



# Full wwPDB X-ray Structure Validation Report i

May 25, 2020 – 08:53 am BST

PDB ID : 3QWT  
Title : The crystal structure of a possible member of GH105 family from *Salmonella enterica* subsp. *enterica* serovar Paratyphi A str. ATCC 9150  
Authors : Tan, K.; Hatzos-Skitges, C.; Bearden, J.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)  
Deposited on : 2011-02-28  
Resolution : 2.18 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 i symbol.

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The following versions of software and data (see [references](#) ①) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.1.3  
EDS : 2.11  
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

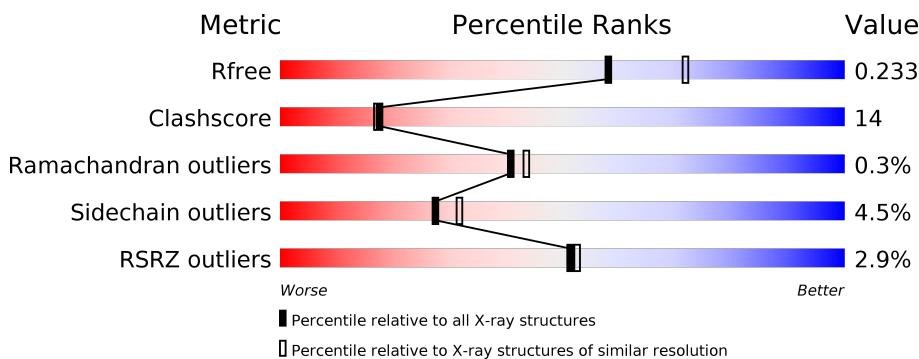
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## X-RAY DIFFRACTION

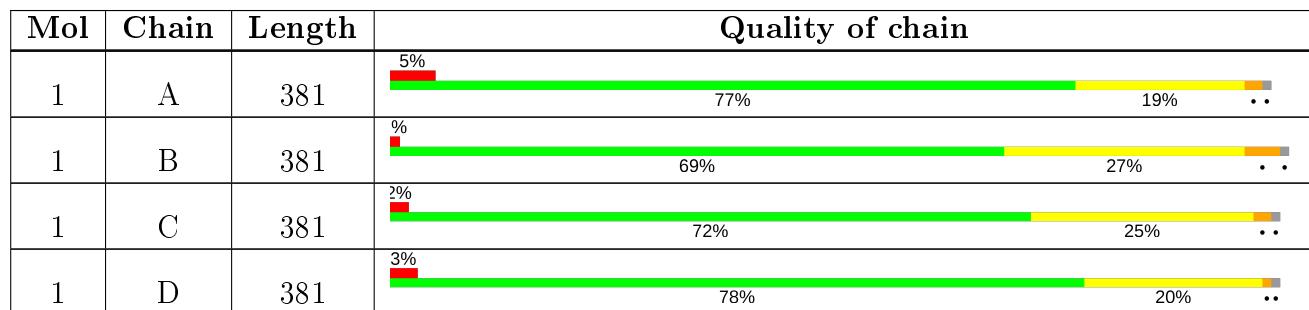
The reported resolution of this entry is 2.18 Å.

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	6864 (2.20-2.16)
Clashscore	141614	7689 (2.20-2.16)
Ramachandran outliers	138981	7564 (2.20-2.16)
Sidechain outliers	138945	7564 (2.20-2.16)
RSRZ outliers	127900	6738 (2.20-2.16)

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.



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	GOL	D	381	-	-	X	-

## 2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 12818 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.

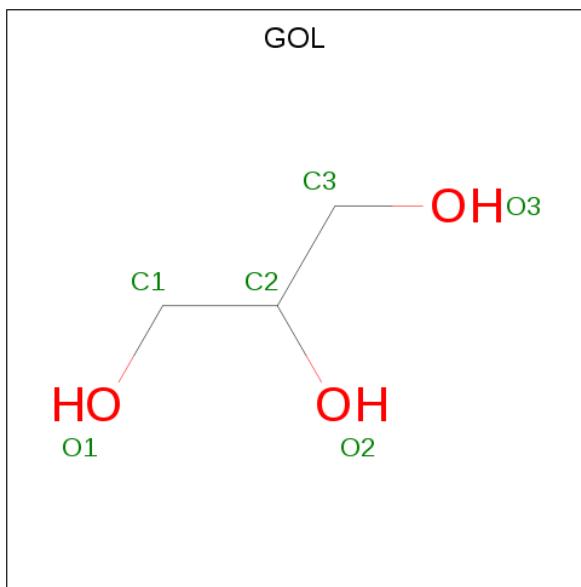
- Molecule 1 is a protein called Putative GH105 family protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	378	Total	C	N	O	S	Se	0	0	0
			3099	1989	534	559	2	15			
1	B	378	Total	C	N	O	S	Se	0	0	0
			3099	1989	534	559	2	15			
1	C	378	Total	C	N	O	S	Se	0	0	0
			3099	1989	534	559	2	15			
1	D	378	Total	C	N	O	S	Se	0	0	0
			3099	1989	534	559	2	15			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP Q5PMY7
A	-1	ASN	-	EXPRESSION TAG	UNP Q5PMY7
A	0	ALA	-	EXPRESSION TAG	UNP Q5PMY7
B	-2	SER	-	EXPRESSION TAG	UNP Q5PMY7
B	-1	ASN	-	EXPRESSION TAG	UNP Q5PMY7
B	0	ALA	-	EXPRESSION TAG	UNP Q5PMY7
C	-2	SER	-	EXPRESSION TAG	UNP Q5PMY7
C	-1	ASN	-	EXPRESSION TAG	UNP Q5PMY7
C	0	ALA	-	EXPRESSION TAG	UNP Q5PMY7
D	-2	SER	-	EXPRESSION TAG	UNP Q5PMY7
D	-1	ASN	-	EXPRESSION TAG	UNP Q5PMY7
D	0	ALA	-	EXPRESSION TAG	UNP Q5PMY7

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	B	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	90	Total O 90 90	0	0

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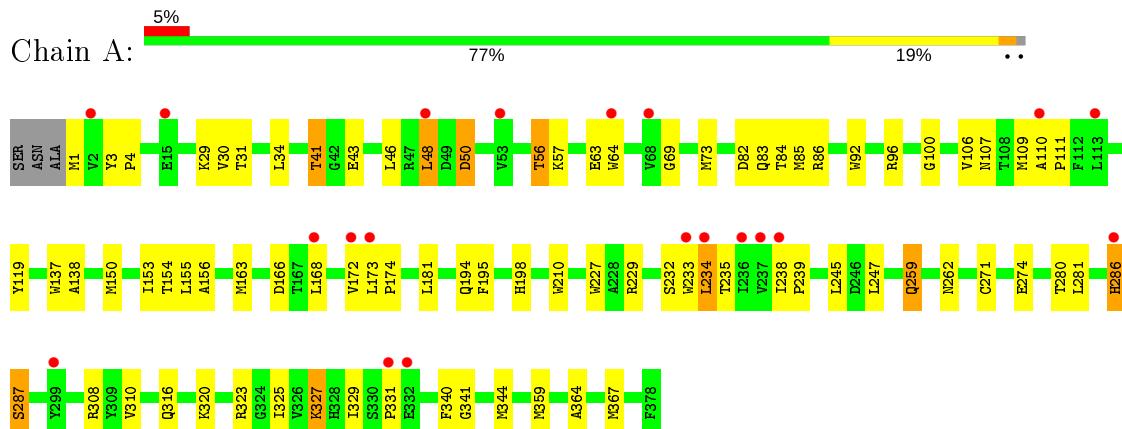
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	107	Total O 107 107	0	0
3	C	54	Total O 54 54	0	0
3	D	105	Total O 105 105	0	0

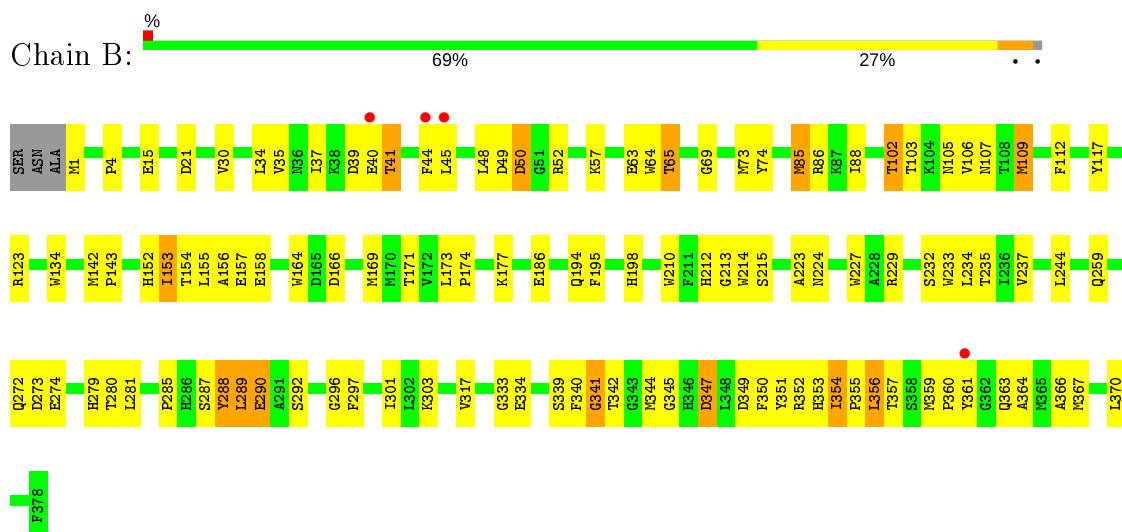
### 3 Residue-property plots

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.

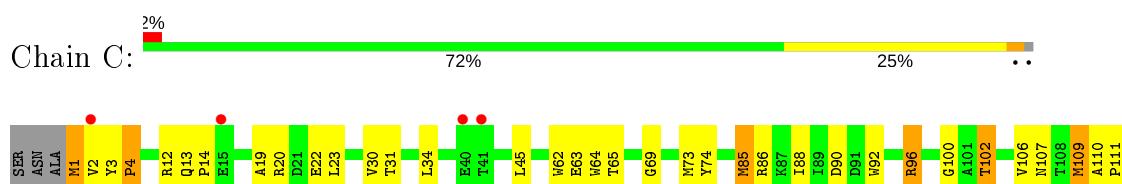
- Molecule 1: Putative GH105 family protein

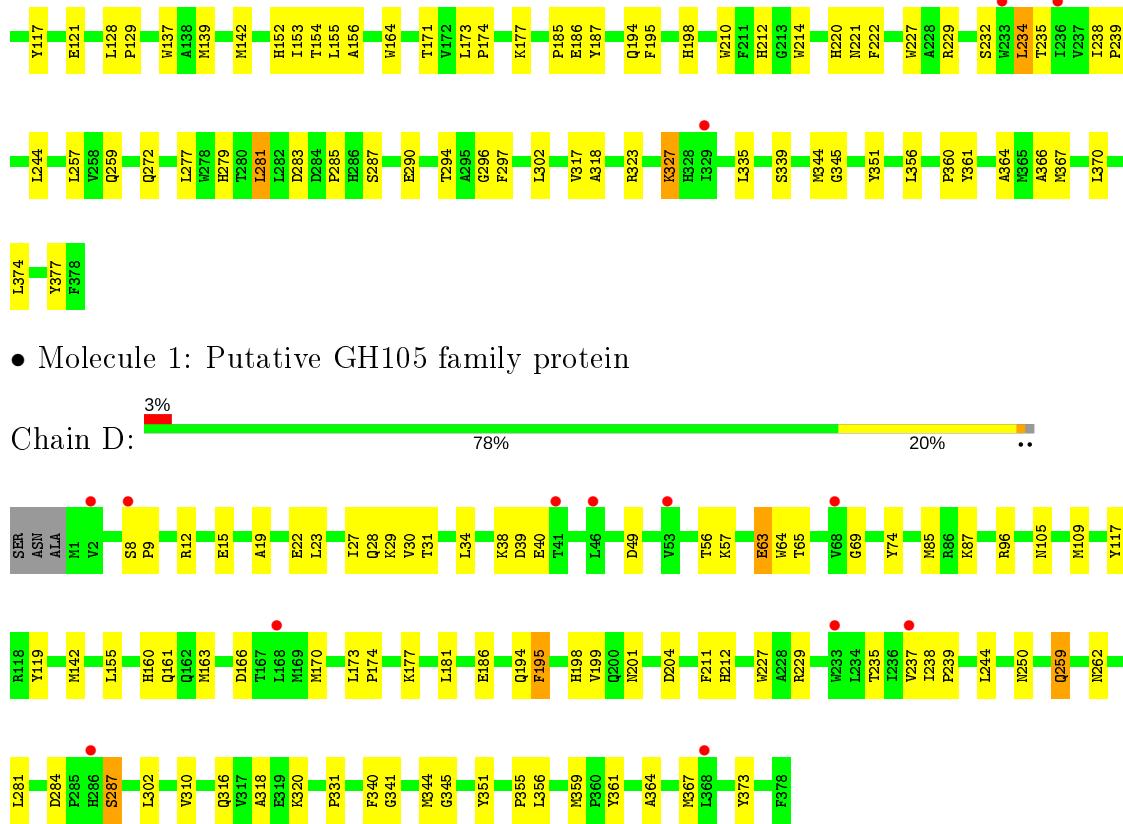


- Molecule 1: Putative GH105 family protein



- Molecule 1: Putative GH105 family protein





## 4 Data and refinement statistics i

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	121.03 Å    159.36 Å    179.12 Å 90.00°    90.00°    90.00°	Depositor
Resolution (Å)	39.84 – 2.18 39.84 – 2.18	Depositor EDS
% Data completeness (in resolution range)	94.5 (39.84-2.18) 97.7 (39.84-2.18)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle^1$	2.12 (at 2.18 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
$R$ , $R_{free}$	0.182 , 0.237 0.180 , 0.233	Depositor DCC
$R_{free}$ test set	4402 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.5	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 44.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	12818	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	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: GOL

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	0.42	0/3176	0.56	0/4294
1	B	0.43	0/3176	0.57	0/4294
1	C	0.37	0/3176	0.54	0/4294
1	D	0.45	0/3176	0.57	0/4294
All	All	0.42	0/12704	0.56	0/17176

There are no bond length outliers.

There are no bond angle outliers.

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	3099	0	2990	54	0
1	B	3099	0	2990	116	0
1	C	3099	0	2990	95	0
1	D	3099	0	2990	73	0
2	A	18	0	24	4	0
2	B	6	0	8	0	0
2	C	18	0	24	1	0
2	D	24	0	32	6	0
3	A	90	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	107	0	0	2	0
3	C	54	0	0	2	0
3	D	105	0	0	5	0
All	All	12818	0	12048	335	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 14.

All (335) 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:D:31:THR:HG21	1:D:85:MSE:HE1	1.38	1.01
1:A:154:THR:HG22	1:A:156:ALA:H	1.25	1.01
1:D:56:THR:HG23	1:D:359:MSE:HE1	1.47	0.96
1:B:354:ILE:HD13	1:B:355:PRO:HD2	1.52	0.90
1:B:339:SER:HB3	1:B:354:ILE:HG21	1.54	0.89
1:A:63:GLU:HG2	1:A:64:TRP:H	1.37	0.89
1:B:360:PRO:O	1:B:364:ALA:HB2	1.75	0.87
1:A:69:GLY:HA2	1:A:364:ALA:HB2	1.60	0.84
1:D:344:MSE:HE1	3:D:410:HOH:O	1.78	0.81
1:D:344:MSE:CE	3:D:410:HOH:O	2.27	0.81
1:D:28:GLN:HG3	2:D:381:GOL:H31	1.62	0.81
1:B:21:ASP:HB2	3:B:450:HOH:O	1.79	0.80
1:C:85:MSE:HA	1:C:85:MSE:CE	2.12	0.80
1:D:344:MSE:HE3	1:D:345:GLY:N	1.99	0.78
1:B:65:THR:OG1	1:B:360:PRO:HB2	1.84	0.77
1:A:31:THR:OG1	1:A:85:MSE:HE1	1.85	0.77
1:B:177:LYS:HA	1:B:244:LEU:HD21	1.68	0.76
1:C:85:MSE:HA	1:C:85:MSE:HE3	1.68	0.75
1:B:229:ARG:NH2	1:B:339:SER:HB2	2.01	0.75
1:B:37:ILE:HG22	1:B:57:LYS:HG2	1.68	0.74
1:C:65:THR:HG21	1:C:360:PRO:HD2	1.68	0.73
1:B:195:PHE:HZ	1:B:237:VAL:HG12	1.53	0.73
1:D:142:MSE:O	2:D:382:GOL:H11	1.87	0.73
1:B:360:PRO:O	1:B:364:ALA:CB	2.38	0.71
1:D:344:MSE:HE3	1:D:345:GLY:H	1.56	0.70
1:A:229:ARG:O	1:A:232:SER:HB3	1.92	0.70
1:A:56:THR:CG2	1:A:359:MSE:HE1	2.20	0.70
1:B:64:TRP:HB3	1:B:107:ASN:HB3	1.72	0.70
1:D:69:GLY:HA2	1:D:364:ALA:HB2	1.74	0.70
1:B:357:THR:HG23	1:B:359:MSE:HG2	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:106:VAL:HG23	1:C:153:ILE:O	1.91	0.69
1:C:45:LEU:HD12	3:C:423:HOH:O	1.91	0.69
1:C:194:GLN:O	1:C:198:HIS:HD2	1.75	0.69
1:B:106:VAL:HG21	1:B:152:HIS:HA	1.76	0.67
1:B:64:TRP:CH2	1:B:361:TYR:CD2	2.82	0.67
1:D:63:GLU:OE1	1:D:63:GLU:HA	1.94	0.67
1:A:107:ASN:OD1	2:A:380:GOL:H32	1.94	0.67
1:C:277:LEU:HD22	1:C:294:THR:HG21	1.76	0.67
1:B:15:GLU:O	1:B:15:GLU:HG3	1.93	0.67
1:C:109:MSE:O	1:C:109:MSE:HE3	1.95	0.67
1:B:85:MSE:HE3	1:B:88:ILE:HD12	1.78	0.66
1:B:194:GLN:O	1:B:198:HIS:HD2	1.79	0.66
1:D:56:THR:HG22	3:D:482:HOH:O	1.95	0.65
1:B:143:PRO:HD2	1:B:153:ILE:HD11	1.78	0.65
1:D:344:MSE:CE	1:D:351:TYR:OH	2.45	0.65
1:B:109:MSE:O	1:B:109:MSE:HE3	1.97	0.64
1:D:356:LEU:N	1:D:356:LEU:HD12	2.12	0.64
1:B:288:TYR:OH	1:B:354:ILE:HG22	1.99	0.63
1:C:154:THR:HG22	1:C:155:LEU:N	2.13	0.63
1:B:37:ILE:CG2	1:B:57:LYS:CG	2.77	0.63
1:A:34:LEU:HD13	1:A:73:MSE:HE3	1.81	0.62
1:A:56:THR:HG22	1:A:359:MSE:HE1	1.80	0.62
1:C:164:TRP:CZ2	1:C:344:MSE:HB2	2.34	0.62
1:C:117:TYR:O	1:C:121:GLU:HG3	1.99	0.62
1:A:323:ARG:O	1:A:327:LYS:HE2	2.00	0.61
1:A:100:GLY:O	2:A:379:GOL:H32	2.00	0.61
1:B:37:ILE:HG22	1:B:57:LYS:CG	2.31	0.61
1:D:177:LYS:HA	1:D:244:LEU:HD21	1.82	0.61
1:B:279:HIS:HB3	1:B:287:SER:HB3	1.82	0.60
1:B:344:MSE:HE3	1:B:345:GLY:H	1.64	0.60
1:D:65:THR:HG22	1:D:361:TYR:HD2	1.66	0.60
1:A:69:GLY:HA2	1:A:364:ALA:CB	2.32	0.60
1:C:85:MSE:HE3	1:C:88:ILE:HD12	1.83	0.60
1:B:212:HIS:HB3	1:B:344:MSE:HE1	1.83	0.60
1:A:63:GLU:HG2	1:A:64:TRP:N	2.11	0.60
1:B:85:MSE:HA	1:B:85:MSE:CE	2.31	0.60
1:D:63:GLU:OE2	1:D:64:TRP:HD1	1.85	0.60
1:C:121:GLU:HB3	3:C:405:HOH:O	2.02	0.59
1:D:19:ALA:HB3	1:D:22:GLU:HG3	1.83	0.59
1:D:344:MSE:HE1	1:D:351:TYR:OH	2.01	0.59
1:D:30:VAL:HG12	1:D:367:MSE:HE2	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:339:SER:HB3	1:B:354:ILE:CG2	2.30	0.59
1:C:1:MSE:HG3	1:C:1:MSE:O	2.03	0.59
1:B:154:THR:HG23	1:B:156:ALA:H	1.68	0.59
1:B:102:THR:HG22	1:B:103:THR:H	1.67	0.58
1:A:150:MSE:HG3	1:A:194:GLN:CD	2.23	0.58
1:C:366:ALA:O	1:C:370:LEU:HG	2.04	0.58
1:C:186:GLU:H	1:C:186:GLU:CD	2.07	0.58
1:B:40:GLU:H	1:B:40:GLU:CD	2.06	0.58
1:B:69:GLY:O	1:B:73:MSE:HG3	2.04	0.58
1:D:57:LYS:HG3	1:D:359:MSE:SE	2.53	0.58
1:D:170:MSE:O	1:D:174:PRO:HG2	2.03	0.57
1:C:73:MSE:SE	1:C:85:MSE:HE1	2.54	0.57
1:B:297:PHE:O	1:B:301:ILE:HG13	2.03	0.57
1:B:34:LEU:HD12	1:B:363:GLN:HB2	1.85	0.57
1:C:364:ALA:HA	1:C:367:MSE:HG2	1.85	0.57
1:B:30:VAL:HG12	1:B:367:MSE:HE2	1.87	0.57
1:C:106:VAL:HG21	1:C:152:HIS:HA	1.87	0.56
1:D:109:MSE:HE3	1:D:174:PRO:HB2	1.87	0.56
1:D:28:GLN:HB2	2:D:381:GOL:H2	1.87	0.56
1:C:154:THR:CG2	1:C:155:LEU:N	2.68	0.56
1:A:41:THR:HB	1:A:43:GLU:H	1.69	0.56
1:B:344:MSE:HE3	1:B:351:TYR:OH	2.05	0.56
1:D:65:THR:HG22	1:D:361:TYR:CD2	2.41	0.56
1:A:154:THR:HG23	3:A:460:HOH:O	2.05	0.56
1:C:229:ARG:O	1:C:232:SER:HB3	2.05	0.56
1:B:361:TYR:HA	1:B:364:ALA:HB3	1.87	0.55
1:B:349:ASP:HA	1:B:352:ARG:CD	2.36	0.55
1:B:64:TRP:HB3	1:B:107:ASN:CB	2.36	0.55
1:C:19:ALA:HB3	1:C:22:GLU:HG3	1.89	0.55
1:C:65:THR:CG2	1:C:360:PRO:HD2	2.37	0.54
1:B:212:HIS:CD2	1:B:351:TYR:OH	2.60	0.54
1:C:86:ARG:HD2	1:C:90:ASP:OD2	2.07	0.54
1:B:4:PRO:HA	1:D:49:ASP:OD1	2.07	0.54
1:B:223:ALA:O	1:B:224:ASN:HB2	2.06	0.54
1:D:238:ILE:HB	1:D:239:PRO:HD3	1.88	0.54
1:A:364:ALA:HA	1:A:367:MSE:HE3	1.90	0.54
1:B:212:HIS:HD2	1:B:351:TYR:OH	1.91	0.54
1:A:154:THR:HG22	1:A:156:ALA:N	2.08	0.53
1:B:37:ILE:CG2	1:B:57:LYS:HG3	2.38	0.53
1:C:63:GLU:HG3	1:C:64:TRP:N	2.23	0.53
1:B:280:THR:HG22	1:B:287:SER:OG	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:273:ASP:OD1	1:B:274:GLU:N	2.41	0.53
1:B:37:ILE:HG21	1:B:57:LYS:HG3	1.90	0.53
1:C:154:THR:HG22	1:C:156:ALA:H	1.72	0.53
1:D:105:ASN:HD21	1:D:155:LEU:HG	1.73	0.53
1:B:229:ARG:HH21	1:B:339:SER:HB2	1.74	0.53
1:C:195:PHE:CE2	1:C:257:LEU:CD2	2.92	0.52
1:D:74:TYR:CZ	1:D:117:TYR:HB3	2.44	0.52
1:B:112:PHE:CE2	1:B:134:TRP:CD2	2.97	0.52
1:D:194:GLN:O	1:D:198:HIS:HD2	1.91	0.52
1:C:64:TRP:HB3	1:C:107:ASN:HB3	1.92	0.52
1:D:344:MSE:HE3	1:D:351:TYR:OH	2.09	0.52
1:D:39:ASP:HB2	1:D:57:LYS:HD3	1.92	0.52
1:A:82:ASP:OD2	1:A:84:THR:HB	2.10	0.52
1:C:194:GLN:O	1:C:198:HIS:CD2	2.61	0.52
1:C:74:TYR:CZ	1:C:117:TYR:HB3	2.45	0.52
1:B:169:MSE:HE1	1:B:173:LEU:HD23	1.90	0.52
1:C:106:VAL:HG22	1:C:142:MSE:HE3	1.91	0.51
1:C:63:GLU:HG3	1:C:64:TRP:H	1.75	0.51
1:B:360:PRO:HG3	3:B:472:HOH:O	2.10	0.51
1:D:316:GLN:O	1:D:320:LYS:HG3	2.10	0.51
1:B:152:HIS:CD2	1:B:164:TRP:CD2	2.99	0.51
1:D:119:TYR:HD2	1:D:181:LEU:HD21	1.74	0.51
1:C:31:THR:HA	1:C:367:MSE:HE1	1.91	0.51
1:D:344:MSE:HE1	1:D:351:TYR:HH	1.74	0.51
1:D:34:LEU:C	1:D:34:LEU:HD23	2.31	0.51
1:B:195:PHE:HZ	1:B:237:VAL:CG1	2.23	0.51
1:C:30:VAL:HG12	1:C:367:MSE:HE3	1.92	0.51
1:C:30:VAL:HG12	1:C:367:MSE:CE	2.39	0.51
1:D:109:MSE:CE	1:D:174:PRO:HB2	2.41	0.51
1:C:74:TYR:HE1	1:C:121:GLU:OE1	1.94	0.51
1:A:106:VAL:HG12	1:A:153:ILE:O	2.12	0.50
1:B:177:LYS:CA	1:B:244:LEU:HD21	2.40	0.50
1:B:347:ASP:O	1:B:350:PHE:HB3	2.12	0.50
1:C:222:PHE:HB2	1:C:344:MSE:HE2	1.93	0.50
1:A:280:THR:HG22	1:A:287:SER:HB3	1.94	0.50
1:C:185:PRO:HD2	1:C:186:GLU:OE2	2.11	0.50
1:D:344:MSE:HE2	3:D:410:HOH:O	2.02	0.50
1:D:364:ALA:HA	1:D:367:MSE:HE3	1.94	0.50
1:D:69:GLY:HA2	1:D:364:ALA:CB	2.40	0.50
1:C:69:GLY:HA2	1:C:364:ALA:HB2	1.94	0.49
1:B:44:PHE:CD2	1:B:57:LYS:HE2	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:MSE:HE2	1:B:174:PRO:HB2	1.94	0.49
1:B:73:MSE:SE	1:B:85:MSE:HE1	2.63	0.49
1:D:28:GLN:HB2	2:D:381:GOL:O1	2.12	0.49
1:A:154:THR:CG2	1:A:155:LEU:N	2.75	0.49
1:D:31:THR:HG21	1:D:85:MSE:CE	2.27	0.49
1:A:245:LEU:HB3	1:A:247:LEU:HG	1.95	0.49
1:B:344:MSE:HE3	1:B:345:GLY:N	2.28	0.49
1:D:166:ASP:OD2	1:D:212:HIS:HE1	1.94	0.49
1:B:340:PHE:N	1:B:355:PRO:O	2.38	0.49
1:C:279:HIS:CD2	1:C:285:PRO:HA	2.48	0.49
1:B:210:TRP:CD2	1:B:234:LEU:HD13	2.48	0.48
1:C:212:HIS:HD2	1:C:351:TYR:OH	1.96	0.48
1:B:49:ASP:OD2	1:C:4:PRO:HA	2.14	0.48
1:D:173:LEU:HB3	1:D:174:PRO:HD3	1.95	0.48
1:B:194:GLN:O	1:B:198:HIS:CD2	2.63	0.48
1:D:195:PHE:HZ	1:D:237:VAL:HG12	1.78	0.48
1:B:65:THR:HB	1:B:361:TYR:HD2	1.79	0.48
1:B:39:ASP:C	1:B:41:THR:H	2.17	0.48
1:C:23:LEU:HD23	1:C:374:LEU:HD23	1.94	0.48
1:A:227:TRP:CZ2	1:A:229:ARG:HB3	2.49	0.48
1:B:213:GLY:HA3	1:B:344:MSE:SE	2.64	0.48
1:B:153:ILE:HG23	1:B:158:GLU:HA	1.94	0.48
1:B:105:ASN:HD21	1:B:155:LEU:HG	1.78	0.48
1:C:20:ARG:HB2	1:C:377:TYR:CD2	2.48	0.48
1:C:335:LEU:HD22	1:C:366:ALA:HB2	1.96	0.48
1:B:154:THR:CG2	1:B:157:GLU:H	2.27	0.48
1:C:214:TRP:HE3	1:C:221:ASN:ND2	2.12	0.48
1:C:31:THR:CA	1:C:367:MSE:HE1	2.44	0.48
1:C:344:MSE:HE3	1:C:345:GLY:N	2.29	0.48
1:B:352:ARG:O	1:B:353:HIS:CD2	2.67	0.47
1:D:23:LEU:HD21	1:D:373:TYR:HD1	1.79	0.47
1:C:64:TRP:HB3	1:C:107:ASN:CB	2.44	0.47
1:A:64:TRP:HB3	1:A:107:ASN:HB3	1.96	0.47
1:A:63:GLU:CG	1:A:64:TRP:H	2.16	0.47
1:A:166:ASP:CB	2:A:380:GOL:H31	2.45	0.47
1:B:50:ASP:CG	1:B:52:ARG:HE	2.17	0.47
1:C:238:ILE:HB	1:C:239:PRO:HD3	1.95	0.47
1:D:262:ASN:ND2	1:D:310:VAL:HG12	2.30	0.47
1:C:227:TRP:CZ2	1:C:229:ARG:HB3	2.49	0.47
1:C:195:PHE:CE2	1:C:257:LEU:HD22	2.49	0.47
1:A:262:ASN:ND2	1:A:310:VAL:HG12	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:GLN:OE1	1:A:86:ARG:NE	2.41	0.47
1:B:156:ALA:HB1	1:C:3:TYR:CG	2.50	0.47
1:B:272:GLN:OE1	1:B:317:VAL:HG13	2.15	0.47
1:D:186:GLU:H	1:D:186:GLU:CD	2.18	0.47
1:D:23:LEU:O	1:D:27:ILE:HG13	2.15	0.47
1:B:210:TRP:CZ2	1:B:234:LEU:HD22	2.50	0.46
1:C:177:LYS:HA	1:C:244:LEU:HD21	1.96	0.46
1:B:285:PRO:C	1:B:287:SER:H	2.19	0.46
1:C:210:TRP:CZ2	1:C:234:LEU:HD22	2.51	0.46
1:C:31:THR:OG1	1:C:367:MSE:HE1	2.16	0.46
1:C:323:ARG:O	1:C:327:LYS:HE3	2.16	0.46
1:C:92:TRP:O	1:C:96:ARG:HB2	2.14	0.46
1:A:238:ILE:HB	1:A:239:PRO:HD3	1.98	0.46
1:A:29:LYS:HD3	1:A:331:PRO:O	2.16	0.46
1:D:87:LYS:HD2	3:D:422:HOH:O	2.16	0.46
1:A:166:ASP:HB3	2:A:380:GOL:H31	1.97	0.46
1:A:325:ILE:O	1:A:329:ILE:HG13	2.16	0.46
1:A:48:LEU:C	1:A:50:ASP:H	2.17	0.46
1:B:344:MSE:CE	1:B:351:TYR:OH	2.64	0.46
1:B:57:LYS:HB2	1:B:57:LYS:NZ	2.30	0.46
1:C:109:MSE:HE2	1:C:109:MSE:HB3	1.82	0.46
1:D:356:LEU:N	1:D:356:LEU:CD1	2.79	0.46
1:B:229:ARG:HH22	1:B:339:SER:HB2	1.78	0.45
1:B:39:ASP:C	1:B:41:THR:N	2.70	0.45
1:B:44:PHE:HD2	1:B:57:LYS:HE2	1.82	0.45
1:C:232:SER:OG	1:C:296:GLY:HA3	2.15	0.45
1:B:233:TRP:HA	1:B:233:TRP:CE3	2.51	0.45
1:B:340:PHE:CG	1:B:341:GLY:N	2.84	0.45
1:B:48:LEU:HB3	1:B:50:ASP:HB2	1.98	0.45
1:D:227:TRP:CZ2	1:D:229:ARG:HB3	2.52	0.45
1:D:56:THR:CG2	1:D:359:MSE:HE1	2.33	0.45
1:C:210:TRP:CD2	1:C:234:LEU:HD13	2.51	0.45
1:C:356:LEU:HD12	1:C:356:LEU:N	2.32	0.45
1:B:214:TRP:CG	1:B:215:SER:N	2.84	0.45
1:D:284:ASP:O	1:D:287:SER:HB3	2.16	0.45
1:D:367:MSE:HB2	1:D:367:MSE:HE3	1.61	0.45
1:B:143:PRO:HD2	1:B:153:ILE:CD1	2.44	0.45
1:C:232:SER:HA	1:C:297:PHE:CE1	2.52	0.45
1:C:23:LEU:HD11	1:C:370:LEU:HD22	1.99	0.45
1:D:302:LEU:HG	1:D:318:ALA:HB1	1.99	0.45
1:B:349:ASP:HA	1:B:352:ARG:HD2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:34:LEU:O	1:C:34:LEU:HD23	2.17	0.45
1:D:8:SER:HA	1:D:9:PRO:HD3	1.87	0.45
1:B:333:GLY:O	1:B:363:GLN:NE2	2.48	0.44
1:C:222:PHE:CB	1:C:344:MSE:HE2	2.47	0.44
1:B:153:ILE:CG2	1:B:158:GLU:HA	2.47	0.44
1:B:227:TRP:CZ2	1:B:229:ARG:HB3	2.52	0.44
1:C:34:LEU:C	1:C:34:LEU:HD23	2.37	0.44
1:C:290:GLU:OE2	1:C:339:SER:HB3	2.17	0.44
1:A:92:TRP:O	1:A:96:ARG:HG2	2.17	0.44
1:B:171:THR:C	1:B:174:PRO:HD2	2.38	0.44
1:B:227:TRP:CZ2	1:B:342:THR:HB	2.53	0.44
1:C:128:LEU:HB3	1:C:129:PRO:HD3	2.00	0.44
1:D:163:MSE:SE	1:D:201:ASN:HD22	2.51	0.44
1:B:356:LEU:HD23	1:B:356:LEU:N	2.33	0.44
1:A:286:HIS:C	1:A:286:HIS:HD1	2.21	0.44
1:B:290:GLU:OE1	1:B:292:SER:OG	2.35	0.44
1:B:63:GLU:OE2	1:B:64:TRP:CD1	2.70	0.44
1:A:286:HIS:C	1:A:286:HIS:ND1	2.70	0.43
1:B:106:VAL:HG22	1:B:142:MSE:CE	2.47	0.43
1:C:171:THR:C	1:C:174:PRO:HD2	2.39	0.43
1:C:377:TYR:N	1:C:377:TYR:CD1	2.86	0.43
1:D:29:LYS:HD2	1:D:331:PRO:O	2.18	0.43
1:A:163:MSE:HE3	1:A:198:HIS:CE1	2.53	0.43
1:B:169:MSE:CE	1:B:173:LEU:HD23	2.48	0.43
1:B:186:GLU:H	1:B:186:GLU:CD	2.22	0.43
1:B:34:LEU:HD22	1:B:73:MSE:CE	2.49	0.43
1:C:327:LYS:HB2	1:C:327:LYS:HE3	1.77	0.43
1:C:281:LEU:HD12	1:C:281:LEU:HA	1.79	0.43
1:C:96:ARG:HA	1:C:96:ARG:HD3	1.51	0.43
1:D:160:HIS:O	1:D:161:GLN:HB2	2.18	0.43
1:B:109:MSE:CE	1:B:174:PRO:HB2	2.49	0.43
1:C:62:TRP:CD2	1:C:102:THR:OG1	2.72	0.43
1:D:173:LEU:HA	1:D:173:LEU:HD12	1.88	0.43
1:B:366:ALA:O	1:B:370:LEU:HG	2.18	0.43
1:B:85:MSE:HA	1:B:85:MSE:HE2	2.01	0.43
1:A:109:MSE:HE3	1:A:138:ALA:HB2	2.00	0.43
1:A:340:PHE:CG	1:A:341:GLY:N	2.86	0.43
1:B:63:GLU:OE1	1:B:155:LEU:HD12	2.18	0.43
1:B:279:HIS:CE1	1:B:289:LEU:HD11	2.54	0.43
1:B:288:TYR:CD1	1:B:288:TYR:C	2.92	0.43
1:C:69:GLY:HA2	1:C:364:ALA:CB	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:63:GLU:OE2	1:D:64:TRP:CD1	2.68	0.43
1:B:232:SER:HB2	1:B:296:GLY:HA3	2.00	0.43
1:C:212:HIS:CD2	1:C:344:MSE:SE	3.22	0.43
1:D:15:GLU:HG3	1:D:15:GLU:O	2.19	0.43
1:B:85:MSE:HA	1:B:85:MSE:HE3	1.99	0.43
1:C:283:ASP:O	1:C:285:PRO:HD3	2.19	0.43
1:D:340:PHE:CG	1:D:341:GLY:N	2.87	0.43
1:D:355:PRO:C	1:D:356:LEU:HD12	2.38	0.43
1:C:139:MSE:HG3	1:C:187:TYR:CE1	2.53	0.43
1:A:30:VAL:HG12	1:A:367:MSE:HE2	2.01	0.42
1:A:233:TRP:HA	1:A:233:TRP:CE3	2.54	0.42
1:C:210:TRP:CZ3	1:C:234:LEU:HD11	2.55	0.42
1:C:86:ARG:HE	1:C:86:ARG:HB3	1.74	0.42
1:D:28:GLN:CG	2:D:381:GOL:H31	2.40	0.42
1:B:285:PRO:C	1:B:287:SER:N	2.72	0.42
1:B:212:HIS:HD2	1:B:344:MSE:CE	2.31	0.42
1:A:344:MSE:HE2	1:A:344:MSE:HB2	1.95	0.42
1:C:220:HIS:ND1	1:C:222:PHE:N	2.60	0.42
1:C:302:LEU:HG	1:C:318:ALA:HB1	2.02	0.42
1:D:8:SER:O	1:D:12:ARG:HG2	2.20	0.42
1:D:259:GLN:HB3	1:D:259:GLN:HE21	1.72	0.42
1:D:40:GLU:HG2	1:D:40:GLU:H	1.66	0.42
1:C:272:GLN:OE1	1:C:317:VAL:HG13	2.19	0.42
1:B:212:HIS:CD2	1:B:344:MSE:CE	3.03	0.42
1:C:63:GLU:CG	1:C:64:TRP:H	2.32	0.42
1:D:204:ASP:HB2	1:D:211:PHE:CE2	2.54	0.42
1:A:57:LYS:HB2	1:A:359:MSE:SE	2.70	0.42
1:B:334:GLU:HA	1:B:363:GLN:HE22	1.84	0.42
1:A:110:ALA:HB3	1:A:111:PRO:HD3	2.02	0.41
1:C:109:MSE:O	1:C:109:MSE:CE	2.67	0.41
1:C:137:TRP:CZ2	1:C:142:MSE:HB2	2.55	0.41
1:B:152:HIS:CD2	1:B:164:TRP:CE3	3.08	0.41
1:A:210:TRP:CD2	1:A:234:LEU:HD13	2.56	0.41
1:C:65:THR:HG23	1:C:361:TYR:HD2	1.86	0.41
1:A:259:GLN:HB3	1:A:259:GLN:HE21	1.72	0.41
1:C:173:LEU:HB3	1:C:174:PRO:HD3	2.02	0.41
1:B:39:ASP:HB3	1:B:57:LYS:HD2	2.03	0.41
1:C:13:GLN:CG	1:C:14:PRO:HD2	2.50	0.41
1:A:3:TYR:HA	1:A:4:PRO:HD3	1.79	0.41
1:C:110:ALA:N	1:C:111:PRO:CD	2.83	0.41
1:C:63:GLU:CG	1:C:64:TRP:N	2.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:TRP:CG	1:B:234:LEU:HD13	2.56	0.41
1:B:367:MSE:HE2	1:B:367:MSE:HB3	1.99	0.41
1:C:212:HIS:CD2	1:C:351:TYR:OH	2.73	0.41
1:D:195:PHE:O	1:D:199:VAL:HG23	2.21	0.41
1:D:28:GLN:HB2	2:D:381:GOL:C2	2.51	0.41
1:A:109:MSE:CE	1:A:137:TRP:CE3	3.04	0.41
1:A:168:LEU:HD23	1:A:172:VAL:HG21	2.03	0.41
1:B:74:TYR:CZ	1:B:117:TYR:HB3	2.56	0.41
1:A:316:GLN:O	1:A:320:LYS:HG3	2.21	0.40
1:C:100:GLY:O	2:C:381:GOL:H32	2.21	0.40
1:D:49:ASP:C	1:D:49:ASP:OD2	2.60	0.40
1:A:119:TYR:HD2	1:A:181:LEU:HD21	1.86	0.40
1:A:173:LEU:HB3	1:A:174:PRO:HD3	2.02	0.40
1:A:210:TRP:CZ3	1:A:234:LEU:HD11	2.57	0.40
1:B:123:ARG:HD3	1:B:123:ARG:HA	1.90	0.40
1:B:39:ASP:OD1	1:B:39:ASP:O	2.39	0.40
1:C:85:MSE:HA	1:C:85:MSE:HE2	1.96	0.40
1:D:212:HIS:HD2	1:D:351:TYR:OH	2.05	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	376/381 (99%)	364 (97%)	11 (3%)	1 (0%)	41 43
1	B	376/381 (99%)	357 (95%)	16 (4%)	3 (1%)	19 17
1	C	376/381 (99%)	367 (98%)	8 (2%)	1 (0%)	41 43
1	D	376/381 (99%)	367 (98%)	9 (2%)	0	100 100
All	All	1504/1524 (99%)	1455 (97%)	44 (3%)	5 (0%)	41 43

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	274	GLU
1	B	341	GLY
1	B	45	LEU
1	B	35	VAL
1	C	4	PRO

### 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	321/308 (104%)	305 (95%)	16 (5%)	24 27
1	B	321/308 (104%)	301 (94%)	20 (6%)	18 19
1	C	321/308 (104%)	308 (96%)	13 (4%)	31 37
1	D	321/308 (104%)	312 (97%)	9 (3%)	43 53
All	All	1284/1232 (104%)	1226 (96%)	58 (4%)	27 32

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	41	THR
1	A	46	LEU
1	A	48	LEU
1	A	50	ASP
1	A	56	THR
1	A	195	PHE
1	A	234	LEU
1	A	235	THR
1	A	259	GLN
1	A	271	CYS
1	A	281	LEU
1	A	286	HIS
1	A	287	SER
1	A	308	ARG
1	A	327	LYS
1	B	1	MSE

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Mol	Chain	Res	Type
1	B	41	THR
1	B	50	ASP
1	B	65	THR
1	B	85	MSE
1	B	86	ARG
1	B	102	THR
1	B	109	MSE
1	B	153	ILE
1	B	166	ASP
1	B	235	THR
1	B	259	GLN
1	B	281	LEU
1	B	288	TYR
1	B	289	LEU
1	B	290	GLU
1	B	303	LYS
1	B	347	ASP
1	B	354	ILE
1	B	356	LEU
1	C	1	MSE
1	C	2	VAL
1	C	12	ARG
1	C	85	MSE
1	C	96	ARG
1	C	102	THR
1	C	109	MSE
1	C	234	LEU
1	C	235	THR
1	C	259	GLN
1	C	281	LEU
1	C	287	SER
1	C	327	LYS
1	D	38	LYS
1	D	63	GLU
1	D	96	ARG
1	D	195	PHE
1	D	235	THR
1	D	250	ASN
1	D	259	GLN
1	D	281	LEU
1	D	287	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	160	HIS
1	A	200	GLN
1	A	259	GLN
1	B	107	ASN
1	B	198	HIS
1	B	200	GLN
1	B	212	HIS
1	B	259	GLN
1	B	353	HIS
1	C	107	ASN
1	C	198	HIS
1	C	212	HIS
1	C	221	ASN
1	C	259	GLN
1	D	16	HIS
1	D	107	ASN
1	D	198	HIS
1	D	212	HIS
1	D	250	ASN
1	D	259	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 [\(i\)](#)

11 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
2	GOL	C	379	-	5,5,5	0.38	0	5,5,5	0.56	0
2	GOL	C	380	-	5,5,5	0.35	0	5,5,5	0.33	0
2	GOL	B	379	-	5,5,5	0.38	0	5,5,5	0.28	0
2	GOL	D	379	-	5,5,5	0.35	0	5,5,5	0.72	0
2	GOL	A	381	-	5,5,5	0.37	0	5,5,5	0.28	0
2	GOL	C	381	-	5,5,5	0.28	0	5,5,5	0.47	0
2	GOL	A	379	-	5,5,5	0.30	0	5,5,5	0.34	0
2	GOL	D	382	-	5,5,5	0.35	0	5,5,5	0.27	0
2	GOL	D	381	-	5,5,5	0.53	0	5,5,5	0.42	0
2	GOL	A	380	-	5,5,5	0.37	0	5,5,5	0.19	0
2	GOL	D	380	-	5,5,5	0.37	0	5,5,5	0.36	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
2	GOL	C	379	-	-	4/4/4/4	-
2	GOL	C	380	-	-	2/4/4/4	-
2	GOL	B	379	-	-	3/4/4/4	-
2	GOL	D	379	-	-	2/4/4/4	-
2	GOL	A	381	-	-	0/4/4/4	-
2	GOL	C	381	-	-	4/4/4/4	-
2	GOL	A	379	-	-	4/4/4/4	-
2	GOL	D	382	-	-	2/4/4/4	-
2	GOL	D	381	-	-	2/4/4/4	-
2	GOL	A	380	-	-	2/4/4/4	-
2	GOL	D	380	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	379	GOL	C1-C2-C3-O3
2	C	380	GOL	O1-C1-C2-C3
2	B	379	GOL	O1-C1-C2-C3
2	C	381	GOL	O1-C1-C2-C3
2	C	381	GOL	C1-C2-C3-O3
2	A	379	GOL	O1-C1-C2-C3
2	A	379	GOL	C1-C2-C3-O3
2	D	381	GOL	O1-C1-C2-C3
2	A	380	GOL	C1-C2-C3-O3
2	C	381	GOL	O2-C2-C3-O3
2	D	379	GOL	O1-C1-C2-C3
2	D	382	GOL	C1-C2-C3-O3
2	C	380	GOL	O1-C1-C2-O2
2	B	379	GOL	O1-C1-C2-O2
2	C	381	GOL	O1-C1-C2-O2
2	A	379	GOL	O1-C1-C2-O2
2	A	379	GOL	O2-C2-C3-O3
2	D	381	GOL	O1-C1-C2-O2
2	A	380	GOL	O2-C2-C3-O3
2	D	382	GOL	O2-C2-C3-O3
2	C	379	GOL	O2-C2-C3-O3
2	D	379	GOL	O1-C1-C2-O2
2	C	379	GOL	O1-C1-C2-O2
2	C	379	GOL	O1-C1-C2-C3
2	B	379	GOL	C1-C2-C3-O3

There are no ring outliers.

5 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	381	GOL	1	0
2	A	379	GOL	1	0
2	D	382	GOL	1	0
2	D	381	GOL	5	0
2	A	380	GOL	3	0

## 5.7 Other polymers [\(i\)](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [\(i\)](#)

There are no chain breaks in this entry.

## 6 Fit of model and data i

### 6.1 Protein, DNA and RNA chains i

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	363/381 (95%)	0.10	20 (5%) 25 26	23, 42, 67, 104	0
1	B	363/381 (95%)	-0.20	4 (1%) 80 80	23, 40, 69, 114	0
1	C	363/381 (95%)	-0.04	7 (1%) 66 67	28, 48, 74, 109	0
1	D	363/381 (95%)	-0.12	11 (3%) 50 51	21, 35, 59, 90	0
All	All	1452/1524 (95%)	-0.06	42 (2%) 51 52	21, 41, 69, 114	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	233	TRP	3.7
1	A	236	ILE	3.5
1	C	236	ILE	3.5
1	A	286	HIS	3.3
1	C	15	GLU	3.2
1	A	237	VAL	3.2
1	B	45	LEU	3.2
1	C	2	VAL	3.0
1	A	168	LEU	3.0
1	A	68	VAL	2.9
1	A	15	GLU	2.8
1	A	113	LEU	2.7
1	A	172	VAL	2.6
1	A	299	TYR	2.6
1	D	41	THR	2.5
1	C	233	TRP	2.5
1	A	110	ALA	2.5
1	A	64	TRP	2.5
1	D	53	VAL	2.4
1	D	68	VAL	2.4
1	C	40	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	234	LEU	2.4
1	D	8	SER	2.4
1	D	168	LEU	2.3
1	A	173	LEU	2.2
1	D	46	LEU	2.2
1	D	368	LEU	2.2
1	A	331	PRO	2.2
1	D	233	TRP	2.2
1	B	40	GLU	2.2
1	A	53	VAL	2.2
1	B	44	PHE	2.2
1	A	332	GLU	2.2
1	A	238	ILE	2.1
1	C	41	THR	2.1
1	B	361	TYR	2.1
1	D	286	HIS	2.1
1	D	2	VAL	2.1
1	D	237	VAL	2.0
1	A	2	VAL	2.0
1	A	48	LEU	2.0
1	C	329	ILE	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<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
2	GOL	D	382	6/6	0.71	0.38	74,77,79,80	0
2	GOL	A	381	6/6	0.79	0.23	47,51,57,62	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	GOL	D	381	6/6	0.81	0.17	52,67,70,72	0
2	GOL	C	380	6/6	0.84	0.22	49,57,69,73	0
2	GOL	C	381	6/6	0.87	0.15	59,68,75,77	0
2	GOL	A	379	6/6	0.90	0.23	38,52,55,61	0
2	GOL	D	379	6/6	0.92	0.17	38,46,55,58	0
2	GOL	C	379	6/6	0.93	0.18	46,52,54,54	0
2	GOL	B	379	6/6	0.94	0.13	40,49,55,69	0
2	GOL	A	380	6/6	0.95	0.28	42,53,56,58	0
2	GOL	D	380	6/6	0.95	0.17	26,40,45,45	0

## 6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.