Full wwPDB X-ray Structure Validation Report (i)

Jun 17, 2024 – 06:03 AM EDT

PDB ID	:	3RCE
Title	:	Bacterial oligosaccharyltransferase PglB
Authors	:	Lizak, C.; Gerber, S.; Numao, S.; Aebi, M.; Locher, K.P.
Deposited on	:	2011-03-31
Resolution	:	3.40 Å(reported)

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

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

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

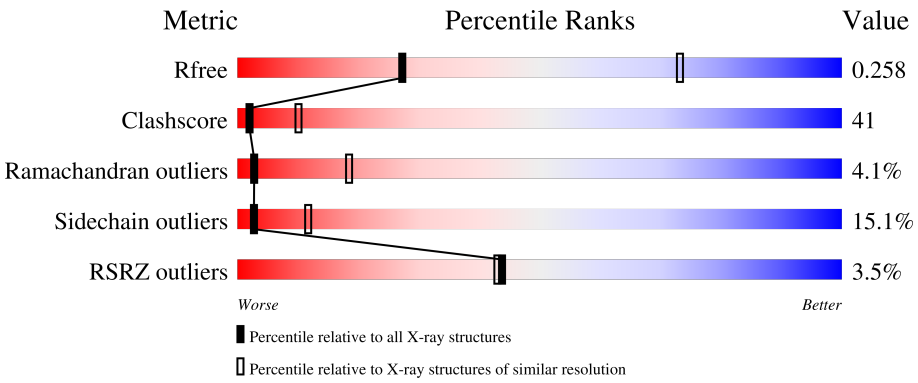
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.40 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	1026 (3.48-3.32)
Clashscore	141614	1055 (3.48-3.32)
Ramachandran outliers	138981	1038 (3.48-3.32)
Sidechain outliers	138945	1038 (3.48-3.32)
RSRZ outliers	127900	2173 (3.50-3.30)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	724	3%; 36%, 47%, 10%, • 6%
2	B	8	62%, 12%, 25%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5668 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 Oligosaccharide transferase to N-glycosylate proteins.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	683	5606	3730	857	990	29	0	0	0

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	GLU	LYS	ENGINEERED MUTATION	UNP B9KDD4
A	108	THR	ALA	SEE REMARK 999	UNP B9KDD4
A	187	ILE	VAL	SEE REMARK 999	UNP B9KDD4
A	227	ALA	THR	SEE REMARK 999	UNP B9KDD4
A	275	SER	LEU	SEE REMARK 999	UNP B9KDD4
A	536	GLN	LYS	SEE REMARK 999	UNP B9KDD4
A	550	ASN	ASP	SEE REMARK 999	UNP B9KDD4
A	681	ALA	SER	SEE REMARK 999	UNP B9KDD4
A	699	VAL	ILE	SEE REMARK 999	UNP B9KDD4
A	713	GLU	-	EXPRESSION TAG	UNP B9KDD4
A	714	PHE	-	EXPRESSION TAG	UNP B9KDD4
A	715	HIS	-	EXPRESSION TAG	UNP B9KDD4
A	716	HIS	-	EXPRESSION TAG	UNP B9KDD4
A	717	HIS	-	EXPRESSION TAG	UNP B9KDD4
A	718	HIS	-	EXPRESSION TAG	UNP B9KDD4
A	719	HIS	-	EXPRESSION TAG	UNP B9KDD4
A	720	HIS	-	EXPRESSION TAG	UNP B9KDD4
A	721	HIS	-	EXPRESSION TAG	UNP B9KDD4
A	722	HIS	-	EXPRESSION TAG	UNP B9KDD4
A	723	HIS	-	EXPRESSION TAG	UNP B9KDD4
A	724	HIS	-	EXPRESSION TAG	UNP B9KDD4

- Molecule 2 is a protein called Substrate Mimic Peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	8	59	33	11	15	5	0	0

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
3	A	2	2	2	0	0

- Molecule 4 is water.

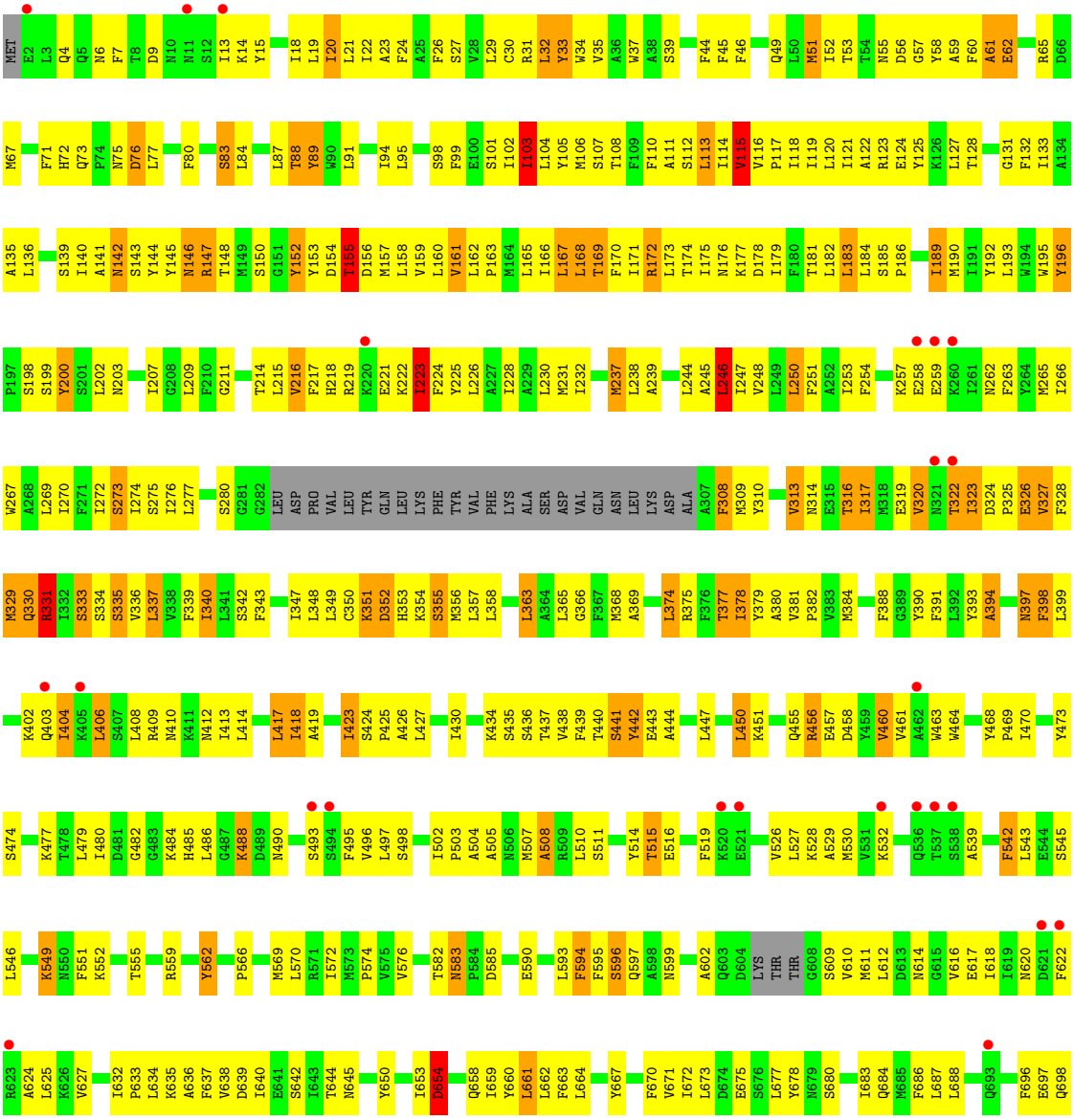
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	1	1	1	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

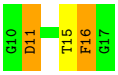
- Molecule 1: Oligosaccharide transferase to N-glycosylate proteins

Chain A: 3% 36% 47% 10% • 6%



• Molecule 2: Substrate Mimic Peptide

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	85.06Å 116.10Å 175.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.98 – 3.40 29.98 – 3.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (29.98-3.40) 97.4 (29.98-3.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.70 (at 3.39Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.238 , 0.271 0.231 , 0.258	Depositor DCC
R_{free} test set	2000 reflections (8.39%)	wwPDB-VP
Wilson B-factor (Å^2)	112.7	Xtriage
Anisotropy	0.015	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.25 , 71.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	5668	wwPDB-VP
Average B, all atoms (Å^2)	129.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.07% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

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

5 Model quality (i)

5.1 Standard geometry (i)

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

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	0.53	0/5752	0.71	2/7791 (0.0%)
2	B	0.53	0/43	0.83	0/55
All	All	0.53	0/5795	0.71	2/7846 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	337	LEU	CA-CB-CG	5.76	128.54	115.30
1	A	246	LEU	CA-CB-CG	5.10	127.03	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5606	0	5628	465	0
2	B	59	0	42	3	0
3	A	2	0	0	0	0
4	A	1	0	0	0	0
All	All	5668	0	5670	466	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 41.

All (466) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:THR:HG21	1:A:185:SER:HB3	1.22	1.14
1:A:488:LYS:O	1:A:488:LYS:HD3	1.47	1.09
1:A:20:ILE:HG13	1:A:119:ILE:HD11	1.25	1.06
1:A:329:MET:HE1	1:A:336:VAL:HG13	1.39	1.03
1:A:173:LEU:HB2	1:A:181:THR:HG21	1.41	0.98
1:A:572:ILE:HG22	1:A:576:VAL:HG23	1.47	0.96
1:A:140:ILE:HG22	1:A:426:ALA:HB1	1.48	0.95
1:A:169:THR:CG2	1:A:185:SER:HB3	1.96	0.95
1:A:456:ARG:HA	1:A:456:ARG:HH11	1.34	0.92
1:A:442:TYR:HD1	1:A:442:TYR:H	1.18	0.92
1:A:237:MET:HG3	1:A:280:SER:HB3	1.51	0.91
1:A:144:TYR:HA	1:A:378:ILE:HD11	1.53	0.89
1:A:323:ILE:HD11	1:A:374:LEU:HD12	1.56	0.87
1:A:329:MET:CE	1:A:336:VAL:HG13	2.06	0.86
1:A:325:PRO:O	1:A:329:MET:HG2	1.79	0.83
1:A:664:LEU:HD21	1:A:687:LEU:HD22	1.61	0.82
1:A:67:MET:HA	1:A:80:PHE:HE1	1.44	0.82
1:A:67:MET:HG2	1:A:80:PHE:CD1	2.13	0.82
1:A:336:VAL:O	1:A:340:ILE:HG12	1.81	0.81
1:A:125:TYR:HE1	1:A:356:MET:HE3	1.45	0.81
1:A:31:ARG:NH2	1:A:139:SER:O	2.14	0.81
1:A:488:LYS:HD3	1:A:488:LYS:C	2.00	0.80
1:A:186:PRO:O	1:A:190:MET:HG3	1.80	0.80
1:A:31:ARG:HB2	1:A:108:THR:HG23	1.63	0.80
1:A:663:PHE:CD1	1:A:670:PHE:CE2	2.70	0.79
1:A:488:LYS:NZ	1:A:526:VAL:N	2.30	0.79
1:A:583:ASN:HD21	1:A:585:ASP:HB2	1.48	0.79
1:A:317:ILE:HD11	1:A:485:HIS:CE1	2.17	0.78
1:A:460:VAL:HG11	1:A:470:ILE:HG21	1.65	0.78
1:A:125:TYR:OH	1:A:171:ILE:HD13	1.85	0.77
1:A:460:VAL:HG23	1:A:562:TYR:HB2	1.67	0.77
1:A:270:ILE:O	1:A:274:ILE:HG13	1.84	0.77
1:A:169:THR:HG21	1:A:185:SER:CB	2.08	0.76
1:A:698:GLN:HE22	1:A:701:ASN:HD22	1.30	0.76
1:A:20:ILE:CG1	1:A:119:ILE:HD11	2.11	0.75
1:A:334:SER:O	1:A:335:SER:HB2	1.84	0.75
1:A:53:THR:HG22	1:A:53:THR:O	1.85	0.75
1:A:207:ILE:HG23	1:A:231:MET:HE2	1.66	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:ARG:NH2	1:A:319:GLU:HG3	2.02	0.74
1:A:263:PHE:HB3	1:A:267:TRP:HE1	1.52	0.74
1:A:179:ILE:HG23	1:A:182:LEU:HD12	1.68	0.74
1:A:322:THR:O	1:A:323:ILE:HG12	1.88	0.74
1:A:313:VAL:HG21	1:A:485:HIS:O	1.88	0.74
1:A:67:MET:HA	1:A:80:PHE:CE1	2.23	0.73
1:A:18:ILE:O	1:A:22:ILE:HG13	1.88	0.73
1:A:317:ILE:HD11	1:A:485:HIS:HE1	1.52	0.73
1:A:663:PHE:HD1	1:A:670:PHE:CD2	2.07	0.73
1:A:88:THR:HG23	1:A:106:MET:SD	2.28	0.72
1:A:507:MET:CE	1:A:511:SER:HB3	2.19	0.72
1:A:663:PHE:HD1	1:A:670:PHE:CE2	2.08	0.72
1:A:83:SER:HB3	1:A:195:TRP:HA	1.71	0.71
1:A:60:PHE:CD1	1:A:84:LEU:HD23	2.25	0.71
1:A:91:LEU:O	1:A:95:LEU:HD12	1.89	0.71
1:A:145:TYR:CZ	1:A:430:ILE:HG23	2.25	0.71
1:A:262:ASN:H	1:A:265:MET:HE3	1.53	0.71
1:A:409:ARG:O	1:A:413:ILE:HG13	1.91	0.70
1:A:33:TYR:C	1:A:33:TYR:CD2	2.63	0.70
1:A:343:PHE:HD1	1:A:384:MET:HE3	1.55	0.70
1:A:73:GLN:NE2	1:A:456:ARG:HB3	2.07	0.70
1:A:595:PHE:CD2	1:A:596:SER:N	2.60	0.69
1:A:661:LEU:HD12	1:A:662:LEU:N	2.07	0.69
1:A:488:LYS:NZ	1:A:526:VAL:H	1.90	0.69
1:A:322:THR:C	1:A:323:ILE:HG12	2.13	0.69
1:A:437:THR:HB	1:A:439:PHE:O	1.93	0.69
1:A:246:LEU:HD12	1:A:250:LEU:HD12	1.75	0.69
1:A:26:PHE:CD2	1:A:136:LEU:HD22	2.28	0.68
1:A:125:TYR:HE1	1:A:356:MET:CE	2.06	0.68
1:A:507:MET:HE3	1:A:511:SER:HB3	1.75	0.68
1:A:20:ILE:HG13	1:A:119:ILE:CD1	2.13	0.68
1:A:637:PHE:HE2	1:A:639:ASP:HB2	1.58	0.68
1:A:158:LEU:HB2	1:A:195:TRP:CH2	2.28	0.68
1:A:618:ILE:HD13	1:A:661:LEU:HD21	1.76	0.68
1:A:323:ILE:CG2	1:A:327:VAL:HB	2.24	0.67
1:A:394:ALA:O	1:A:397:ASN:HB2	1.94	0.67
1:A:124:GLU:HA	1:A:124:GLU:OE1	1.95	0.67
1:A:507:MET:O	1:A:508:ALA:C	2.31	0.67
1:A:262:ASN:O	1:A:265:MET:HG2	1.95	0.67
1:A:147:ARG:HH21	1:A:319:GLU:HG3	1.60	0.67
1:A:207:ILE:HG12	1:A:231:MET:HE3	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:528:LYS:HE2	1:A:532:LYS:HE3	1.77	0.67
1:A:642:SER:HB3	1:A:645:ASN:HD22	1.59	0.67
1:A:172:ARG:HH11	1:A:172:ARG:HB3	1.60	0.66
1:A:39:SER:HA	1:A:45:PHE:CZ	2.30	0.66
1:A:323:ILE:HG21	1:A:327:VAL:HB	1.78	0.66
1:A:75:ASN:ND2	1:A:519:PHE:HB3	2.10	0.66
1:A:140:ILE:HG22	1:A:426:ALA:CB	2.25	0.66
1:A:111:ALA:CB	1:A:148:THR:CG2	2.73	0.66
1:A:461:VAL:HG22	1:A:479:LEU:HD12	1.78	0.66
1:A:569:MET:O	1:A:569:MET:HG2	1.95	0.66
1:A:148:THR:O	1:A:148:THR:HG22	1.95	0.65
1:A:111:ALA:HB3	1:A:148:THR:CG2	2.27	0.65
1:A:488:LYS:HZ3	1:A:527:LEU:H	1.43	0.65
1:A:624:ALA:HA	1:A:632:ILE:O	1.97	0.65
1:A:698:GLN:NE2	1:A:701:ASN:HB2	2.11	0.65
1:A:510:LEU:HB3	1:A:530:MET:HE1	1.77	0.65
1:A:640:ILE:CD1	1:A:688:LEU:HD23	2.27	0.65
1:A:442:TYR:CD1	1:A:442:TYR:N	2.66	0.64
1:A:349:LEU:CD2	1:A:356:MET:HG2	2.27	0.64
1:A:413:ILE:HG22	1:A:417:LEU:CD1	2.28	0.64
1:A:195:TRP:CD2	1:A:195:TRP:O	2.51	0.64
1:A:503:PRO:HA	1:A:546:LEU:O	1.97	0.64
1:A:4:GLN:O	1:A:4:GLN:HG2	1.98	0.64
1:A:193:LEU:HD13	1:A:203:ASN:ND2	2.13	0.64
1:A:514:TYR:HE1	1:A:555:THR:HG21	1.63	0.64
1:A:514:TYR:CE1	1:A:555:THR:HG21	2.32	0.64
1:A:119:ILE:HG22	1:A:131:GLY:C	2.18	0.63
1:A:355:SER:O	1:A:358:LEU:HB2	1.97	0.63
1:A:640:ILE:HD11	1:A:688:LEU:HD23	1.79	0.63
1:A:664:LEU:HD21	1:A:687:LEU:CD2	2.27	0.63
1:A:37:TRP:CE2	1:A:434:LYS:HD3	2.34	0.63
1:A:71:PHE:CD2	1:A:72:HIS:O	2.52	0.63
1:A:488:LYS:HZ2	1:A:526:VAL:H	1.46	0.63
1:A:354:LYS:O	1:A:357:LEU:HB2	1.99	0.63
1:A:413:ILE:O	1:A:417:LEU:HD12	1.99	0.62
1:A:111:ALA:CB	1:A:148:THR:HG21	2.30	0.62
1:A:542:PHE:C	1:A:542:PHE:CD2	2.73	0.62
1:A:195:TRP:O	1:A:195:TRP:CE3	2.52	0.62
1:A:230:LEU:HD22	1:A:247:ILE:HD13	1.80	0.61
1:A:272:ILE:HG22	1:A:273:SER:N	2.14	0.61
1:A:57:GLY:HA2	1:A:153:TYR:O	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:118:ILE:HB	1:A:135:ALA:HB2	1.83	0.60
1:A:616:VAL:HG13	1:A:627:VAL:HG22	1.83	0.60
1:A:203:ASN:O	1:A:207:ILE:HG13	2.01	0.60
1:A:155:THR:HG23	1:A:155:THR:O	2.02	0.60
1:A:158:LEU:H	1:A:195:TRP:HH2	1.49	0.60
1:A:202:LEU:HD23	1:A:365:LEU:HD12	1.84	0.60
1:A:698:GLN:NE2	1:A:701:ASN:HD22	2.00	0.60
1:A:632:ILE:HG21	1:A:659:ILE:HD12	1.84	0.59
1:A:37:TRP:CH2	1:A:434:LYS:HB3	2.37	0.59
1:A:450:LEU:O	1:A:450:LEU:HD22	2.02	0.59
1:A:622:PHE:O	1:A:622:PHE:CD1	2.55	0.59
1:A:526:VAL:O	1:A:529:ALA:HB3	2.03	0.59
1:A:582:THR:HG22	1:A:582:THR:O	2.01	0.59
1:A:632:ILE:HG23	1:A:633:PRO:HD2	1.84	0.59
1:A:179:ILE:HG22	1:A:251:PHE:CE1	2.38	0.58
1:A:595:PHE:CD2	1:A:595:PHE:C	2.76	0.58
1:A:167:LEU:O	1:A:171:ILE:HG13	2.03	0.58
1:A:640:ILE:HD11	1:A:688:LEU:CD2	2.33	0.58
1:A:625:LEU:HD21	1:A:634:LEU:HD21	1.84	0.58
1:A:75:ASN:HD21	1:A:519:PHE:HB3	1.67	0.58
1:A:263:PHE:HB3	1:A:267:TRP:NE1	2.18	0.58
1:A:329:MET:HE3	1:A:339:PHE:CD2	2.38	0.58
1:A:330:GLN:O	1:A:334:SER:N	2.31	0.58
1:A:104:LEU:HD23	1:A:105:TYR:CE1	2.39	0.57
1:A:31:ARG:CB	1:A:108:THR:HG23	2.32	0.57
1:A:365:LEU:HD13	1:A:379:TYR:CD1	2.40	0.57
1:A:468:TYR:HB2	1:A:469:PRO:HD3	1.87	0.57
1:A:88:THR:CG2	1:A:106:MET:SD	2.93	0.57
1:A:331:ARG:NH1	1:A:377:THR:OG1	2.38	0.57
1:A:543:LEU:C	1:A:545:SER:H	2.07	0.56
1:A:53:THR:O	1:A:53:THR:CG2	2.52	0.56
1:A:342:SER:OG	1:A:388:PHE:HB2	2.06	0.56
1:A:348:LEU:O	1:A:352:ASP:HB2	2.06	0.56
1:A:530:MET:HE2	1:A:542:PHE:CZ	2.40	0.56
1:A:638:VAL:HB	1:A:662:LEU:CD2	2.35	0.56
1:A:460:VAL:CG2	1:A:562:TYR:HB2	2.35	0.56
1:A:350:CYS:C	1:A:352:ASP:H	2.08	0.56
1:A:238:LEU:HD11	1:A:276:ILE:HG21	1.88	0.56
1:A:464:TRP:CE3	1:A:482:GLY:HA2	2.39	0.56
1:A:595:PHE:HB2	1:A:677:LEU:CD1	2.36	0.56
1:A:15:TYR:CZ	1:A:19:LEU:HD11	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:ARG:HB2	1:A:128:THR:OG1	2.05	0.56
1:A:44:PHE:HE2	1:A:436:SER:HB2	1.71	0.56
1:A:614:ASN:OD1	1:A:614:ASN:N	2.37	0.56
1:A:488:LYS:HZ3	1:A:526:VAL:N	2.03	0.55
1:A:133:ILE:HD13	1:A:419:ALA:HB2	1.86	0.55
1:A:121:ILE:CG1	1:A:168:LEU:HB2	2.37	0.55
1:A:61:ALA:CB	1:A:103:ILE:HD11	2.36	0.55
1:A:635:LYS:HD3	1:A:660:TYR:CZ	2.41	0.55
1:A:207:ILE:HD13	1:A:231:MET:HB3	1.88	0.55
1:A:451:LYS:HG3	1:A:474:SER:O	2.06	0.55
1:A:479:LEU:HD11	1:A:511:SER:OG	2.05	0.55
1:A:185:SER:O	1:A:189:ILE:HB	2.07	0.55
1:A:196:TYR:CE2	1:A:198:SER:HB2	2.42	0.55
1:A:168:LEU:O	1:A:168:LEU:HG	2.07	0.55
1:A:510:LEU:HB3	1:A:530:MET:CE	2.36	0.55
1:A:146:ASN:N	1:A:146:ASN:HD22	2.03	0.54
1:A:313:VAL:HG21	1:A:486:LEU:HA	1.89	0.54
1:A:67:MET:HG2	1:A:80:PHE:CE1	2.41	0.54
1:A:51:MET:SD	1:A:437:THR:HG21	2.48	0.54
1:A:61:ALA:H	1:A:153:TYR:HE2	1.55	0.54
1:A:686:PHE:HD1	1:A:706:LYS:HG3	1.72	0.54
1:A:33:TYR:C	1:A:33:TYR:HD2	2.07	0.54
1:A:406:LEU:HD22	1:A:410:ASN:HB2	1.89	0.54
1:A:31:ARG:NH1	1:A:148:THR:HB	2.23	0.54
1:A:155:THR:O	1:A:195:TRP:CH2	2.60	0.54
1:A:334:SER:O	1:A:335:SER:CB	2.53	0.54
1:A:653:ILE:HG22	1:A:654:ASP:N	2.22	0.54
1:A:198:SER:C	1:A:200:TYR:H	2.09	0.54
2:B:16:PPN:CD1	2:B:16:PPN:C	2.78	0.54
1:A:140:ILE:CG2	1:A:426:ALA:HB1	2.31	0.54
1:A:166:ILE:HD13	1:A:189:ILE:HG12	1.88	0.54
1:A:365:LEU:HD23	1:A:368:MET:HE2	1.90	0.54
1:A:637:PHE:C	1:A:637:PHE:CD2	2.81	0.54
1:A:6:ASN:ND2	1:A:9:ASP:HB2	2.23	0.54
1:A:257:LYS:O	1:A:259:GLU:N	2.41	0.54
1:A:352:ASP:O	1:A:353:HIS:CD2	2.61	0.54
1:A:121:ILE:HG12	1:A:168:LEU:HB2	1.89	0.53
1:A:634:LEU:O	1:A:653:ILE:HB	2.07	0.53
1:A:269:LEU:O	1:A:273:SER:HB3	2.08	0.53
1:A:350:CYS:O	1:A:352:ASP:N	2.39	0.53
1:A:61:ALA:N	1:A:153:TYR:HE2	2.06	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:516:GLU:HG2	1:A:559:ARG:NH1	2.23	0.53
1:A:160:LEU:HD21	1:A:379:TYR:HD2	1.74	0.53
1:A:238:LEU:O	1:A:239:ALA:C	2.47	0.53
1:A:184:LEU:HD23	1:A:184:LEU:N	2.22	0.53
1:A:398:PHE:O	1:A:402:LYS:HB2	2.08	0.53
1:A:625:LEU:HD21	1:A:634:LEU:HD11	1.91	0.53
1:A:23:ALA:HB1	1:A:115:VAL:CG2	2.39	0.53
1:A:87:LEU:HD21	1:A:110:PHE:CZ	2.44	0.53
1:A:460:VAL:CG1	1:A:470:ILE:HG21	2.36	0.53
1:A:530:MET:HE2	1:A:542:PHE:HZ	1.73	0.52
1:A:457:GLU:CG	1:A:559:ARG:HH21	2.22	0.52
1:A:542:PHE:CE2	1:A:546:LEU:HD21	2.44	0.52
1:A:273:SER:HA	1:A:276:ILE:HG13	1.91	0.52
1:A:442:TYR:CB	1:A:704:ARG:HD2	2.39	0.52
1:A:463:TRP:CE2	1:A:576:VAL:HG22	2.44	0.52
1:A:505:ALA:HB2	1:A:696:PHE:CZ	2.45	0.52
1:A:20:ILE:HA	1:A:119:ILE:HD11	1.92	0.52
1:A:23:ALA:HB1	1:A:115:VAL:HG22	1.92	0.52
1:A:209:LEU:HD23	1:A:358:LEU:HA	1.91	0.52
1:A:173:LEU:HD11	1:A:223:ILE:HG21	1.92	0.51
1:A:263:PHE:O	1:A:267:TRP:CD1	2.63	0.51
1:A:154:ASP:C	1:A:156:ASP:H	2.14	0.51
1:A:354:LYS:O	1:A:357:LEU:CB	2.58	0.51
1:A:366:GLY:O	1:A:369:ALA:CB	2.57	0.51
1:A:413:ILE:HG22	1:A:417:LEU:HD12	1.91	0.51
1:A:542:PHE:O	1:A:545:SER:HB3	2.10	0.51
1:A:58:TYR:CE2	1:A:473:TYR:CE1	2.98	0.51
1:A:88:THR:HG21	1:A:153:TYR:CE1	2.46	0.51
1:A:172:ARG:HH11	1:A:172:ARG:CB	2.23	0.51
1:A:308:PHE:CE2	1:A:527:LEU:HD23	2.45	0.51
1:A:329:MET:HE3	1:A:339:PHE:HD2	1.75	0.51
1:A:331:ARG:NH2	2:B:11:ASP:OD2	2.43	0.51
1:A:15:TYR:CE2	1:A:19:LEU:HD11	2.45	0.51
1:A:158:LEU:HB2	1:A:195:TRP:CZ3	2.46	0.51
1:A:320:VAL:HG13	1:A:320:VAL:O	2.10	0.51
1:A:399:LEU:N	1:A:399:LEU:HD23	2.24	0.51
1:A:128:THR:O	1:A:131:GLY:N	2.44	0.51
1:A:108:THR:OG1	1:A:150:SER:HA	2.11	0.51
1:A:366:GLY:HA3	1:A:380:ALA:HB2	1.91	0.51
1:A:442:TYR:CD2	1:A:704:ARG:HD2	2.46	0.51
1:A:39:SER:HA	1:A:45:PHE:HZ	1.73	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:365:LEU:HD13	1:A:379:TYR:CE1	2.46	0.50
1:A:551:PHE:CD2	1:A:552:LYS:O	2.64	0.50
1:A:381:VAL:HG12	1:A:382:PRO:N	2.26	0.50
1:A:660:TYR:HE1	1:A:675:GLU:HG2	1.76	0.50
1:A:662:LEU:HB2	1:A:671:VAL:HB	1.92	0.50
1:A:7:PHE:HD1	1:A:412:ASN:HD21	1.59	0.50
1:A:323:ILE:CD1	1:A:374:LEU:HD12	2.36	0.50
1:A:502:ILE:HD12	1:A:502:ILE:H	1.77	0.50
1:A:595:PHE:HB2	1:A:677:LEU:HD12	1.93	0.50
1:A:638:VAL:HB	1:A:662:LEU:HD23	1.93	0.50
1:A:404:ILE:O	1:A:404:ILE:HG22	2.11	0.50
1:A:142:ASN:HB2	1:A:331:ARG:HA	1.94	0.50
1:A:634:LEU:HD13	1:A:661:LEU:HB3	1.93	0.49
1:A:147:ARG:HA	1:A:152:TYR:CE1	2.47	0.49
1:A:310:TYR:CZ	1:A:488:LYS:HE3	2.47	0.49
1:A:653:ILE:N	1:A:653:ILE:HD12	2.28	0.49
1:A:172:ARG:NH2	1:A:178:ASP:OD2	2.46	0.49
1:A:182:LEU:HD22	1:A:230:LEU:HD12	1.94	0.49
1:A:198:SER:C	1:A:200:TYR:N	2.65	0.49
1:A:225:TYR:HB2	1:A:266:ILE:HG21	1.94	0.49
1:A:451:LYS:O	1:A:456:ARG:NH2	2.46	0.49
1:A:468:TYR:HB2	1:A:469:PRO:CD	2.42	0.49
1:A:542:PHE:C	1:A:542:PHE:HD2	2.14	0.49
1:A:37:TRP:CZ2	1:A:434:LYS:HD3	2.48	0.49
1:A:166:ILE:HD11	1:A:189:ILE:HA	1.93	0.49
1:A:424:SER:HA	1:A:427:LEU:HD12	1.94	0.49
1:A:457:GLU:HG2	1:A:559:ARG:HH21	1.77	0.49
1:A:617:GLU:O	1:A:625:LEU:HA	2.13	0.49
1:A:175:ILE:HG22	1:A:176:ASN:ND2	2.28	0.49
1:A:329:MET:HG3	1:A:336:VAL:HG22	1.95	0.49
1:A:572:ILE:CG2	1:A:576:VAL:HG23	2.31	0.49
1:A:664:LEU:CD2	1:A:687:LEU:HD22	2.36	0.49
1:A:464:TRP:HB3	1:A:480:ILE:HG13	1.95	0.49
1:A:207:ILE:HG23	1:A:231:MET:CE	2.40	0.49
1:A:156:ASP:HA	1:A:159:VAL:CG2	2.43	0.49
1:A:272:ILE:O	1:A:275:SER:HB2	2.13	0.49
1:A:570:LEU:HB3	1:A:597:GLN:OE1	2.13	0.49
1:A:599:ASN:O	1:A:612:LEU:HD22	2.13	0.49
1:A:308:PHE:CZ	1:A:539:ALA:HB2	2.48	0.48
1:A:172:ARG:HH11	1:A:172:ARG:CG	2.26	0.48
1:A:653:ILE:O	1:A:654:ASP:C	2.51	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:LEU:HD12	1:A:113:LEU:HA	1.62	0.48
1:A:217:PHE:N	1:A:217:PHE:CD1	2.81	0.48
1:A:317:ILE:CD1	1:A:485:HIS:CE1	2.94	0.48
1:A:343:PHE:HA	1:A:384:MET:CE	2.43	0.48
1:A:508:ALA:HB3	1:A:710:LEU:HD11	1.95	0.48
1:A:169:THR:HG22	1:A:170:PHE:N	2.28	0.48
1:A:99:PHE:CE2	1:A:103:ILE:HD12	2.48	0.48
1:A:122:ALA:CB	1:A:131:GLY:HA3	2.44	0.48
1:A:125:TYR:OH	1:A:355:SER:HB2	2.13	0.48
1:A:223:ILE:HG12	1:A:254:PHE:CZ	2.48	0.48
1:A:349:LEU:HD23	1:A:356:MET:HG2	1.96	0.48
1:A:457:GLU:HG3	1:A:559:ARG:HE	1.78	0.48
1:A:638:VAL:HG22	1:A:650:TYR:HD1	1.79	0.48
1:A:329:MET:HG2	1:A:329:MET:H	1.50	0.48
1:A:667:TYR:OH	1:A:703:THR:HG22	2.14	0.48
1:A:51:MET:HE1	1:A:473:TYR:OH	2.14	0.48
1:A:322:THR:HG22	1:A:323:ILE:H	1.79	0.48
1:A:103:ILE:HG22	1:A:104:LEU:N	2.29	0.48
1:A:161:VAL:HG12	1:A:162:LEU:N	2.28	0.47
1:A:397:ASN:O	1:A:398:PHE:C	2.52	0.47
1:A:450:LEU:HD21	1:A:562:TYR:CE2	2.49	0.47
1:A:424:SER:OG	1:A:425:PRO:HD3	2.15	0.47
1:A:496:VAL:HG23	1:A:507:MET:HB3	1.96	0.47
1:A:686:PHE:CD1	1:A:706:LYS:HG3	2.48	0.47
1:A:143:SER:OG	1:A:331:ARG:HB3	2.13	0.47
1:A:437:THR:HG22	1:A:438:VAL:N	2.29	0.47
1:A:562:TYR:N	1:A:562:TYR:CD1	2.83	0.47
1:A:583:ASN:ND2	1:A:585:ASP:HB2	2.24	0.47
1:A:33:TYR:HD2	1:A:34:TRP:N	2.11	0.47
1:A:30:CYS:C	1:A:32:LEU:H	2.17	0.47
1:A:154:ASP:O	1:A:156:ASP:N	2.48	0.47
1:A:393:TYR:O	1:A:394:ALA:C	2.52	0.47
1:A:111:ALA:CB	1:A:148:THR:HG23	2.44	0.47
1:A:125:TYR:CE1	1:A:356:MET:CE	2.93	0.47
1:A:323:ILE:HG22	1:A:324:ASP:N	2.30	0.47
1:A:594:PHE:HD2	1:A:595:PHE:N	2.13	0.47
1:A:160:LEU:HD11	1:A:378:ILE:O	2.15	0.46
1:A:162:LEU:HD13	1:A:192:TYR:HA	1.97	0.46
1:A:482:GLY:C	1:A:485:HIS:HD2	2.19	0.46
1:A:595:PHE:HA	1:A:672:ILE:O	2.16	0.46
1:A:104:LEU:HD23	1:A:105:TYR:CD1	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:ILE:HG12	1:A:116:VAL:HG22	1.97	0.46
1:A:111:ALA:HB1	1:A:148:THR:HG21	1.96	0.46
1:A:460:VAL:HG11	1:A:470:ILE:CG2	2.40	0.46
1:A:87:LEU:HG	1:A:91:LEU:HD12	1.98	0.46
1:A:308:PHE:HE2	1:A:527:LEU:HD23	1.80	0.46
1:A:324:ASP:O	1:A:326:GLU:N	2.48	0.46
1:A:216:VAL:HG12	1:A:217:PHE:N	2.31	0.46
1:A:495:PHE:CD2	1:A:507:MET:HG2	2.51	0.46
1:A:636:ALA:HA	1:A:653:ILE:HD12	1.97	0.46
1:A:104:LEU:O	1:A:150:SER:O	2.33	0.46
1:A:94:ILE:O	1:A:94:ILE:HG22	2.16	0.46
1:A:444:ALA:O	1:A:447:LEU:N	2.50	0.45
1:A:58:TYR:HE2	1:A:473:TYR:CE1	2.34	0.45
1:A:144:TYR:HA	1:A:378:ILE:CD1	2.36	0.45
1:A:193:LEU:HD13	1:A:203:ASN:HD22	1.79	0.45
1:A:314:ASN:C	1:A:316:THR:H	2.19	0.45
1:A:566:PRO:HA	1:A:705:ALA:HA	1.97	0.45
1:A:27:SER:O	1:A:31:ARG:HG2	2.15	0.45
1:A:58:TYR:HA	1:A:61:ALA:HB3	1.99	0.45
1:A:246:LEU:CD1	1:A:250:LEU:HD12	2.44	0.45
1:A:440:THR:O	1:A:441:SER:C	2.53	0.45
1:A:496:VAL:HG22	1:A:508:ALA:H	1.81	0.45
1:A:363:LEU:HD21	1:A:384:MET:HG2	1.98	0.45
1:A:211:GLY:HA2	1:A:228:ILE:HD11	1.98	0.45
1:A:698:GLN:HE22	1:A:701:ASN:ND2	2.07	0.45
1:A:217:PHE:CE2	1:A:354:LYS:HD3	2.52	0.45
1:A:61:ALA:HB2	1:A:103:ILE:HD11	1.99	0.44
1:A:222:LYS:O	1:A:224:PHE:N	2.51	0.44
1:A:237:MET:CG	1:A:280:SER:HB3	2.36	0.44
1:A:333:SER:O	1:A:334:SER:HB2	2.16	0.44
1:A:495:PHE:CD2	1:A:507:MET:CG	3.00	0.44
1:A:99:PHE:CE2	1:A:103:ILE:CD1	3.01	0.44
1:A:323:ILE:CD1	1:A:331:ARG:HD2	2.47	0.44
1:A:636:ALA:HA	1:A:653:ILE:CD1	2.46	0.44
1:A:120:LEU:HB3	1:A:168:LEU:CD1	2.47	0.44
1:A:253:ILE:O	1:A:257:LYS:HB2	2.17	0.44
1:A:440:THR:OG1	1:A:443:GLU:HG3	2.17	0.44
1:A:498:SER:HA	1:A:680:SER:HB2	1.98	0.44
1:A:632:ILE:CG2	1:A:633:PRO:HD2	2.46	0.44
1:A:673:LEU:HD11	1:A:683:ILE:HD12	1.98	0.44
1:A:122:ALA:HB1	1:A:127:LEU:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:209:LEU:HD12	1:A:209:LEU:HA	1.82	0.44
1:A:328:PHE:O	1:A:329:MET:C	2.56	0.44
1:A:343:PHE:HA	1:A:384:MET:HE1	1.97	0.44
1:A:455:GLN:O	1:A:458:ASP:HB2	2.17	0.44
1:A:177:LYS:NZ	1:A:221:GLU:HG2	2.33	0.44
1:A:195:TRP:O	1:A:195:TRP:CG	2.70	0.44
1:A:437:THR:HG22	1:A:439:PHE:H	1.82	0.44
1:A:511:SER:O	1:A:515:THR:HB	2.17	0.44
1:A:640:ILE:HD13	1:A:687:LEU:O	2.17	0.44
1:A:52:ILE:HD11	1:A:153:TYR:CD2	2.53	0.44
1:A:61:ALA:O	1:A:62:GLU:C	2.55	0.44
1:A:219:ARG:HA	1:A:225:TYR:OH	2.18	0.44
1:A:232:ILE:HG22	1:A:273:SER:OG	2.18	0.44
1:A:549:LYS:HB2	1:A:549:LYS:HE3	1.76	0.44
1:A:661:LEU:HD12	1:A:661:LEU:C	2.38	0.44
1:A:160:LEU:HA	1:A:160:LEU:HD23	1.53	0.43
1:A:24:PHE:N	1:A:115:VAL:HG11	2.33	0.43
1:A:192:TYR:CE2	1:A:199:SER:HB2	2.54	0.43
1:A:163:PRO:HG3	1:A:192:TYR:CE2	2.53	0.43
1:A:179:ILE:HG22	1:A:251:PHE:HE1	1.81	0.43
1:A:316:THR:HG23	2:B:15:THR:HG23	1.98	0.43
1:A:566:PRO:HA	1:A:705:ALA:CB	2.48	0.43
1:A:160:LEU:HD13	1:A:382:PRO:HG3	2.00	0.43
1:A:456:ARG:HH11	1:A:456:ARG:CA	2.19	0.43
1:A:366:GLY:O	1:A:369:ALA:HB3	2.18	0.43
1:A:595:PHE:HD1	1:A:677:LEU:HD13	1.84	0.43
1:A:347:ILE:O	1:A:351:LYS:HG2	2.19	0.43
1:A:594:PHE:C	1:A:594:PHE:CD2	2.92	0.43
1:A:49:GLN:NE2	1:A:104:LEU:HD22	2.34	0.43
1:A:625:LEU:N	1:A:625:LEU:HD23	2.33	0.43
1:A:673:LEU:HD22	1:A:677:LEU:HB3	2.00	0.43
1:A:62:GLU:HG3	1:A:65:ARG:NH2	2.34	0.43
1:A:562:TYR:N	1:A:562:TYR:HD1	2.17	0.43
1:A:44:PHE:CZ	1:A:436:SER:HA	2.54	0.43
1:A:35:VAL:O	1:A:39:SER:OG	2.34	0.43
1:A:417:LEU:O	1:A:418:ILE:C	2.57	0.43
1:A:441:SER:O	1:A:442:TYR:C	2.58	0.43
1:A:378:ILE:HG22	1:A:379:TYR:N	2.33	0.42
1:A:76:ASP:HB3	1:A:477:LYS:HE2	2.00	0.42
1:A:310:TYR:OH	1:A:488:LYS:HE3	2.20	0.42
1:A:76:ASP:OD1	1:A:76:ASP:O	2.37	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:SER:O	1:A:99:PHE:C	2.57	0.42
1:A:51:MET:CE	1:A:473:TYR:OH	2.67	0.42
1:A:156:ASP:HA	1:A:159:VAL:HG23	2.00	0.42
1:A:516:GLU:HG2	1:A:559:ARG:HH12	1.85	0.42
1:A:222:LYS:O	1:A:223:ILE:C	2.57	0.42
1:A:658:GLN:O	1:A:658:GLN:HG2	2.20	0.42
1:A:327:VAL:HA	1:A:330:GLN:HB2	2.02	0.42
1:A:347:ILE:HD13	1:A:347:ILE:HA	1.87	0.42
1:A:424:SER:N	1:A:425:PRO:CD	2.82	0.42
1:A:684:GLN:NE2	1:A:688:LEU:HD12	2.34	0.42
1:A:543:LEU:C	1:A:545:SER:N	2.72	0.42
1:A:218:HIS:HA	1:A:221:GLU:OE1	2.19	0.42
1:A:640:ILE:HD13	1:A:688:LEU:HD23	2.01	0.42
1:A:155:THR:O	1:A:159:VAL:HG23	2.20	0.42
1:A:611:MET:O	1:A:612:LEU:HD23	2.20	0.42
1:A:44:PHE:CE2	1:A:436:SER:HB2	2.53	0.42
1:A:398:PHE:O	1:A:402:LYS:CB	2.68	0.42
1:A:340:ILE:HG12	1:A:340:ILE:H	1.54	0.41
1:A:457:GLU:O	1:A:559:ARG:HD3	2.20	0.41
1:A:60:PHE:O	1:A:61:ALA:C	2.59	0.41
1:A:165:LEU:HD23	1:A:165:LEU:HA	1.82	0.41
1:A:698:GLN:O	1:A:698:GLN:HG2	2.19	0.41
1:A:119:ILE:CG2	1:A:132:PHE:HA	2.51	0.41
1:A:183:LEU:HA	1:A:183:LEU:HD23	1.57	0.41
1:A:625:LEU:CD2	1:A:634:LEU:HD21	2.50	0.41
1:A:158:LEU:HD12	1:A:195:TRP:CE2	2.56	0.41
1:A:508:ALA:CB	1:A:710:LEU:HD11	2.51	0.41
1:A:20:ILE:CA	1:A:119:ILE:HD11	2.49	0.41
1:A:116:VAL:HG12	1:A:117:PRO:N	2.35	0.41
1:A:595:PHE:CD1	1:A:673:LEU:HD23	2.56	0.41
1:A:71:PHE:HD2	1:A:72:HIS:O	2.01	0.41
1:A:163:PRO:HG3	1:A:192:TYR:CZ	2.56	0.41
1:A:365:LEU:CD2	1:A:368:MET:HE2	2.49	0.41
1:A:486:LEU:O	1:A:490:ASN:ND2	2.53	0.41
1:A:493:SER:O	1:A:497:LEU:HD12	2.20	0.41
1:A:496:VAL:O	1:A:504:ALA:HB1	2.20	0.41
1:A:119:ILE:HG22	1:A:132:PHE:N	2.36	0.41
1:A:602:ALA:O	1:A:610:VAL:HA	2.20	0.41
1:A:594:PHE:HD2	1:A:594:PHE:C	2.24	0.41
1:A:14:LYS:HB2	1:A:14:LYS:HE3	1.95	0.41
1:A:67:MET:HE2	1:A:89:TYR:CG	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:671:VAL:HG12	1:A:672:ILE:N	2.36	0.41
1:A:698:GLN:NE2	1:A:701:ASN:ND2	2.67	0.41
1:A:127:LEU:HD11	1:A:390:TYR:CE1	2.56	0.41
1:A:574:PRO:HG3	1:A:595:PHE:CD2	2.56	0.41
1:A:49:GLN:HE21	1:A:104:LEU:HD22	1.85	0.40
1:A:437:THR:O	1:A:438:VAL:C	2.59	0.40
1:A:464:TRP:HB3	1:A:480:ILE:CD1	2.51	0.40
1:A:468:TYR:CB	1:A:469:PRO:CD	3.00	0.40
1:A:46:PHE:HB2	1:A:51:MET:HE3	2.02	0.40
1:A:56:ASP:O	1:A:59:ALA:HB3	2.21	0.40
1:A:115:VAL:HG22	1:A:135:ALA:HB1	2.02	0.40
1:A:391:PHE:HA	1:A:394:ALA:HB3	2.02	0.40
1:A:673:LEU:HD13	1:A:678:TYR:HA	2.02	0.40
1:A:77:LEU:HD11	1:A:484:LYS:HD2	2.03	0.40
1:A:87:LEU:O	1:A:91:LEU:HD12	2.21	0.40
1:A:349:LEU:HG	1:A:356:MET:HG2	2.04	0.40
1:A:118:ILE:CB	1:A:135:ALA:HB2	2.50	0.40
1:A:136:LEU:HB3	1:A:423:ILE:CD1	2.52	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	677/724 (94%)	533 (79%)	117 (17%)	27 (4%)	3 18
2	B	5/8 (62%)	3 (60%)	1 (20%)	1 (20%)	0 0
All	All	682/732 (93%)	536 (79%)	118 (17%)	28 (4%)	3 18

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	335	SER
1	A	549	LYS
1	A	590	GLU
1	A	103	ILE
1	A	142	ASN
1	A	155	THR
1	A	308	PHE
1	A	403	GLN
1	A	620	ASN
1	A	654	ASP
1	A	200	TYR
1	A	245	ALA
1	A	258	GLU
1	A	394	ALA
2	B	11	ASP
1	A	61	ALA
1	A	216	VAL
1	A	250	LEU
1	A	330	GLN
1	A	331	ARG
1	A	351	LYS
1	A	508	ALA
1	A	62	GLU
1	A	223	ILE
1	A	141	ALA
1	A	404	ILE
1	A	102	ILE
1	A	115	VAL

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	607/646 (94%)	515 (85%)	92 (15%)	3 11
2	B	4/4 (100%)	4 (100%)	0	100 100
All	All	611/650 (94%)	519 (85%)	92 (15%)	3 12

All (92) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ILE
1	A	20	ILE
1	A	21	LEU
1	A	29	LEU
1	A	32	LEU
1	A	33	TYR
1	A	51	MET
1	A	55	ASN
1	A	76	ASP
1	A	83	SER
1	A	88	THR
1	A	89	TYR
1	A	101	SER
1	A	103	ILE
1	A	107	SER
1	A	112	SER
1	A	113	LEU
1	A	114	ILE
1	A	115	VAL
1	A	146	ASN
1	A	147	ARG
1	A	152	TYR
1	A	155	THR
1	A	157	MET
1	A	161	VAL
1	A	167	LEU
1	A	168	LEU
1	A	169	THR
1	A	172	ARG
1	A	174	THR
1	A	183	LEU
1	A	189	ILE
1	A	196	TYR
1	A	214	THR
1	A	215	LEU
1	A	223	ILE
1	A	226	LEU
1	A	237	MET
1	A	244	LEU
1	A	246	LEU
1	A	248	VAL
1	A	273	SER

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Mol	Chain	Res	Type
1	A	277	LEU
1	A	309	MET
1	A	313	VAL
1	A	316	THR
1	A	317	ILE
1	A	320	VAL
1	A	322	THR
1	A	323	ILE
1	A	326	GLU
1	A	327	VAL
1	A	329	MET
1	A	331	ARG
1	A	333	SER
1	A	337	LEU
1	A	340	ILE
1	A	352	ASP
1	A	355	SER
1	A	363	LEU
1	A	374	LEU
1	A	375	ARG
1	A	377	THR
1	A	378	ILE
1	A	397	ASN
1	A	398	PHE
1	A	406	LEU
1	A	408	LEU
1	A	414	LEU
1	A	417	LEU
1	A	418	ILE
1	A	423	ILE
1	A	435	SER
1	A	441	SER
1	A	442	TYR
1	A	450	LEU
1	A	456	ARG
1	A	460	VAL
1	A	488	LYS
1	A	515	THR
1	A	542	PHE
1	A	562	TYR
1	A	583	ASN
1	A	593	LEU

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Mol	Chain	Res	Type
1	A	594	PHE
1	A	596	SER
1	A	609	SER
1	A	644	THR
1	A	654	ASP
1	A	661	LEU
1	A	697	GLU
1	A	700	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	142	ASN
1	A	146	ASN
1	A	176	ASN
1	A	203	ASN
1	A	353	HIS
1	A	410	ASN
1	A	485	HIS
1	A	583	ASN
1	A	645	ASN
1	A	698	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)

1 non-standard protein/DNA/RNA residue is 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	PPN	B	16	2	13,14,15	0.98	1 (7%)	11,18,20	1.16	1 (9%)

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	PPN	B	16	2	-	1/7/10/12	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	16	PPN	CB-CA	-2.51	1.48	1.53

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	16	PPN	CE1-CZ-N1	3.15	122.09	119.34

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	16	PPN	C-CA-CB-CG

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	16	PPN	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	683/724 (94%)	-0.12	24 (3%) 44 43	73, 119, 191, 321	0
2	B	7/8 (87%)	-0.03	0 100 100	88, 113, 133, 140	2 (28%)
All	All	690/732 (94%)	-0.12	24 (3%) 44 43	73, 119, 191, 321	2 (0%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	259	GLU	5.7
1	A	493	SER	4.6
1	A	622	PHE	3.7
1	A	258	GLU	3.2
1	A	537	THR	3.1
1	A	11	ASN	3.1
1	A	532	LYS	3.0
1	A	220	LYS	3.0
1	A	536	GLN	2.8
1	A	405	LYS	2.7
1	A	403	GLN	2.7
1	A	621	ASP	2.6
1	A	494	SER	2.4
1	A	538	SER	2.3
1	A	321	ASN	2.3
1	A	520	LYS	2.3
1	A	462	ALA	2.2
1	A	521	GLU	2.2
1	A	693	GLN	2.2
1	A	623	ARG	2.2
1	A	322	THR	2.1
1	A	260	LYS	2.1
1	A	13	ILE	2.0
1	A	2	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	PPN	B	16	14/15	0.90	0.28	111,118,131,135	0

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	MG	A	725	1/1	0.96	0.76	78,78,78,78	0
3	MG	A	726	1/1	0.98	0.18	65,65,65,65	0

6.5 Other polymers (i)

There are no such residues in this entry.