: Glucuronidase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	96.07Å 96.23Å 96.19Å 88.27° 74.36° 71.13°	Depositor
Resolution (Å)	46.23 – 2.62 47.13 – 2.62	Depositor EDS
% Data completeness (in resolution range)	98.0 (46.23-2.62) 97.2 (47.13-2.62)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	8.04 (at 2.61Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, $R_{free}$	0.187 , 0.248 0.187 , 0.248	Depositor DCC
$R_{free}$ test set	4540 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	18.8	Xtriage
Anisotropy	0.050	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 33.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.017 for -h,-l,-k	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	18844	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.54% 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

### 5.1 Standard geometry

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

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	0.73	1/4769 (0.0%)	0.91	13/6496 (0.2%)
1	B	0.70	0/4758	0.89	8/6481 (0.1%)
1	C	0.70	1/4769 (0.0%)	0.88	8/6496 (0.1%)
1	D	0.70	0/4758	0.90	19/6481 (0.3%)
All	All	0.71	2/19054 (0.0%)	0.89	48/25954 (0.2%)

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	1
1	B	0	2
1	D	0	2
All	All	0	5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	244	PRO	N-CD	-6.35	1.39	1.47
1	C	107	GLY	N-CA	5.36	1.54	1.46

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	304	ARG	NE-CZ-NH2	-10.06	115.27	120.30
1	B	304	ARG	NE-CZ-NH1	8.89	124.75	120.30
1	A	45	VAL	C-N-CD	8.71	146.70	128.40
1	A	304	ARG	NE-CZ-NH2	-8.51	116.05	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	141	ILE	C-N-CD	8.04	145.27	128.40
1	C	399	ARG	NE-CZ-NH1	7.82	124.21	120.30
1	D	304	ARG	NE-CZ-NH2	-7.69	116.45	120.30
1	A	329	ARG	NE-CZ-NH1	7.46	124.03	120.30
1	C	304	ARG	NE-CZ-NH2	-7.17	116.72	120.30
1	A	399	ARG	NE-CZ-NH1	7.14	123.87	120.30
1	D	329	ARG	NE-CZ-NH2	-7.13	116.74	120.30
1	D	39	THR	CB-CA-C	-6.88	93.01	111.60
1	A	440	ARG	NE-CZ-NH1	6.84	123.72	120.30
1	D	505	ASP	CB-CG-OD1	6.38	124.04	118.30
1	B	399	ARG	NE-CZ-NH1	6.34	123.47	120.30
1	C	45	VAL	C-N-CD	6.23	141.49	128.40
1	A	304	ARG	NE-CZ-NH1	6.20	123.40	120.30
1	D	304	ARG	NE-CZ-NH1	6.15	123.38	120.30
1	D	141	ILE	C-N-CD	6.14	141.30	128.40
1	C	304	ARG	NE-CZ-NH1	6.11	123.35	120.30
1	B	141	ILE	C-N-CD	6.10	141.20	128.40
1	B	575	ASP	CB-CG-OD1	6.09	123.78	118.30
1	A	70	ASP	CB-CG-OD1	6.01	123.71	118.30
1	A	141	ILE	C-N-CD	5.96	140.93	128.40
1	D	329	ARG	NE-CZ-NH1	5.79	123.20	120.30
1	A	141	ILE	C-N-CA	-5.71	98.02	122.00
1	A	60	ASP	CB-CG-OD1	5.66	123.39	118.30
1	D	39	THR	N-CA-C	5.57	126.04	111.00
1	A	45	VAL	C-N-CA	-5.51	98.84	122.00
1	D	274	ARG	NE-CZ-NH2	-5.49	117.56	120.30
1	D	45	VAL	C-N-CA	-5.47	99.01	122.00
1	D	274	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	C	141	ILE	C-N-CA	-5.43	99.18	122.00
1	D	80	GLU	N-CA-CB	5.43	120.37	110.60
1	D	274	ARG	CG-CD-NE	5.42	123.19	111.80
1	C	60	ASP	CB-CG-OD1	5.37	123.13	118.30
1	D	522	TRP	CB-CA-C	-5.32	99.75	110.40
1	D	345	ARG	NE-CZ-NH2	-5.30	117.65	120.30
1	B	304	ARG	CG-CD-NE	-5.30	100.67	111.80
1	B	414	ASP	CB-CG-OD1	5.23	123.01	118.30
1	D	459	ASP	CB-CG-OD1	5.23	123.01	118.30
1	C	434	ARG	NE-CZ-NH1	5.21	122.91	120.30
1	B	345	ARG	NE-CZ-NH2	-5.20	117.70	120.30
1	A	505	ASP	CB-CG-OD1	5.18	122.97	118.30
1	A	380	ARG	N-CA-C	5.10	124.77	111.00
1	D	345	ARG	NE-CZ-NH1	5.07	122.83	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	399	ARG	CG-CD-NE	5.04	122.39	111.80
1	D	274	ARG	CD-NE-CZ	5.01	130.61	123.60

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	141	ILE	Peptide
1	B	245	GLY	Peptide
1	B	295	PHE	Peptide
1	D	141	ILE	Peptide
1	D	210	GLY	Peptide

## 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4648	0	4472	110	0
1	B	4638	0	4465	71	0
1	C	4648	0	4472	93	0
1	D	4638	0	4465	78	0
2	A	34	0	44	19	0
2	C	34	0	44	25	0
3	A	12	0	7	11	0
3	C	12	0	7	10	0
4	A	49	0	0	0	0
4	B	44	0	0	0	0
4	C	41	0	0	1	0
4	D	46	0	0	1	0
All	All	18844	0	17976	367	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 10.

All (367) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:449:LEU:HD23	2:C:701:CBW:C34	1.75	1.16
2:A:701:CBW:H191	2:A:701:CBW:H263	1.30	1.13
1:C:449:LEU:CD2	2:C:701:CBW:H342	1.79	1.11
2:C:701:CBW:H343	2:C:701:CBW:H161	1.22	1.10
2:C:701:CBW:H263	2:C:701:CBW:H191	1.30	1.08
2:A:701:CBW:H192	2:A:701:CBW:H253	1.36	1.07
1:C:45:VAL:O	1:C:47:ALA:N	1.87	1.07
1:D:90:THR:HG22	1:D:108:GLY:O	1.51	1.07
2:C:701:CBW:H253	2:C:701:CBW:H192	1.36	1.05
2:A:701:CBW:H252	3:A:702:GCU:H2	1.39	1.04
1:C:197:GLY:HA3	1:C:239:VAL:HG21	1.41	1.02
1:A:423:ARG:HA	1:A:457:ILE:CD1	1.91	0.99
1:A:141:ILE:HG23	1:A:142:PRO:CD	1.95	0.96
1:A:349:VAL:HG13	1:A:407:VAL:HG11	1.47	0.95
1:A:162:ASP:HB2	3:A:702:GCU:H3	1.48	0.95
1:D:19:TRP:CD1	1:D:45:VAL:HG21	2.02	0.94
1:B:423:ARG:HA	1:B:457:ILE:CD1	2.00	0.92
1:D:39:THR:O	1:D:39:THR:HG23	1.68	0.92
2:C:701:CBW:H343	2:C:701:CBW:C16	2.01	0.91
1:A:423:ARG:CA	1:A:457:ILE:HD11	2.03	0.89
2:C:701:CBW:C34	2:C:701:CBW:H161	2.02	0.89
1:D:423:ARG:HA	1:D:457:ILE:HD11	1.55	0.89
1:D:141:ILE:CD1	1:D:355:PRO:HB2	2.04	0.88
1:A:45:VAL:HG13	1:A:46:PRO:HD3	1.56	0.88
1:D:349:VAL:HG13	1:D:407:VAL:HG11	1.56	0.87
1:D:19:TRP:CD1	1:D:45:VAL:CG2	2.58	0.86
1:A:16:ASP:HA	1:A:45:VAL:HG11	1.58	0.85
1:C:449:LEU:CD2	2:C:701:CBW:C34	2.45	0.85
1:C:449:LEU:HD23	2:C:701:CBW:H342	0.90	0.84
1:A:423:ARG:HA	1:A:457:ILE:HD11	1.59	0.84
1:C:141:ILE:HG23	1:C:142:PRO:CD	2.09	0.81
1:B:274:ARG:HG2	1:B:274:ARG:HH11	1.46	0.81
2:A:701:CBW:H252	3:A:702:GCU:C2	2.10	0.81
1:A:357:VAL:HG12	1:A:413:ASN:HB3	1.63	0.80
1:A:197:GLY:HA3	1:A:239:VAL:HG11	1.64	0.80
1:D:274:ARG:HG2	1:D:274:ARG:HH11	1.45	0.79
1:C:563:ARG:NH2	3:C:702:GCU:O6A	2.15	0.79
1:A:162:ASP:CB	3:A:702:GCU:H3	2.14	0.77
1:A:521:PRO:O	1:A:522:TRP:HB2	1.85	0.77
1:C:509:ASP:O	1:C:522:TRP:HA	1.85	0.77
1:C:550:TRP:CE2	3:C:702:GCU:O4	2.37	0.77
1:C:29:ASN:O	1:C:30:THR:C	2.25	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:423:ARG:CA	1:B:457:ILE:HD11	2.16	0.75
1:A:423:ARG:HA	1:A:457:ILE:HD12	1.68	0.75
1:A:141:ILE:HG23	1:A:142:PRO:HD3	1.69	0.73
2:A:701:CBW:C19	2:A:701:CBW:H263	2.15	0.73
1:A:405:SER:O	1:A:407:VAL:HG13	1.89	0.73
2:A:701:CBW:C19	2:A:701:CBW:H253	2.16	0.73
1:D:141:ILE:HD12	1:D:142:PRO:HD3	1.70	0.73
1:C:162:ASP:HB2	3:C:702:GCU:H3	1.70	0.72
1:B:258:ASP:O	1:B:260:SER:O	2.07	0.72
2:A:701:CBW:C26	2:A:701:CBW:H191	2.16	0.72
2:C:701:CBW:H253	2:C:701:CBW:C19	2.16	0.72
1:B:423:ARG:HA	1:B:457:ILE:HD11	1.70	0.72
2:C:701:CBW:H263	2:C:701:CBW:C19	2.15	0.72
1:C:32:PRO:HG2	1:C:66:TYR:CE2	2.25	0.72
1:A:309:ASP:OD2	1:B:304:ARG:NH2	2.23	0.72
1:D:141:ILE:CD1	1:D:142:PRO:HD3	2.20	0.72
1:C:335:TYR:O	1:C:399:ARG:NH2	2.22	0.71
2:C:701:CBW:H191	2:C:701:CBW:C26	2.17	0.71
1:D:399:ARG:HD3	1:D:400:ASP:OD1	1.91	0.70
1:A:309:ASP:CG	1:B:304:ARG:HH22	1.94	0.70
1:A:349:VAL:HG13	1:A:407:VAL:CG1	2.22	0.70
1:D:405:SER:O	1:D:407:VAL:HG13	1.91	0.69
1:A:45:VAL:HG13	1:A:46:PRO:CD	2.22	0.69
1:D:141:ILE:HD12	1:D:355:PRO:HB2	1.74	0.69
1:C:423:ARG:HA	1:C:457:ILE:HD11	1.74	0.69
1:D:207:THR:OG1	1:D:209:GLN:HG2	1.92	0.69
1:C:550:TRP:CZ2	3:C:702:GCU:O4	2.44	0.69
1:B:64:TRP:CH2	1:B:92:HIS:CD2	2.81	0.69
1:C:141:ILE:HG23	1:C:142:PRO:HD3	1.75	0.68
1:B:33:TRP:O	1:B:127:ARG:NH1	2.24	0.67
1:B:64:TRP:CH2	1:B:92:HIS:HD2	2.11	0.67
1:C:197:GLY:CA	1:C:239:VAL:HG21	2.22	0.67
1:D:361:PHE:HE1	1:D:381:ILE:HD13	1.59	0.67
1:D:39:THR:O	1:D:39:THR:CG2	2.42	0.67
1:C:258:ASP:HB3	1:C:264:ILE:HD11	1.77	0.67
1:D:575:ASP:O	1:D:576:ARG:HB2	1.95	0.67
1:A:277:LYS:NZ	1:A:287:ASP:OD1	2.28	0.67
1:A:16:ASP:HA	1:A:45:VAL:CG1	2.23	0.67
1:B:521:PRO:O	1:B:522:TRP:HB2	1.93	0.67
1:B:423:ARG:CA	1:B:457:ILE:CD1	2.74	0.66
2:A:701:CBW:C25	2:A:701:CBW:H192	2.22	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:29:ASN:C	1:C:30:THR:O	2.29	0.65
1:C:78:SER:O	1:C:79:GLU:HB2	1.94	0.65
4:C:813:HOH:O	1:D:75:LYS:HE2	1.96	0.65
1:A:559:VAL:HG13	1:C:521:PRO:HG2	1.78	0.65
1:B:141:ILE:HG23	1:B:142:PRO:HD3	1.78	0.65
1:B:141:ILE:O	1:B:143:PRO:HD3	1.97	0.65
1:D:19:TRP:CG	1:D:45:VAL:HG21	2.32	0.65
1:A:274:ARG:HH12	1:A:276:VAL:HG22	1.62	0.65
1:D:350:VAL:O	1:D:407:VAL:HG22	1.97	0.65
1:D:521:PRO:O	1:D:522:TRP:HB2	1.96	0.64
1:A:197:GLY:H	1:A:239:VAL:HG13	1.62	0.64
1:C:521:PRO:O	1:C:522:TRP:HB2	1.98	0.64
1:C:45:VAL:O	1:C:46:PRO:C	2.32	0.63
2:C:701:CBW:H252	3:C:702:GCU:C1	2.28	0.63
1:A:423:ARG:CA	1:A:457:ILE:CD1	2.65	0.63
1:A:423:ARG:N	1:A:457:ILE:HD11	2.13	0.63
1:B:423:ARG:N	1:B:457:ILE:HD11	2.13	0.63
2:A:701:CBW:H252	3:A:702:GCU:C1	2.28	0.63
1:B:335:TYR:O	1:B:399:ARG:NH2	2.23	0.63
1:D:19:TRP:CD1	1:D:45:VAL:HG23	2.34	0.63
1:C:162:ASP:HB2	3:C:702:GCU:C3	2.28	0.62
1:A:138:TYR:HB3	1:A:380:ARG:O	1.98	0.62
1:B:91:HIS:O	1:B:106:VAL:O	2.16	0.62
1:D:349:VAL:HG13	1:D:407:VAL:CG1	2.28	0.62
1:A:559:VAL:CG1	1:C:521:PRO:HG2	2.30	0.61
1:A:102:VAL:HG12	1:A:114:ALA:HB1	1.82	0.61
1:A:45:VAL:O	1:A:47:ALA:N	2.34	0.61
1:C:469:TYR:CE1	2:C:701:CBW:H21C	2.37	0.60
1:B:457:ILE:HG22	1:B:461:PHE:HE1	1.66	0.60
1:A:564:VAL:O	1:A:568:LYS:NZ	2.35	0.59
1:C:304:ARG:NH2	1:D:309:ASP:OD2	2.34	0.59
1:D:19:TRP:NE1	1:D:45:VAL:HG21	2.17	0.59
1:B:423:ARG:HA	1:B:457:ILE:HD12	1.83	0.59
1:A:350:VAL:O	1:A:407:VAL:HG22	2.02	0.59
1:A:474:THR:OG1	1:A:475:GLN:OE1	2.21	0.59
1:A:257:ILE:HG23	1:A:261:LYS:HA	1.84	0.59
1:B:399:ARG:HD3	1:B:400:ASP:OD1	2.03	0.59
2:C:701:CBW:C25	2:C:701:CBW:H192	2.22	0.59
1:C:178:VAL:CG1	1:C:182:HIS:CG	2.86	0.58
1:C:304:ARG:HH22	1:D:309:ASP:CG	2.06	0.58
1:A:304:ARG:NH2	1:B:309:ASP:OD2	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:274:ARG:HH12	1:A:276:VAL:CG2	2.15	0.58
1:D:141:ILE:HD11	1:D:355:PRO:HB2	1.85	0.58
1:B:349:VAL:CG1	1:B:407:VAL:HG21	2.33	0.58
1:C:309:ASP:OD2	1:D:304:ARG:NH2	2.37	0.58
1:B:141:ILE:CG2	1:B:142:PRO:HD3	2.33	0.58
1:C:178:VAL:HG11	1:C:182:HIS:CG	2.39	0.58
1:D:302:ASN:OD1	1:D:303:ILE:HG13	2.04	0.58
1:D:12:LEU:HD12	1:D:176:TYR:HB3	1.86	0.58
1:A:181:GLN:HB2	1:A:264:ILE:HD11	1.86	0.57
1:B:181:GLN:HG2	1:B:207:THR:CG2	2.34	0.57
1:D:21:PHE:O	1:D:39:THR:HG21	2.04	0.57
1:C:467:ASN:HB3	1:C:505:ASP:HB2	1.88	0.56
1:A:349:VAL:CG1	1:A:407:VAL:HG21	2.35	0.56
1:C:472:TRP:CZ2	1:C:509:ASP:HB2	2.41	0.56
1:A:533:MET:CE	1:A:534:TYR:CD1	2.90	0.56
1:A:467:ASN:HB3	1:A:505:ASP:HB2	1.87	0.55
1:B:258:ASP:HB3	1:B:264:ILE:HD11	1.89	0.55
1:C:178:VAL:CG1	1:C:182:HIS:CB	2.85	0.55
2:A:701:CBW:C25	3:A:702:GCU:C1	2.85	0.55
1:A:242:TRP:CE3	1:A:274:ARG:HD3	2.43	0.54
1:A:162:ASP:OD2	3:A:702:GCU:C3	2.55	0.54
2:C:701:CBW:C25	3:C:702:GCU:C1	2.85	0.54
1:D:423:ARG:HA	1:D:457:ILE:CD1	2.34	0.54
1:A:188:VAL:HG13	1:A:199:ILE:HG23	1.90	0.54
1:C:277:LYS:NZ	1:C:287:ASP:OD1	2.36	0.54
1:C:309:ASP:CG	1:D:304:ARG:HH22	2.11	0.54
1:B:141:ILE:CG2	1:B:142:PRO:CD	2.86	0.54
1:D:303:ILE:O	1:D:303:ILE:HG22	2.06	0.53
1:C:216:VAL:O	1:C:224:VAL:HG22	2.08	0.53
1:A:141:ILE:HG23	1:A:142:PRO:HD2	1.86	0.53
1:A:141:ILE:CG2	1:A:142:PRO:CD	2.79	0.53
1:A:22:ALA:HB2	1:A:39:THR:HG21	1.89	0.53
1:C:423:ARG:HA	1:C:457:ILE:CD1	2.38	0.53
1:D:461:PHE:O	1:D:499:LYS:NZ	2.36	0.53
1:B:188:VAL:HG13	1:B:199:ILE:HG23	1.90	0.53
1:A:17:GLY:H	1:A:45:VAL:HG12	1.73	0.53
1:C:188:VAL:HG13	1:C:199:ILE:HG23	1.91	0.53
1:C:509:ASP:HB3	1:C:522:TRP:CZ3	2.44	0.53
1:A:304:ARG:HH22	1:B:309:ASP:CG	2.11	0.52
1:B:141:ILE:HG23	1:B:142:PRO:CD	2.38	0.52
1:A:335:TYR:O	1:A:399:ARG:NH2	2.31	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:378:PRO:O	1:A:382:ASN:ND2	2.42	0.52
1:C:45:VAL:O	1:C:169:LEU:O	2.28	0.52
1:D:181:GLN:HG2	1:D:207:THR:HG21	1.90	0.52
1:B:405:SER:O	1:B:407:VAL:HG13	2.08	0.52
1:B:16:ASP:OD2	1:B:172:SER:OG	2.21	0.52
1:C:222:THR:O	1:C:224:VAL:HG13	2.10	0.52
1:A:274:ARG:HG2	1:A:275:THR:N	2.25	0.52
1:A:45:VAL:CG1	1:A:46:PRO:HD3	2.35	0.52
1:B:334:PRO:HG3	1:B:396:LEU:HD13	1.91	0.52
1:A:533:MET:HE1	1:A:534:TYR:CD1	2.45	0.52
1:D:45:VAL:O	1:D:169:LEU:O	2.28	0.52
1:D:181:GLN:HE21	1:D:207:THR:HG23	1.74	0.52
1:D:78:SER:C	1:D:79:GLU:HG2	2.31	0.52
1:D:349:VAL:CG1	1:D:407:VAL:HG21	2.39	0.51
1:C:19:TRP:CD1	1:C:45:VAL:HG13	2.45	0.51
1:D:352:ASP:OD2	1:D:400:ASP:OD2	2.27	0.51
1:A:457:ILE:HG22	1:A:461:PHE:HE1	1.76	0.51
1:A:427:ALA:HB3	1:A:428:PRO:HD3	1.92	0.51
2:A:701:CBW:H271	2:A:701:CBW:H202	1.92	0.51
1:C:258:ASP:O	1:C:260:SER:O	2.28	0.51
1:D:81:ARG:NH1	1:D:115:ASP:OD2	2.44	0.51
1:A:521:PRO:O	1:A:522:TRP:CB	2.57	0.51
1:A:472:TRP:CZ2	1:A:509:ASP:HB2	2.46	0.51
2:C:701:CBW:H271	2:C:701:CBW:H202	1.93	0.51
1:A:45:VAL:O	1:A:169:LEU:O	2.30	0.50
1:C:197:GLY:HA3	1:C:239:VAL:CG2	2.28	0.50
1:A:244:PRO:O	1:A:347:GLY:O	2.30	0.50
1:B:456:ARG:O	1:B:457:ILE:HD13	2.12	0.50
1:C:399:ARG:NH1	1:C:400:ASP:OD2	2.45	0.50
1:C:452:TYR:CD1	1:C:492:GLY:HA3	2.47	0.50
1:B:349:VAL:HG13	1:B:407:VAL:HG21	1.93	0.50
1:D:572:PHE:HA	1:D:577:LYS:O	2.12	0.50
1:C:33:TRP:O	1:C:127:ARG:NH1	2.45	0.49
1:C:533:MET:O	1:C:533:MET:HG2	2.04	0.49
2:C:701:CBW:C25	2:C:701:CBW:C19	2.85	0.49
1:B:178:VAL:HG22	1:B:182:HIS:HB3	1.92	0.49
1:A:181:GLN:HB2	1:A:264:ILE:CD1	2.41	0.49
1:A:78:SER:O	1:A:79:GLU:HB2	2.12	0.49
1:C:29:ASN:O	1:C:30:THR:O	2.31	0.49
1:D:509:ASP:HB3	1:D:522:TRP:CZ3	2.48	0.49
1:C:461:PHE:O	1:C:499:LYS:NZ	2.38	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:141:ILE:CD1	1:D:355:PRO:CB	2.86	0.49
1:D:258:ASP:HB3	1:D:264:ILE:HD11	1.94	0.49
1:A:533:MET:HE1	1:A:534:TYR:CE1	2.47	0.49
1:D:321:HIS:NE2	1:D:346:GLN:OE1	2.43	0.49
1:A:31:GLN:HA	1:A:33:TRP:CZ3	2.48	0.48
1:B:178:VAL:HG22	1:B:182:HIS:CB	2.43	0.48
2:A:701:CBW:C19	2:A:701:CBW:C25	2.85	0.48
1:B:274:ARG:HG2	1:B:274:ARG:NH1	2.16	0.48
1:D:4:PRO:HB2	1:D:12:LEU:HD13	1.95	0.48
1:A:44:PRO:HB3	1:B:310:ASP:HB3	1.95	0.48
1:B:349:VAL:HG13	1:B:407:VAL:HG11	1.96	0.48
1:C:332:HIS:O	1:C:355:PRO:HA	2.14	0.48
1:C:430:PRO:HG3	1:C:461:PHE:CE2	2.48	0.48
1:C:472:TRP:CE3	1:C:508:ALA:HB1	2.48	0.48
1:B:90:THR:HA	1:B:91:HIS:HA	1.66	0.47
1:C:427:ALA:HB3	1:C:428:PRO:HD3	1.96	0.47
1:D:326:ASN:C	1:D:326:ASN:OD1	2.52	0.47
1:C:81:ARG:HG3	1:C:117:THR:OG1	2.14	0.47
1:C:16:ASP:OD2	1:C:172:SER:OG	2.28	0.47
1:C:191:ASP:HB3	1:C:198:LEU:HD11	1.96	0.47
1:A:469:TYR:OH	1:A:505:ASP:HB3	2.14	0.47
1:A:274:ARG:NH1	1:A:276:VAL:HG22	2.28	0.47
1:C:4:PRO:HB2	1:C:12:LEU:HD13	1.97	0.47
1:D:274:ARG:HA	1:D:286:ASN:OD1	2.13	0.47
1:C:349:VAL:CG1	1:C:407:VAL:HG11	2.44	0.47
1:D:96:TYR:HA	1:D:100:ASN:O	2.15	0.47
1:A:162:ASP:OD2	3:A:702:GCU:O3	2.32	0.47
1:B:64:TRP:HH2	1:B:92:HIS:CD2	2.32	0.47
1:A:573:THR:OG1	1:A:577:LYS:HB2	2.15	0.47
1:B:181:GLN:HG2	1:B:207:THR:HG22	1.97	0.47
1:C:27:ASP:C	1:C:29:ASN:H	2.19	0.46
1:B:1:MET:O	1:B:85:ARG:NH2	2.47	0.46
1:C:386:ARG:HB2	1:C:425:TYR:CE2	2.50	0.46
1:A:53:PHE:HA	1:B:318:GLN:HG3	1.97	0.46
1:A:197:GLY:CA	1:A:239:VAL:HG11	2.38	0.46
1:A:28:ASN:C	1:A:30:THR:H	2.19	0.46
1:C:181:GLN:NE2	1:C:209:GLN:O	2.42	0.46
1:C:181:GLN:HG2	1:C:207:THR:CG2	2.46	0.46
2:C:701:CBW:C19	2:C:701:CBW:C26	2.85	0.46
1:D:32:PRO:HD2	1:D:33:TRP:CE3	2.50	0.46
1:A:467:ASN:O	1:A:468:ARG:NH1	2.39	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:302:ASN:CG	1:D:303:ILE:HG13	2.36	0.46
1:D:331:SER:HA	1:D:332:HIS:HA	1.77	0.46
1:B:406:VAL:O	1:B:440:ARG:NE	2.33	0.46
1:B:465:CYS:HA	1:B:502:VAL:O	2.16	0.46
1:A:89:ALA:HB1	1:A:132:VAL:CG2	2.46	0.45
1:B:141:ILE:HD12	1:B:141:ILE:HA	1.71	0.45
1:B:260:SER:OG	1:B:261:LYS:N	2.49	0.45
1:B:257:ILE:HD13	1:B:263:THR:HA	1.97	0.45
1:D:337:GLU:HG3	4:D:732:HOH:O	2.16	0.45
1:B:564:VAL:O	1:B:568:LYS:NZ	2.50	0.45
1:C:452:TYR:CE1	1:C:492:GLY:HA3	2.50	0.45
1:A:443:THR:OG1	1:A:444:PHE:N	2.50	0.45
2:A:701:CBW:C26	2:A:701:CBW:C19	2.85	0.45
1:C:533:MET:CE	1:C:534:TYR:CD1	3.00	0.45
1:D:426:PHE:O	1:D:430:PRO:HD2	2.17	0.45
1:C:274:ARG:C	1:C:274:ARG:HD2	2.36	0.45
1:D:208:THR:O	1:D:210:GLY:O	2.34	0.45
1:B:181:GLN:HG2	1:B:207:THR:HG21	1.99	0.45
1:C:32:PRO:HG2	1:C:66:TYR:CD2	2.52	0.45
1:D:19:TRP:CE2	1:D:45:VAL:HG21	2.52	0.45
1:A:447:VAL:CG2	2:A:701:CBW:H5	2.47	0.45
1:B:244:PRO:O	1:B:245:GLY:C	2.54	0.45
1:D:521:PRO:O	1:D:522:TRP:CB	2.64	0.45
1:D:509:ASP:HB3	1:D:522:TRP:CH2	2.52	0.45
1:B:94:ARG:HB2	1:B:131:ALA:HB3	1.99	0.44
1:C:162:ASP:CB	3:C:702:GCU:H3	2.41	0.44
1:C:90:THR:HA	1:C:91:HIS:HA	1.68	0.44
1:A:25:SER:OG	1:A:26:ASP:N	2.50	0.44
1:A:162:ASP:CG	3:A:702:GCU:H3	2.37	0.44
1:A:90:THR:HA	1:A:91:HIS:HA	1.73	0.44
1:A:4:PRO:HB2	1:A:12:LEU:HD13	1.99	0.44
1:C:321:HIS:NE2	1:C:346:GLN:OE1	2.41	0.44
1:B:45:VAL:O	1:B:169:LEU:O	2.36	0.44
1:C:414:ASP:OD2	2:C:701:CBW:H242	2.18	0.44
1:C:449:LEU:HD21	2:C:701:CBW:C34	2.44	0.44
1:D:207:THR:OG1	1:D:208:THR:N	2.51	0.44
1:B:573:THR:OG1	1:B:577:LYS:HB2	2.16	0.44
1:B:92:HIS:HB3	1:B:133:ASP:HB3	1.98	0.44
1:A:141:ILE:CG2	1:A:142:PRO:HD2	2.47	0.44
2:C:701:CBW:H11C	2:C:701:CBW:O11	2.18	0.44
1:A:32:PRO:HG2	1:A:66:TYR:CE2	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:GLN:NE2	1:B:72:ILE:O	2.50	0.43
1:B:577:LYS:HG3	1:C:515:HIS:O	2.18	0.43
1:C:414:ASP:OD2	2:C:701:CBW:C24	2.66	0.43
1:D:470:PHE:HD2	1:D:474:THR:CG2	2.31	0.43
1:A:449:LEU:HB2	2:A:701:CBW:H271	2.00	0.43
1:A:456:ARG:O	1:A:457:ILE:HD13	2.18	0.43
1:C:178:VAL:HG11	1:C:182:HIS:CD2	2.52	0.43
1:C:550:TRP:NE1	3:C:702:GCU:O4	2.36	0.43
1:D:335:TYR:O	1:D:399:ARG:NH2	2.38	0.43
1:B:509:ASP:HB3	1:B:522:TRP:CZ3	2.53	0.43
1:A:21:PHE:HA	1:A:66:TYR:O	2.19	0.43
1:D:292:PHE:HB2	1:D:546:GLY:HA3	2.01	0.43
1:A:33:TRP:O	1:A:127:ARG:NH1	2.51	0.43
1:B:427:ALA:HB3	1:B:428:PRO:HD3	2.01	0.43
1:C:294:GLY:HA3	1:C:327:SER:O	2.18	0.43
2:A:701:CBW:H11C	2:A:701:CBW:O11	2.18	0.43
1:C:474:THR:OG1	1:C:475:GLN:OE1	2.30	0.43
1:D:427:ALA:HB3	1:D:428:PRO:HD3	2.00	0.43
1:B:381:ILE:CD1	1:B:381:ILE:N	2.82	0.43
1:A:326:ASN:C	1:A:326:ASN:OD1	2.57	0.42
2:A:701:CBW:H193	2:A:701:CBW:O11	2.19	0.42
1:C:53:PHE:HA	1:D:318:GLN:HG3	2.00	0.42
2:C:701:CBW:H193	2:C:701:CBW:O11	2.19	0.42
1:A:274:ARG:NH1	1:A:276:VAL:CG2	2.82	0.42
1:A:426:PHE:O	1:A:430:PRO:HD2	2.20	0.42
1:A:43:CYS:HB2	1:A:53:PHE:HZ	1.85	0.42
1:A:474:THR:HG1	1:A:475:GLN:CD	2.21	0.42
1:A:236:ILE:O	1:A:239:VAL:HG13	2.20	0.42
1:A:89:ALA:HB1	1:A:132:VAL:HG22	2.00	0.42
1:A:45:VAL:HG13	1:A:46:PRO:N	2.34	0.42
1:A:81:ARG:HG3	1:A:117:THR:OG1	2.20	0.42
1:B:326:ASN:C	1:B:326:ASN:OD1	2.58	0.42
1:A:159:TYR:CD2	1:A:161:HIS:CE1	3.07	0.42
1:A:73:VAL:O	1:A:123:GLY:N	2.48	0.42
1:C:217:ILE:HD12	1:C:253:HIS:CE1	2.55	0.42
1:D:78:SER:O	1:D:79:GLU:CB	2.68	0.42
1:B:64:TRP:CZ2	1:B:92:HIS:HD2	2.37	0.42
1:A:197:GLY:H	1:A:239:VAL:CG1	2.28	0.41
1:A:469:TYR:CE1	2:A:701:CBW:H21C	2.54	0.41
1:B:300:ASP:OD2	1:B:555:PHE:HA	2.19	0.41
1:D:97:VAL:HG21	1:D:116:ILE:HG21	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:90:THR:HA	1:D:91:HIS:HA	1.64	0.41
1:B:274:ARG:HA	1:B:286:ASN:OD1	2.21	0.41
1:C:68:GLN:NE2	1:C:127:ARG:HD2	2.34	0.41
1:C:141:ILE:HA	1:C:141:ILE:HD12	1.56	0.41
1:C:162:ASP:HB2	3:C:702:GCU:O3	2.19	0.41
1:C:357:VAL:HG22	1:C:358:GLY:N	2.34	0.41
1:A:92:HIS:O	1:A:132:VAL:HA	2.20	0.41
1:A:413:ASN:HD22	1:A:445:ALA:HB3	1.85	0.41
1:B:231:ASN:OD1	1:B:231:ASN:C	2.57	0.41
1:A:345:ARG:NH1	1:B:42:GLU:OE2	2.53	0.41
1:C:197:GLY:N	1:C:239:VAL:HG23	2.35	0.41
1:C:283:PHE:HZ	1:C:407:VAL:HG22	1.85	0.41
2:A:701:CBW:H152	2:A:701:CBW:H262	1.79	0.41
1:D:577:LYS:HD3	1:D:577:LYS:HA	1.77	0.41
1:A:38:LYS:O	1:A:39:THR:HG23	2.20	0.41
1:A:473:TYR:HH	3:A:702:GCU:C6	2.34	0.41
1:A:345:ARG:HD2	1:B:42:GLU:OE1	2.21	0.41
1:A:569:LYS:NZ	3:A:702:GCU:O6A	2.52	0.41
1:A:1:MET:O	1:A:85:ARG:NH2	2.46	0.41
1:A:68:GLN:NE2	1:A:127:ARG:HD2	2.36	0.41
1:B:487:GLU:HG2	1:B:491:ARG:NH2	2.36	0.41
1:B:521:PRO:O	1:B:522:TRP:CB	2.64	0.41
1:D:465:CYS:HA	1:D:502:VAL:O	2.20	0.41
2:C:701:CBW:H262	2:C:701:CBW:H152	1.80	0.41
1:D:141:ILE:HG23	1:D:388:ALA:HB1	2.02	0.41
1:D:349:VAL:CG1	1:D:407:VAL:HG11	2.40	0.41
1:D:564:VAL:O	1:D:568:LYS:NZ	2.53	0.41
1:A:521:PRO:HG2	1:C:559:VAL:HG13	2.02	0.40
1:C:380:ARG:O	1:C:381:ILE:C	2.58	0.40
1:A:399:ARG:NH1	1:A:400:ASP:OD2	2.54	0.40
1:C:141:ILE:CG2	1:C:142:PRO:CD	2.90	0.40
1:D:141:ILE:HD12	1:D:142:PRO:CD	2.44	0.40
1:D:256:ILE:HG22	1:D:264:ILE:HD12	2.02	0.40
1:B:329:ARG:NH1	1:B:504:THR:HB	2.37	0.40
1:C:1:MET:HB2	1:C:113:GLU:OE2	2.21	0.40
1:D:236:ILE:HA	1:D:237:PRO:HD2	1.87	0.40
1:C:181:GLN:HG2	1:C:207:THR:HG22	2.03	0.40
1:D:361:PHE:CE1	1:D:381:ILE:HD13	2.48	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	579/637 (91%)	547 (94%)	28 (5%)	4 (1%)	22	41
1	B	578/637 (91%)	548 (95%)	28 (5%)	2 (0%)	41	62
1	C	579/637 (91%)	538 (93%)	35 (6%)	6 (1%)	15	30
1	D	578/637 (91%)	548 (95%)	28 (5%)	2 (0%)	41	62
All	All	2314/2548 (91%)	2181 (94%)	119 (5%)	14 (1%)	25	45

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	303	ILE
1	B	39	THR
1	C	332	HIS
1	D	209	GLN
1	A	29	ASN
1	A	331	SER
1	A	332	HIS
1	B	260	SER
1	C	28	ASN
1	C	331	SER
1	C	30	THR
1	C	381	ILE
1	C	46	PRO
1	A	244	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	493/531 (93%)	490 (99%)	3 (1%)	86	94
1	B	492/531 (93%)	487 (99%)	5 (1%)	76	89
1	C	493/531 (93%)	489 (99%)	4 (1%)	81	92
1	D	492/531 (93%)	489 (99%)	3 (1%)	86	94
All	All	1970/2124 (93%)	1955 (99%)	15 (1%)	81	92

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	198	LEU
1	A	200	ASP
1	A	222	THR
1	B	52	ILE
1	B	162	ASP
1	B	274	ARG
1	B	399	ARG
1	B	571	VAL
1	C	39	THR
1	C	231	ASN
1	C	257	ILE
1	C	274	ARG
1	D	162	ASP
1	D	274	ARG
1	D	495	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	181	GLN
1	A	318	GLN
1	A	413	ASN
1	B	92	HIS
1	C	29	ASN
1	C	332	HIS
1	D	160	GLN
1	D	181	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 monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

4 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	GCU	C	702	2	9,12,13	0.86	0	12,17,19	0.75	0
2	CBW	A	701	3	35,38,38	2.95	14 (40%)	61,67,67	2.26	14 (22%)
2	CBW	C	701	3	35,38,38	2.97	14 (40%)	61,67,67	2.37	16 (26%)
3	GCU	A	702	2	9,12,13	0.84	0	12,17,19	0.76	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	GCU	C	702	2	-	0/0/21/24	0/1/1/1
2	CBW	A	701	3	-	0/0/100/100	0/5/5/5
2	CBW	C	701	3	-	0/0/100/100	0/5/5/5
3	GCU	A	702	2	-	0/0/21/24	0/1/1/1

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	CBW	C9-C11	-8.89	1.41	1.52
2	C	701	CBW	C9-C11	-8.88	1.41	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	701	CBW	C7-C8	-8.31	1.39	1.54
2	A	701	CBW	C7-C8	-8.27	1.39	1.54
2	C	701	CBW	C15-C14	-5.14	1.46	1.54
2	A	701	CBW	C15-C14	-5.02	1.47	1.54
2	A	701	CBW	C1-C10	-5.00	1.45	1.54
2	C	701	CBW	C1-C10	-4.95	1.45	1.54
2	A	701	CBW	C7-C6	-4.68	1.43	1.53
2	C	701	CBW	C7-C6	-4.67	1.43	1.53
2	A	701	CBW	C8-C14	3.85	1.65	1.58
2	C	701	CBW	C8-C14	3.85	1.65	1.58
2	A	701	CBW	C8-C9	-3.83	1.52	1.56
2	C	701	CBW	C8-C9	-3.79	1.52	1.56
2	A	701	CBW	C10-C9	3.67	1.62	1.56
2	C	701	CBW	C10-C9	3.63	1.62	1.56
2	C	701	CBW	C22-C21	-2.45	1.51	1.54
2	C	701	CBW	C17-C18	-2.44	1.51	1.55
2	A	701	CBW	C17-C18	-2.35	1.51	1.55
2	A	701	CBW	C14-C13	-2.18	1.49	1.53
2	C	701	CBW	C14-C13	-2.14	1.49	1.53
2	A	701	CBW	C22-C21	-2.11	1.51	1.54
2	A	701	CBW	C21-C33	2.07	1.54	1.50
2	C	701	CBW	C16-C15	2.05	1.59	1.53
2	A	701	CBW	C16-C15	2.02	1.59	1.53
2	C	701	CBW	C21-C33	2.02	1.54	1.50
2	A	701	CBW	C10-C5	-2.01	1.53	1.56
2	C	701	CBW	C10-C5	-2.00	1.53	1.56

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	701	CBW	C8-C9-C10	-10.03	110.78	118.10
2	A	701	CBW	C8-C9-C10	-10.03	110.78	118.10
2	C	701	CBW	C34-C21-C33	-6.21	98.70	109.22
2	A	701	CBW	C34-C21-C33	-5.66	99.62	109.22
2	C	701	CBW	C4-C5-C10	-5.33	111.92	117.17
2	A	701	CBW	C4-C5-C10	-5.32	111.92	117.17
2	C	701	CBW	C20-C18-C17	-4.01	109.95	113.13
2	C	701	CBW	C21-C20-C18	3.89	119.48	113.62
2	A	701	CBW	C20-C18-C17	-3.46	110.39	113.13
2	A	701	CBW	C15-C16-C17	-3.41	106.30	113.88
2	C	701	CBW	C6-C7-C8	3.39	118.68	112.84
2	A	701	CBW	C6-C7-C8	3.38	118.66	112.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	701	CBW	C15-C16-C17	-3.38	106.35	113.88
2	A	701	CBW	C22-C21-C20	-3.27	102.67	109.21
2	C	701	CBW	C27-C14-C8	-3.21	109.33	112.33
2	A	701	CBW	C27-C14-C8	-3.18	109.36	112.33
2	C	701	CBW	C22-C21-C20	-2.81	103.60	109.21
2	C	701	CBW	C1-C10-C9	2.73	111.15	108.19
2	A	701	CBW	C1-C10-C9	2.73	111.15	108.19
2	A	701	CBW	C10-C9-C11	-2.59	113.27	115.50
2	C	701	CBW	C10-C9-C11	-2.58	113.28	115.50
2	C	701	CBW	C2-C3-C4	-2.41	110.84	113.32
2	A	701	CBW	C2-C3-C4	-2.41	110.85	113.32
2	A	701	CBW	C27-C14-C13	2.31	109.36	106.95
2	C	701	CBW	C27-C14-C13	2.26	109.31	106.95
2	A	701	CBW	C1-C2-C3	-2.25	108.11	111.51
2	C	701	CBW	C1-C2-C3	-2.24	108.13	111.51
2	C	701	CBW	C22-C23-C17	2.05	118.44	113.88
2	C	701	CBW	O11-C11-C9	-2.03	120.59	123.19
2	A	701	CBW	O11-C11-C9	-2.02	120.60	123.19

There are no chirality outliers.

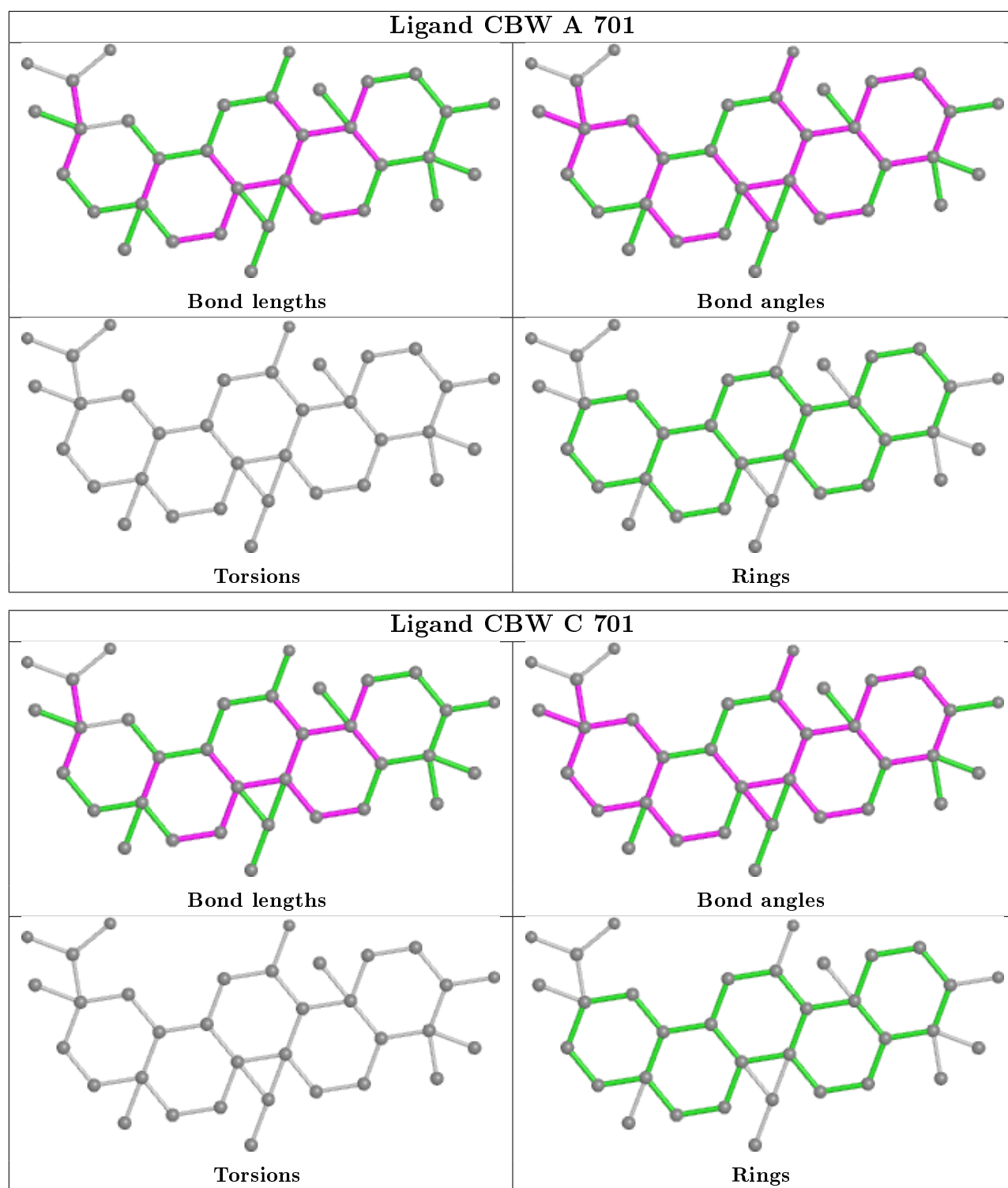
There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 59 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	702	GCU	10	0
2	A	701	CBW	19	0
2	C	701	CBW	25	0
3	A	702	GCU	11	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.



## 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<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	583/637 (91%)	-0.61	9 (1%) 73 70	4, 11, 29, 105	0
1	B	582/637 (91%)	-0.59	13 (2%) 62 57	4, 13, 32, 95	0
1	C	583/637 (91%)	-0.51	9 (1%) 73 70	3, 12, 32, 99	0
1	D	582/637 (91%)	-0.64	6 (1%) 82 80	3, 11, 28, 64	0
All	All	2330/2548 (91%)	-0.59	37 (1%) 72 68	3, 12, 31, 105	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	28	ASN	11.5
1	B	28	ASN	10.8
1	A	28	ASN	10.1
1	A	593	HIS	7.3
1	A	29	ASN	6.3
1	C	593	HIS	5.7
1	B	29	ASN	5.2
1	C	208	THR	4.9
1	C	29	ASN	4.3
1	D	208	THR	4.0
1	B	261	LYS	3.7
1	B	38	LYS	3.6
1	B	27	ASP	3.6
1	B	208	THR	3.4
1	D	261	LYS	3.4
1	B	30	THR	3.3
1	C	592	LEU	3.2
1	C	26	ASP	3.2
1	D	141	ILE	3.0
1	A	38	LYS	3.0
1	B	259	SER	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	208	THR	2.9
1	A	592	LEU	2.8
1	C	27	ASP	2.8
1	A	362	SER	2.5
1	D	475	GLN	2.5
1	B	26	ASP	2.5
1	A	30	THR	2.4
1	B	592	LEU	2.3
1	B	245	GLY	2.2
1	C	373	PRO	2.2
1	B	262	LYS	2.2
1	D	260	SER	2.2
1	C	259	SER	2.1
1	A	262	LYS	2.1
1	B	209	GLN	2.0
1	D	27	ASP	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 monosaccharides 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<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.

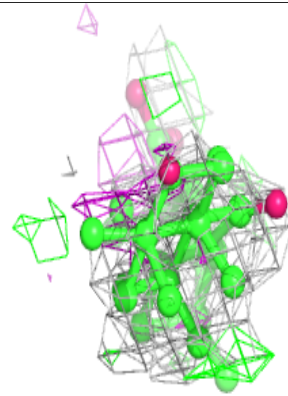
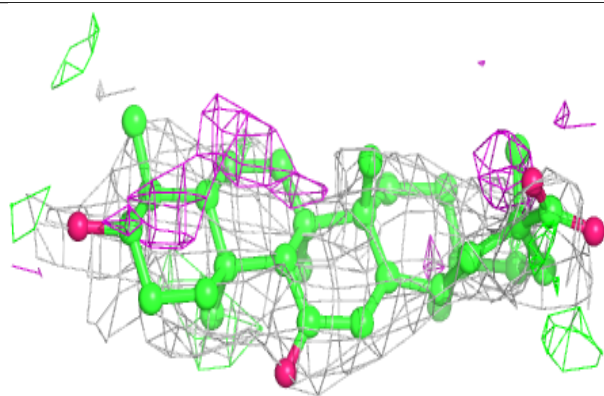
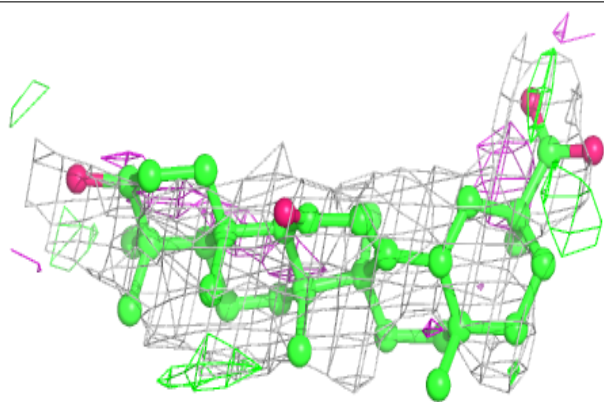
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	CBW	C	701	34/34	0.66	0.40	45,70,86,87	0
3	GCU	C	702	12/13	0.72	0.39	28,42,56,56	0
2	CBW	A	701	34/34	0.73	0.35	49,72,84,85	0
3	GCU	A	702	12/13	0.76	0.41	22,42,70,78	0

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.

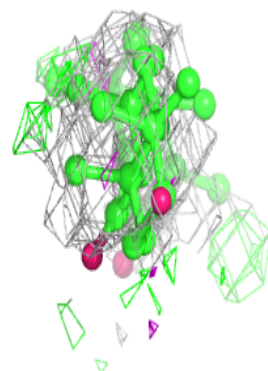
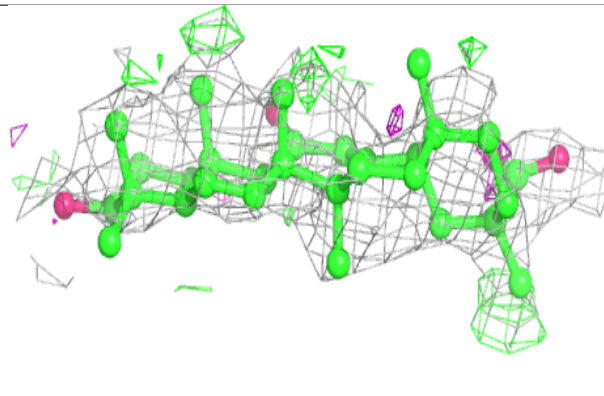
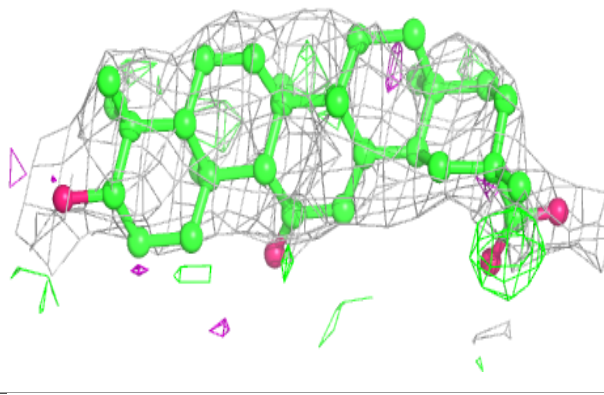
**Electron density around CBW C 701:**

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



**Electron density around CBW A 701:**

$2mF_o-DF_c$  (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.