



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:41 PM UTC

PDB ID : 7QYH / pdb_00007qyh
Title : Structure of plasmepsin II in complex with 2-aminoquinazolin-4(3H)-one based open-flap inhibitor
Authors : Bobrovs, R.; Jaudzems, K.
Deposited on : 2022-01-28
Resolution : 3.33 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

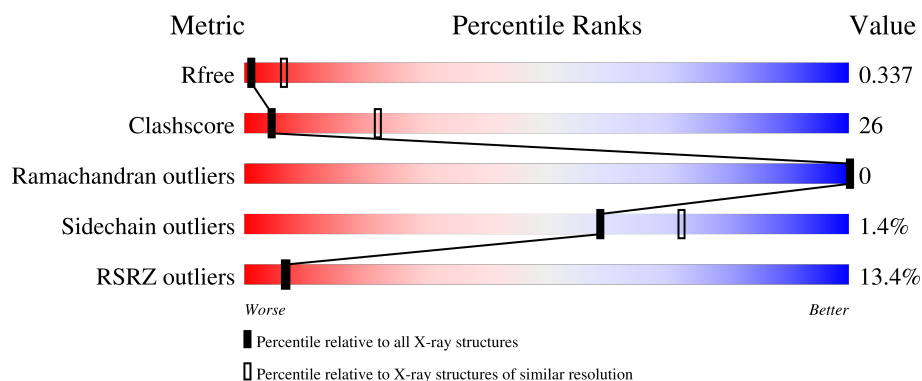
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.33 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}	180053	1434 (3.38-3.30)
Clashscore	190562	1479 (3.38-3.30)
Ramachandran outliers	187476	1456 (3.38-3.30)
Sidechain outliers	187428	1455 (3.38-3.30)
RSRZ outliers	180081	1434 (3.38-3.30)

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 $\leq 5\%$. 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	329	11% 51% 49%
1	B	329	13% 58% 42% .
1	C	329	14% 68% 31% .
1	D	329	16% 60% 40% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10478 atoms, of which 29 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.

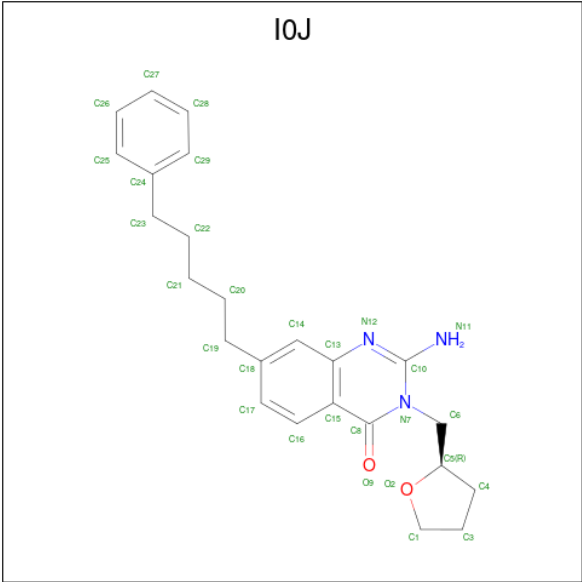
- Molecule 1 is a protein called Plasmepsin II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	0	0	0
			2605	1687	404	504	10			
1	B	329	Total	C	N	O	S	0	0	0
			2605	1687	404	504	10			
1	C	329	Total	C	N	O	S	0	0	0
			2605	1687	404	504	10			
1	D	329	Total	C	N	O	S	0	0	0
			2605	1687	404	504	10			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	205	SER	MET	conflict	UNP Q8I6V3
A	318	HIS	GLN	variant	UNP Q8I6V3
B	205	SER	MET	conflict	UNP Q8I6V3
B	318	HIS	GLN	variant	UNP Q8I6V3
C	205	SER	MET	conflict	UNP Q8I6V3
C	318	HIS	GLN	variant	UNP Q8I6V3
D	205	SER	MET	conflict	UNP Q8I6V3
D	318	HIS	GLN	variant	UNP Q8I6V3

- Molecule 2 is 2-azanyl-3-[[(2 {R})-oxolan-2-yl]methyl]-7-(5-phenylpentyl)quinazolin-4-one (CCD ID: I0J) (formula: C₂₄H₂₉N₃O₂) (labeled as "Ligand of Interest" by depositor).

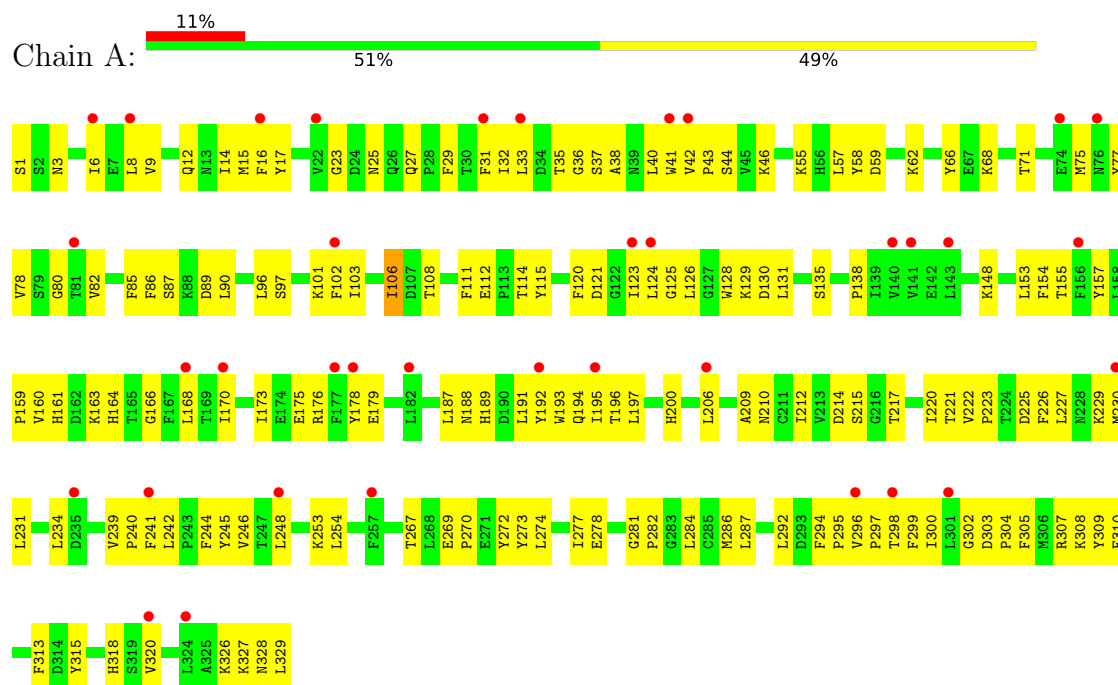


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	0	0
			58	24	29	3	2		

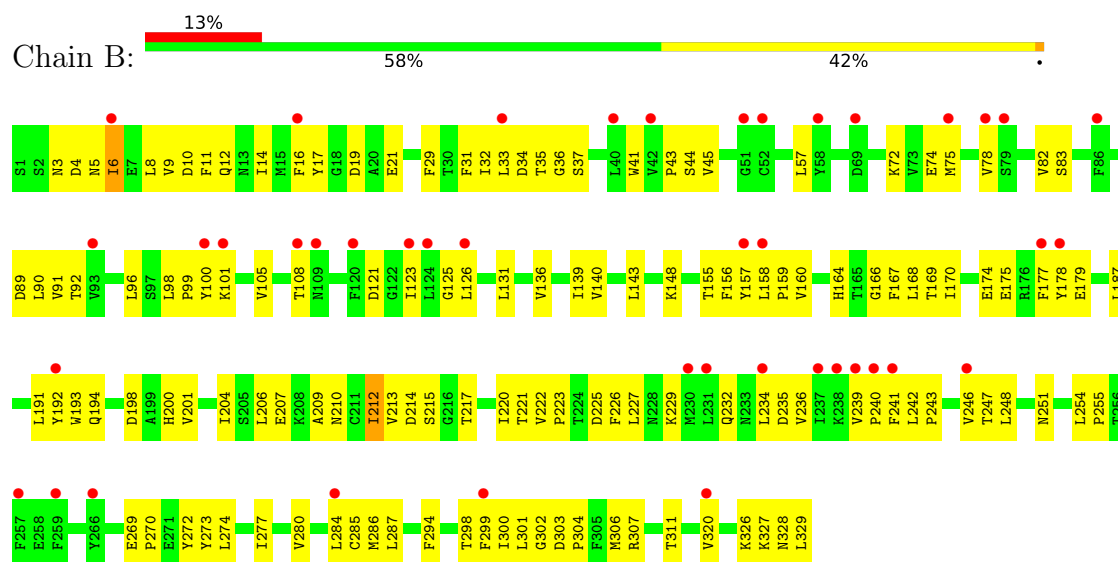
3 Residue-property plots [i](#)

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.

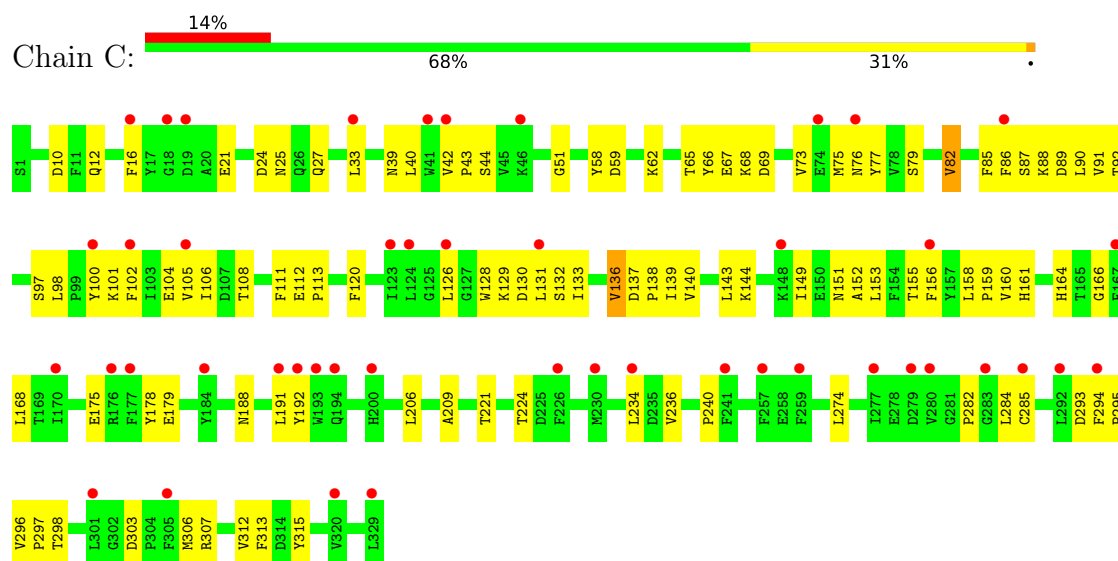
• Molecule 1: Plasmepsin II



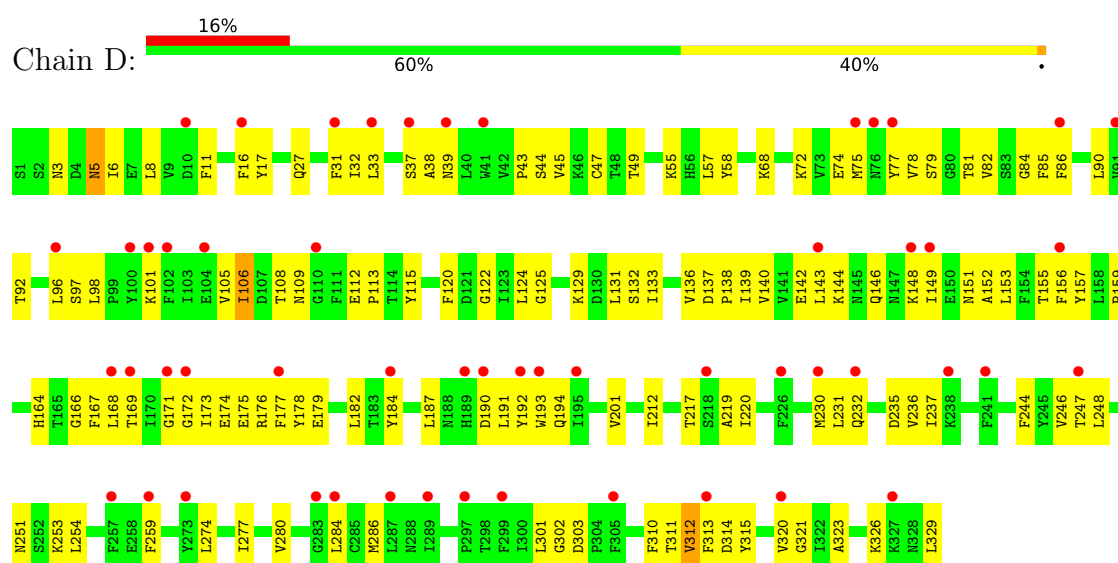
• Molecule 1: Plasmepsin II



● Molecule 1: Plasmepsin II



● Molecule 1: Plasmepsin II



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	70.82Å 274.07Å 273.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	137.03 – 3.33 137.03 – 3.33	Depositor EDS
% Data completeness (in resolution range)	38.6 (137.03-3.33) 38.6 (137.03-3.33)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 3.33Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.310 , 0.336 0.309 , 0.337	Depositor DCC
R_{free} test set	834 reflections (2.10%)	wwPDB-VP
Wilson B-factor (Å ²)	68.3	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 82.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.80	EDS
Total number of atoms	10478	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.69% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: I0J

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	0.93	0/2672	1.18	3/3636 (0.1%)
1	B	0.93	0/2672	1.18	2/3636 (0.1%)
1	C	0.93	1/2672 (0.0%)	1.28	2/3636 (0.1%)
1	D	0.95	1/2672 (0.0%)	1.25	2/3636 (0.1%)
All	All	0.94	2/10688 (0.0%)	1.22	9/14544 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	303	ASP	CG-OD1	5.94	1.36	1.25
1	D	303	ASP	CG-OD1	5.85	1.36	1.25

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	23	GLY	CA-C-O	-5.80	118.23	122.23
1	C	120	PHE	CA-CB-CG	5.78	119.58	113.80
1	D	303	ASP	CA-CB-CG	5.38	117.98	112.60
1	A	106	ILE	N-CA-C	-5.33	108.31	113.53
1	B	251	ASN	CA-CB-CG	5.22	117.82	112.60
1	C	303	ASP	CA-CB-CG	5.17	117.77	112.60
1	B	328	ASN	CA-CB-CG	5.10	117.70	112.60
1	D	106	ILE	N-CA-C	-5.07	108.56	113.53
1	A	200	HIS	CE1-NE2-CD2	-5.04	103.96	109.00

There are no chirality outliers.

There are no planarity outliers.

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	2605	0	2532	194	0
1	B	2605	0	2532	131	0
1	C	2605	0	2532	92	0
1	D	2605	0	2532	135	0
2	A	29	29	0	1	0
All	All	10449	29	10128	542	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 26.

All (542) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:PHE:HZ	1:B:215:SER:HB2	1.35	0.89
1:D:5:ASN:HB3	1:D:167:PHE:HB3	1.57	0.86
1:B:239:VAL:HG11	1:B:246:VAL:HG12	1.57	0.86
1:B:10:ASP:HB2	1:B:159:PRO:HG2	1.56	0.83
1:B:204:ILE:HD12	1:B:229:LYS:HZ3	1.44	0.83
1:D:153:LEU:HG	1:D:171:GLY:HA2	1.61	0.82
1:A:43:PRO:HB2	1:A:57:LEU:HD23	1.60	0.82
1:B:10:ASP:HA	1:B:16:PHE:HA	1.60	0.82
1:A:214:ASP:OD2	1:A:217:THR:HG23	1.79	0.81
1:A:29:PHE:HB2	1:A:31:PHE:CE2	2.15	0.81
1:A:246:VAL:HG13	1:A:277:ILE:HD11	1.61	0.81
1:A:71:THR:HB	1:A:86:PHE:HD2	1.46	0.81
1:A:46:LYS:HB2	1:A:106:ILE:HD12	1.62	0.81
1:A:80:GLY:HA3	1:A:111:PHE:CZ	2.16	0.80
1:A:277:ILE:HD13	1:A:286:MET:HB2	1.64	0.79
1:B:159:PRO:HA	1:B:164:HIS:CD2	2.17	0.79
1:B:198:ASP:HB3	1:B:200:HIS:HE1	1.48	0.79
1:D:129:LYS:HB2	1:D:138:PRO:HD3	1.64	0.79
1:A:240:PRO:HG2	1:C:284:LEU:HD21	1.65	0.78
1:A:195:ILE:HD11	1:A:197:LEU:HD11	1.65	0.78
1:A:17:TYR:CE2	1:A:32:ILE:HD11	2.19	0.78
1:B:198:ASP:HB3	1:B:200:HIS:CE1	2.18	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:27:GLN:HE22	1:D:58:TYR:HA	1.48	0.78
1:A:194:GLN:HB2	1:A:210:ASN:HD21	1.49	0.78
1:D:140:VAL:HA	1:D:143:LEU:HD12	1.65	0.77
1:C:90:LEU:HD11	1:C:97:SER:HB3	1.66	0.76
1:D:8:LEU:HB2	1:D:166:GLY:HA3	1.65	0.75
1:D:232:GLN:HA	1:D:236:VAL:HG21	1.68	0.75
1:A:189:HIS:HB3	1:A:194:GLN:NE2	2.01	0.75
1:D:182:LEU:HD21	1:D:312:VAL:HG11	1.67	0.75
1:C:44:SER:HB2	1:C:104:GLU:HB3	1.68	0.75
1:A:25:ASN:HD21	1:A:27:GLN:HE21	1.34	0.74
1:A:38:ALA:O	1:A:125:GLY:HA3	1.88	0.74
1:B:101:LYS:HD3	1:B:136:VAL:HG22	1.70	0.74
1:A:226:PHE:HA	1:A:229:LYS:HE2	1.69	0.74
1:D:90:LEU:HD11	1:D:97:SER:HB3	1.69	0.74
1:D:237:ILE:HB	1:D:247:THR:HA	1.70	0.74
1:B:8:LEU:HB2	1:B:159:PRO:HD3	1.71	0.73
1:D:184:TYR:HA	1:D:321:GLY:HA2	1.69	0.73
1:B:234:LEU:HD11	1:B:255:PRO:HD3	1.69	0.73
1:A:159:PRO:HB3	1:A:166:GLY:N	2.04	0.72
1:D:6:ILE:HB	1:D:168:LEU:HB3	1.70	0.72
1:A:278:GLU:HB3	1:A:282:PRO:HA	1.71	0.72
1:A:192:TYR:O	1:A:194:GLN:HG2	1.89	0.72
1:D:159:PRO:HB3	1:D:166:GLY:H	1.54	0.72
1:A:8:LEU:HB2	1:A:166:GLY:HA3	1.71	0.72
1:D:277:ILE:HD13	1:D:286:MET:HE3	1.70	0.71
1:C:77:TYR:HB3	1:C:111:PHE:HZ	1.55	0.71
1:C:88:LYS:HA	1:C:101:LYS:HA	1.71	0.71
1:A:75:MET:HE3	1:A:75:MET:O	1.91	0.71
1:A:159:PRO:HA	1:A:164:HIS:ND1	2.05	0.71
1:B:16:PHE:CZ	1:B:215:SER:HB2	2.23	0.71
1:C:66:TYR:CE1	1:C:87:SER:HB3	2.26	0.71
1:A:231:LEU:HD23	1:A:234:LEU:HD12	1.71	0.71
1:D:248:LEU:H	1:D:248:LEU:HD12	1.56	0.70
1:C:12:GLN:H	1:C:161:HIS:CE1	2.10	0.70
1:A:189:HIS:CE1	1:A:191:LEU:HB2	2.25	0.70
1:B:206:LEU:HD22	1:B:209:ALA:HB2	1.73	0.70
1:B:159:PRO:HA	1:B:164:HIS:HD2	1.55	0.69
1:A:221:THR:HG22	1:A:222:VAL:H	1.58	0.69
1:D:37:SER:O	1:D:125:GLY:HA2	1.94	0.68
1:C:66:TYR:HE1	1:C:87:SER:HB3	1.58	0.68
1:B:33:LEU:HB3	1:B:126:LEU:HD11	1.74	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:5:ASN:HB3	1:B:167:PHE:HB3	1.76	0.68
1:D:159:PRO:HB3	1:D:166:GLY:N	2.09	0.68
1:A:187:LEU:HD21	1:A:320:VAL:HG23	1.76	0.67
1:B:43:PRO:HG3	1:B:121:ASP:HB3	1.75	0.67
1:C:76:ASN:HB3	1:D:75:MET:HE3	1.77	0.67
1:A:35:THR:OG1	1:A:214:ASP:HA	1.95	0.67
1:A:294:PHE:HB3	1:A:295:PRO:HD2	1.76	0.67
1:A:192:TYR:HE2	1:A:212:ILE:HD13	1.60	0.67
1:A:192:TYR:CE2	1:A:212:ILE:HD13	2.30	0.66
1:D:68:LYS:HG3	1:D:85:PHE:HE2	1.60	0.66
1:D:45:VAL:HA	1:D:57:LEU:HD13	1.78	0.66
1:A:35:THR:HG23	1:A:215:SER:H	1.60	0.66
1:C:98:LEU:HD11	1:C:139:ILE:HG23	1.77	0.66
1:D:173:ILE:HG23	1:D:178:TYR:HE2	1.60	0.66
1:C:39:ASN:HD21	1:C:132:SER:HB2	1.61	0.66
1:A:239:VAL:HG12	1:A:242:LEU:HB2	1.76	0.65
1:B:16:PHE:HE2	1:B:33:LEU:HD12	1.61	0.65
1:D:143:LEU:HD13	1:D:149:ILE:HD13	1.77	0.65
1:A:191:LEU:HD22	1:A:192:TYR:HD1	1.60	0.65
1:D:27:GLN:NE2	1:D:58:TYR:HA	2.11	0.65
1:B:35:THR:HA	1:B:126:LEU:HB2	1.78	0.65
1:A:66:TYR:HE1	1:A:87:SER:HB3	1.62	0.65
1:A:15:MET:CG	1:A:32:ILE:HD13	2.27	0.64
1:A:130:ASP:OD2	1:A:191:LEU:HA	1.97	0.64
1:A:303:ASP:O	1:A:307:ARG:HG3	1.98	0.64
1:B:96:LEU:HD13	1:B:143:LEU:HD13	1.78	0.64
1:B:226:PHE:HA	1:B:229:LYS:HZ2	1.63	0.64
1:C:144:LYS:HD2	1:C:152:ALA:HB2	1.80	0.64
1:A:157:TYR:O	1:A:166:GLY:HA2	1.98	0.64
1:B:254:LEU:HD13	1:B:274:LEU:HD11	1.79	0.64
1:D:133:ILE:HD12	1:D:133:ILE:H	1.63	0.63
1:B:155:THR:O	1:B:168:LEU:HD12	1.98	0.63
1:B:247:THR:HG22	1:B:287:LEU:HD11	1.79	0.63
1:A:25:ASN:OD1	1:A:27:GLN:HG3	1.99	0.63
1:B:43:PRO:HB2	1:B:57:LEU:HD23	1.81	0.62
1:B:232:GLN:HA	1:B:236:VAL:HA	1.82	0.62
1:A:195:ILE:CD1	1:A:197:LEU:HD11	2.29	0.62
1:A:214:ASP:HB3	1:A:217:THR:OG1	2.00	0.62
1:C:27:GLN:NE2	1:C:58:TYR:HA	2.15	0.61
1:B:270:PRO:HA	1:B:273:TYR:CZ	2.35	0.61
1:A:128:TRP:CH2	1:A:315:TYR:HA	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:LEU:HB3	1:A:244:PHE:CD2	2.35	0.61
1:C:75:MET:HG3	1:C:82:VAL:HG13	1.81	0.61
1:D:44:SER:HA	1:D:58:TYR:H	1.66	0.61
1:B:226:PHE:HA	1:B:229:LYS:NZ	2.16	0.61
1:D:173:ILE:HG23	1:D:178:TYR:CE2	2.35	0.61
1:B:139:ILE:HD12	1:B:139:ILE:H	1.65	0.60
1:A:191:LEU:C	1:A:191:LEU:HD23	2.26	0.60
1:A:27:GLN:OE1	1:A:58:TYR:HA	2.01	0.60
1:D:182:LEU:HA	1:D:323:ALA:HB2	1.84	0.60
1:A:43:PRO:HB2	1:A:57:LEU:CD2	2.31	0.60
1:B:280:VAL:HG12	1:B:284:LEU:HD22	1.84	0.60
1:D:39:ASN:ND2	1:D:132:SER:HB2	2.16	0.60
1:B:175:GLU:HA	1:B:178:TYR:CE1	2.37	0.60
1:A:278:GLU:HA	1:A:281:GLY:O	2.01	0.60
1:A:55:LYS:HG3	1:A:115:TYR:CE2	2.37	0.60
1:B:3:ASN:HB2	1:B:169:THR:HG23	1.84	0.60
1:D:16:PHE:HE2	1:D:33:LEU:HD12	1.66	0.60
1:D:86:PHE:CD1	1:D:101:LYS:HE2	2.36	0.60
1:A:227:LEU:HD12	1:A:230:MET:HE2	1.82	0.59
1:B:140:VAL:HA	1:B:143:LEU:HD12	1.83	0.59
1:A:128:TRP:HB2	1:A:191:LEU:O	2.01	0.59
1:A:129:LYS:HB2	1:A:138:PRO:HD3	1.83	0.59
1:A:124:LEU:CD2	1:A:126:LEU:HD11	2.32	0.59
1:A:33:LEU:HA	1:A:124:LEU:HB3	1.85	0.59
1:D:72:LYS:HE2	1:D:74:GLU:HB2	1.84	0.59
1:D:84:GLY:HA2	1:D:106:ILE:HG12	1.83	0.59
1:D:6:ILE:H	1:D:6:ILE:HD12	1.68	0.59
1:A:128:TRP:CE3	1:A:315:TYR:HD1	2.21	0.59
1:D:3:ASN:HB3	1:D:169:THR:HG23	1.85	0.59
1:A:12:GLN:HB2	1:A:14:ILE:HG22	1.84	0.59
1:A:17:TYR:CZ	1:A:32:ILE:HD11	2.38	0.58
1:B:36:GLY:HA2	1:B:192:TYR:HE1	1.68	0.58
1:D:43:PRO:HA	1:D:105:VAL:HB	1.85	0.58
1:A:130:ASP:HB3	1:A:191:LEU:HG	1.85	0.58
1:A:209:ALA:HB1	1:A:299:PHE:CD1	2.39	0.58
1:A:220:ILE:HD11	1:A:304:PRO:HB2	1.86	0.58
1:D:230:MET:C	1:D:232:GLN:H	2.12	0.58
1:A:71:THR:HG21	1:B:72:LYS:HE2	1.86	0.58
1:B:131:LEU:HD13	1:B:191:LEU:HD21	1.86	0.58
1:C:144:LYS:HE2	1:C:151:ASN:HA	1.85	0.58
1:D:47:CYS:HA	1:D:106:ILE:HA	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:15:MET:SD	1:A:32:ILE:HG21	2.44	0.57
1:A:124:LEU:HD21	1:A:126:LEU:HD11	1.84	0.57
1:A:124:LEU:HD11	1:A:126:LEU:HD21	1.86	0.57
1:C:156:PHE:CE1	1:C:168:LEU:HD13	2.39	0.57
1:A:29:PHE:HB2	1:A:31:PHE:CZ	2.40	0.57
1:C:159:PRO:HB3	1:C:166:GLY:N	2.19	0.57
1:B:35:THR:OG1	1:B:214:ASP:HA	2.04	0.57
1:A:124:LEU:HG	1:A:126:LEU:HG	1.86	0.57
1:D:248:LEU:HB2	1:D:251:ASN:HB2	1.87	0.57
1:C:77:TYR:HB3	1:C:111:PHE:CZ	2.39	0.57
1:B:98:LEU:HD23	1:B:98:LEU:H	1.70	0.56
1:A:75:MET:HE1	1:A:77:TYR:HB3	1.87	0.56
1:D:77:TYR:CD2	1:D:78:VAL:HG12	2.40	0.56
1:D:98:LEU:HD11	1:D:148:LYS:HD2	1.87	0.56
1:A:155:THR:O	1:A:168:LEU:HD12	2.05	0.56
1:A:234:LEU:HD22	1:A:253:LYS:HB3	1.86	0.56
1:A:287:LEU:HD12	1:A:287:LEU:H	1.70	0.56
1:A:25:ASN:HD21	1:A:27:GLN:NE2	2.02	0.56
1:D:49:THR:HB	1:D:112:GLU:CD	2.31	0.56
1:C:89:ASP:HB3	1:C:102:PHE:HE1	1.71	0.56
1:A:254:LEU:HD13	1:A:274:LEU:HD11	1.88	0.56
1:D:231:LEU:O	1:D:253:LYS:HD3	2.06	0.56
1:C:132:SER:HB2	1:C:136:VAL:HB	1.87	0.56
1:B:43:PRO:HA	1:B:105:VAL:HG22	1.88	0.56
1:D:68:LYS:HG3	1:D:85:PHE:CE2	2.39	0.55
1:B:242:LEU:HD12	1:B:243:PRO:HD2	1.88	0.55
1:C:24:ASP:HB2	1:C:65:THR:HG23	1.88	0.55
1:B:98:LEU:HD22	1:B:143:LEU:HD21	1.88	0.55
1:A:86:PHE:CG	1:A:101:LYS:HE2	2.42	0.55
1:A:8:LEU:HD22	1:A:33:LEU:HD11	1.89	0.55
1:C:39:ASN:ND2	1:C:132:SER:HB2	2.21	0.55
1:A:130:ASP:HB2	1:A:191:LEU:HA	1.88	0.54
1:A:123:ILE:HG22	1:A:124:LEU:H	1.71	0.54
1:A:75:MET:HE1	1:A:77:TYR:H	1.72	0.54
1:C:126:LEU:HD13	1:C:156:PHE:HZ	1.71	0.54
1:D:140:VAL:HG12	1:D:152:ALA:HA	1.88	0.54
1:D:11:PHE:HB3	1:D:17:TYR:HE1	1.72	0.54
1:A:55:LYS:HB3	1:A:121:ASP:OD1	2.08	0.54
1:D:231:LEU:HB3	1:D:253:LYS:HG3	1.88	0.54
1:A:33:LEU:HA	1:A:124:LEU:HD23	1.89	0.54
1:C:40:LEU:HD23	1:C:102:PHE:HB3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:155:THR:HG22	1:C:312:VAL:HG13	1.89	0.54
1:D:176:ARG:HH12	1:D:329:LEU:HD13	1.73	0.54
1:A:129:LYS:HE3	1:A:135:SER:C	2.33	0.53
1:A:155:THR:HB	1:A:310:PHE:CZ	2.43	0.53
1:A:175:GLU:HA	1:A:178:TYR:CE1	2.43	0.53
1:B:90:LEU:HA	1:B:99:PRO:HA	1.90	0.53
1:C:153:LEU:HD12	1:C:153:LEU:O	2.08	0.53
1:D:155:THR:OG1	1:D:169:THR:HB	2.09	0.53
1:A:206:LEU:HD22	1:A:299:PHE:HE1	1.74	0.53
1:D:39:ASN:OD1	1:D:133:ILE:HB	2.08	0.53
1:A:123:ILE:HG22	1:A:124:LEU:N	2.23	0.53
1:A:220:ILE:HD11	1:A:305:PHE:H	1.74	0.53
1:B:83:SER:H	1:B:108:THR:HG22	1.73	0.53
1:B:98:LEU:HD21	1:B:139:ILE:HG23	1.89	0.53
1:A:187:LEU:HG	1:A:318:HIS:O	2.09	0.53
1:D:97:SER:O	1:D:98:LEU:HD13	2.08	0.53
1:A:192:TYR:CE2	1:A:194:GLN:HB3	2.43	0.53
1:A:55:LYS:HB3	1:A:121:ASP:CG	2.33	0.53
1:A:96:LEU:HD22	1:A:148:LYS:HB3	1.91	0.53
1:A:163:LYS:HB3	1:A:328:ASN:HD21	1.74	0.52
1:B:91:VAL:HG21	1:B:100:TYR:HB3	1.91	0.52
1:B:43:PRO:HA	1:B:105:VAL:CG2	2.40	0.52
1:B:74:GLU:HG3	1:B:82:VAL:O	2.09	0.52
1:A:77:TYR:CE1	1:A:78:VAL:HG12	2.43	0.52
1:B:3:ASN:HB3	1:B:170:ILE:O	2.09	0.52
1:B:29:PHE:HB2	1:B:31:PHE:CE2	2.45	0.52
1:D:253:LYS:O	1:D:254:LEU:C	2.52	0.52
1:A:173:ILE:HG22	1:A:178:TYR:CE1	2.44	0.52
1:D:8:LEU:HD21	1:D:168:LEU:HB2	1.91	0.52
1:D:259:PHE:CZ	1:D:301:LEU:HD21	2.45	0.52
1:A:281:GLY:HA3	1:A:284:LEU:HD12	1.92	0.52
1:B:82:VAL:HG23	1:B:108:THR:CG2	2.40	0.52
1:C:98:LEU:H	1:C:98:LEU:HD23	1.75	0.52
1:C:160:VAL:HB	1:C:164:HIS:H	1.74	0.52
1:B:33:LEU:HD13	1:B:168:LEU:HD22	1.91	0.52
1:A:221:THR:HG22	1:A:222:VAL:N	2.24	0.52
1:A:221:THR:OG1	1:A:300:ILE:HD12	2.10	0.52
1:C:88:LYS:HA	1:C:101:LYS:CA	2.40	0.52
1:D:5:ASN:N	1:D:5:ASN:HD22	2.07	0.51
1:A:115:TYR:OH	1:A:121:ASP:HA	2.10	0.51
1:B:5:ASN:OD1	1:B:169:THR:HA	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:246:VAL:HG22	1:D:286:MET:HE2	1.92	0.51
1:D:43:PRO:HB2	1:D:57:LEU:HD23	1.92	0.51
1:B:8:LEU:HB2	1:B:166:GLY:HA3	1.92	0.51
1:B:306:MET:HG2	1:B:311:THR:HG21	1.93	0.51
1:D:112:GLU:HB3	1:D:113:PRO:HD3	1.91	0.51
1:D:17:TYR:HD2	1:D:32:ILE:HG12	1.75	0.51
1:A:209:ALA:HB1	1:A:299:PHE:HD1	1.75	0.51
1:C:67:GLU:HB3	1:C:88:LYS:HB3	1.93	0.51
1:C:21:GLU:HB2	1:C:92:THR:HB	1.93	0.51
1:C:294:PHE:CD1	1:C:298:THR:HB	2.46	0.51
1:A:130:ASP:CB	1:A:191:LEU:HA	2.40	0.51
1:B:36:GLY:HA2	1:B:192:TYR:CE1	2.46	0.51
1:D:251:ASN:HB3	1:D:254:LEU:HG	1.93	0.51
1:D:153:LEU:CG	1:D:171:GLY:HA2	2.37	0.51
1:D:194:GLN:CB	1:D:212:ILE:HG12	2.41	0.51
1:A:242:LEU:HB3	1:A:244:PHE:HD2	1.74	0.50
1:B:83:SER:O	1:B:108:THR:HG23	2.11	0.50
1:A:226:PHE:HA	1:A:229:LYS:CE	2.40	0.50
1:B:11:PHE:O	1:B:14:ILE:HG22	2.11	0.50
1:A:128:TRP:HD1	1:A:192:TYR:HA	1.76	0.50
1:A:270:PRO:HA	1:A:273:TYR:CE2	2.47	0.50
1:B:75:MET:HB3	1:B:82:VAL:CG1	2.42	0.50
1:C:66:TYR:HE2	1:C:68:LYS:HG2	1.77	0.50
1:C:175:GLU:HA	1:C:178:TYR:CZ	2.46	0.50
1:A:44:SER:HB3	1:A:106:ILE:HD13	1.92	0.50
1:A:115:TYR:CE1	1:A:121:ASP:HA	2.46	0.50
1:A:239:VAL:CG1	1:A:242:LEU:HB2	2.42	0.50
1:A:15:MET:HG2	1:A:32:ILE:HD13	1.93	0.50
1:D:244:PHE:HB3	1:D:286:MET:SD	2.51	0.50
1:B:270:PRO:HA	1:B:273:TYR:CE2	2.47	0.50
1:C:130:ASP:HB2	1:C:191:LEU:HG	1.94	0.50
1:B:37:SER:OG	1:B:125:GLY:HA2	2.11	0.50
1:B:83:SER:N	1:B:108:THR:HG22	2.27	0.50
1:B:179:GLU:CD	1:B:326:LYS:HB3	2.37	0.50
1:D:97:SER:H	1:D:148:LYS:HE2	1.77	0.50
1:D:96:LEU:HD22	1:D:148:LYS:HD3	1.93	0.49
1:D:98:LEU:N	1:D:98:LEU:HD22	2.27	0.49
1:B:34:ASP:C	1:B:36:GLY:H	2.20	0.49
1:C:221:THR:HG21	1:C:294:PHE:CE1	2.46	0.49
1:A:160:VAL:HG13	1:A:307:ARG:HD3	1.94	0.49
1:C:108:THR:OG1	1:C:112:GLU:HB2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:174:GLU:HB3	1:B:177:PHE:HD2	1.76	0.49
1:C:88:LYS:CA	1:C:101:LYS:HA	2.40	0.49
1:B:126:LEU:HD13	1:B:156:PHE:HZ	1.78	0.49
1:C:43:PRO:HA	1:C:105:VAL:HB	1.95	0.49
1:C:295:PRO:HG2	1:D:79:SER:HB3	1.95	0.49
1:D:193:TRP:HZ2	1:D:315:TYR:CZ	2.30	0.49
1:B:187:LEU:HD21	1:B:320:VAL:HG23	1.94	0.49
1:B:246:VAL:HG23	1:B:285:CYS:O	2.13	0.49
1:A:75:MET:HE2	2:A:401:I0J:C19	2.43	0.49
1:A:220:ILE:CG1	1:A:304:PRO:HB2	2.43	0.49
1:C:133:ILE:HG23	1:D:72:LYS:NZ	2.28	0.49
1:A:66:TYR:CE1	1:A:87:SER:HB3	2.45	0.48
1:B:246:VAL:HA	1:B:286:MET:HA	1.95	0.48
1:C:27:GLN:HE22	1:C:58:TYR:HA	1.77	0.48
1:D:192:TYR:O	1:D:212:ILE:HG23	2.13	0.48
1:B:34:ASP:C	1:B:36:GLY:N	2.71	0.48
1:B:160:VAL:H	1:B:164:HIS:CG	2.32	0.48
1:A:1:SER:C	1:A:3:ASN:H	2.20	0.48
1:A:8:LEU:H	1:A:166:GLY:C	2.22	0.48
1:B:240:PRO:O	1:B:241:PHE:C	2.56	0.48
1:D:173:ILE:HG22	1:D:175:GLU:HG2	1.95	0.48
1:B:194:GLN:HB3	1:B:212:ILE:HG23	1.96	0.48
1:C:69:ASP:OD2	1:C:101:LYS:HD2	2.13	0.48
1:A:9:VAL:O	1:A:16:PHE:HA	2.13	0.48
1:B:96:LEU:HD22	1:B:148:LYS:HB3	1.95	0.48
1:B:206:LEU:HB2	1:B:226:PHE:HZ	1.79	0.48
1:D:144:LYS:HD3	1:D:151:ASN:HA	1.96	0.48
1:B:206:LEU:HD23	1:B:207:GLU:N	2.29	0.48
1:A:161:HIS:HB3	1:A:164:HIS:HB2	1.96	0.48
1:A:1:SER:C	1:A:3:ASN:N	2.72	0.48
1:A:25:ASN:ND2	1:A:27:GLN:HE21	2.08	0.48
1:A:176:ARG:O	1:A:326:LYS:HD2	2.12	0.48
1:C:69:ASP:HB3	1:C:86:PHE:O	2.14	0.48
1:C:126:LEU:HD13	1:C:156:PHE:CZ	2.48	0.48
1:A:6:ILE:N	1:A:6:ILE:HD12	2.29	0.47
1:A:75:MET:SD	1:A:75:MET:C	2.97	0.47
1:A:294:PHE:CD2	1:A:298:THR:HB	2.49	0.47
1:B:21:GLU:HB2	1:B:92:THR:HB	1.95	0.47
1:D:131:LEU:HB2	1:D:191:LEU:HB3	1.95	0.47
1:A:86:PHE:CD1	1:A:101:LYS:HE2	2.49	0.47
1:C:85:PHE:CE1	1:C:104:GLU:HB2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:97:SER:C	1:D:98:LEU:HD22	2.39	0.47
1:C:42:VAL:HG13	1:C:43:PRO:HD2	1.97	0.47
1:C:77:TYR:HD2	1:C:111:PHE:HE1	1.63	0.47
1:A:108:THR:O	1:A:112:GLU:HB2	2.14	0.47
1:A:221:THR:CB	1:A:300:ILE:HB	2.44	0.47
1:B:12:GLN:HB2	1:B:14:ILE:HG22	1.96	0.47
1:B:159:PRO:HD3	1:B:166:GLY:HA3	1.96	0.47
1:C:131:LEU:HB2	1:C:191:LEU:HD21	1.97	0.47
1:B:221:THR:OG1	1:B:300:ILE:HB	2.14	0.47
1:C:85:PHE:CE1	1:C:106:ILE:HD11	2.49	0.47
1:A:187:LEU:HD22	1:A:194:GLN:O	2.14	0.47
1:C:73:VAL:HG22	1:D:72:LYS:HB3	1.96	0.47
1:D:131:LEU:HD13	1:D:191:LEU:HB3	1.95	0.47
1:B:82:VAL:HG23	1:B:108:THR:HG22	1.97	0.47
1:D:176:ARG:NH1	1:D:329:LEU:HD13	2.29	0.47
1:D:254:LEU:HD13	1:D:274:LEU:HD11	1.95	0.47
1:A:191:LEU:HD23	1:A:191:LEU:O	2.14	0.47
1:A:193:TRP:O	1:A:212:ILE:HG23	2.15	0.47
1:D:155:THR:HA	1:D:312:VAL:HA	1.96	0.47
1:C:77:TYR:CD2	1:C:111:PHE:HE1	2.33	0.47
1:A:77:TYR:CZ	1:A:78:VAL:CG1	2.98	0.46
1:A:188:ASN:HB3	1:A:196:THR:OG1	2.14	0.46
1:B:98:LEU:HD23	1:B:98:LEU:N	2.29	0.46
1:B:187:LEU:HD11	1:B:193:TRP:CE3	2.50	0.46
1:B:232:GLN:HG3	1:B:236:VAL:HG12	1.97	0.46
1:D:153:LEU:HB2	1:D:313:PHE:O	2.15	0.46
1:A:36:GLY:O	1:A:131:LEU:HD23	2.14	0.46
1:A:40:LEU:HD12	1:A:41:TRP:H	1.79	0.46
1:B:174:GLU:HB3	1:B:177:PHE:CD2	2.50	0.46
1:A:130:ASP:CB	1:A:191:LEU:HG	2.45	0.46
1:A:220:ILE:HD13	1:A:302:GLY:H	1.81	0.46
1:B:210:ASN:O	1:B:298:THR:HA	2.16	0.46
1:A:29:PHE:HB2	1:A:31:PHE:HE2	1.78	0.46
1:D:5:ASN:ND2	1:D:169:THR:HG23	2.31	0.46
1:A:35:THR:HG1	1:A:214:ASP:HA	1.78	0.46
1:C:91:VAL:CG2	1:C:100:TYR:HB3	2.46	0.46
1:A:220:ILE:CD1	1:A:304:PRO:HB2	2.46	0.46
1:A:245:TYR:O	1:A:286:MET:HG3	2.16	0.46
1:D:231:LEU:HB3	1:D:253:LYS:CG	2.46	0.46
1:B:82:VAL:CG2	1:B:108:THR:HG22	2.46	0.46
1:D:156:PHE:HB2	1:D:311:THR:OG1	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:ILE:HG12	1:A:304:PRO:HG2	1.98	0.46
1:B:157:TYR:CZ	1:B:329:LEU:HD12	2.51	0.46
1:C:282:PRO:HB2	1:C:284:LEU:HD23	1.98	0.46
1:A:193:TRP:CE3	1:A:313:PHE:HD1	2.33	0.45
1:A:241:PHE:C	1:A:242:LEU:HD12	2.41	0.45
1:A:128:TRP:CD1	1:A:192:TYR:HA	2.51	0.45
1:A:210:ASN:O	1:A:212:ILE:HG13	2.16	0.45
1:A:269:GLU:H	1:A:272:TYR:HD2	1.63	0.45
1:A:248:LEU:HD22	1:A:248:LEU:H	1.82	0.45
1:C:143:LEU:HB3	1:C:149:ILE:HG23	1.99	0.45
1:A:326:LYS:HE2	1:A:329:LEU:OXT	2.16	0.45
1:A:8:LEU:CD2	1:A:33:LEU:HD11	2.47	0.45
1:A:292:LEU:HD21	1:B:78:VAL:HG12	1.97	0.45
1:C:144:LYS:HD2	1:C:152:ALA:CB	2.46	0.45
1:A:187:LEU:CD2	1:A:320:VAL:HG23	2.43	0.45
1:B:157:TYR:CZ	1:B:164:HIS:CE1	3.05	0.45
1:A:75:MET:HE1	1:A:77:TYR:CB	2.46	0.45
1:D:98:LEU:HD12	1:D:146:GLN:OE1	2.17	0.45
1:A:27:GLN:HB2	1:A:29:PHE:HE1	1.82	0.44
1:D:31:PHE:HA	1:D:122:GLY:O	2.17	0.44
1:A:173:ILE:HG22	1:A:178:TYR:HE1	1.82	0.44
1:B:222:VAL:HG22	1:B:223:PRO:HD2	1.99	0.44
1:D:143:LEU:HB3	1:D:149:ILE:HD13	2.00	0.44
1:A:59:ASP:OD2	1:A:62:LYS:HG3	2.17	0.44
1:B:6:ILE:HD11	1:B:19:ASP:H	1.82	0.44
1:D:5:ASN:N	1:D:5:ASN:ND2	2.65	0.44
1:A:17:TYR:CE2	1:A:120:PHE:HB3	2.52	0.44
1:B:44:SER:O	1:B:57:LEU:HD22	2.17	0.44
1:D:139:ILE:O	1:D:143:LEU:HG	2.18	0.44
1:A:75:MET:O	1:A:82:VAL:HB	2.17	0.44
1:B:126:LEU:HD13	1:B:156:PHE:CZ	2.52	0.44
1:B:213:VAL:HG22	1:B:301:LEU:HD12	2.00	0.44
1:C:100:TYR:HE1	1:C:136:VAL:HG13	1.81	0.44
1:D:246:VAL:HG11	1:D:284:LEU:HD11	1.99	0.44
1:A:33:LEU:CD2	1:A:124:LEU:HD23	2.47	0.44
1:B:89:ASP:O	1:B:99:PRO:HA	2.17	0.44
1:A:46:LYS:HB2	1:A:106:ILE:CD1	2.40	0.44
1:A:55:LYS:HG3	1:A:115:TYR:CZ	2.53	0.44
1:C:191:LEU:C	1:C:191:LEU:HD23	2.43	0.44
1:D:176:ARG:O	1:D:326:LYS:HD2	2.18	0.44
1:A:87:SER:HB2	1:A:102:PHE:CE1	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:10:ASP:HB3	1:C:161:HIS:HA	1.99	0.44
1:C:133:ILE:HG23	1:D:72:LYS:HZ3	1.83	0.44
1:D:157:TYR:CD1	1:D:177:PHE:HB3	2.53	0.44
1:B:222:VAL:CG1	1:B:227:LEU:HG	2.47	0.43
1:A:154:PHE:HE1	1:A:315:TYR:CE2	2.36	0.43
1:A:159:PRO:HB3	1:A:166:GLY:H	1.79	0.43
1:A:187:LEU:HD23	1:A:195:ILE:HG22	2.00	0.43
1:B:16:PHE:HB3	1:B:307:ARG:HH12	1.83	0.43
1:C:112:GLU:N	1:C:113:PRO:HD2	2.32	0.43
1:C:128:TRP:CH2	1:C:315:TYR:HA	2.53	0.43
1:D:153:LEU:CB	1:D:314:ASP:HA	2.48	0.43
1:A:303:ASP:N	1:A:304:PRO:CD	2.80	0.43
1:B:269:GLU:HB2	1:B:272:TYR:CE2	2.54	0.43
1:C:59:ASP:OD2	1:C:62:LYS:HG3	2.17	0.43
1:D:81:THR:HB	1:D:109:ASN:O	2.18	0.43
1:D:155:THR:HB	1:D:310:PHE:CZ	2.53	0.43
1:A:77:TYR:CZ	1:A:78:VAL:HG12	2.53	0.43
1:C:144:LYS:HB2	1:C:149:ILE:HD11	2.00	0.43
1:D:182:LEU:HD11	1:D:312:VAL:HG12	2.01	0.43
1:A:77:TYR:CG	1:A:78:VAL:N	2.86	0.43
1:A:248:LEU:HD11	1:C:240:PRO:HB2	2.01	0.43
1:B:4:ASP:HB3	1:B:170:ILE:HG12	2.00	0.43
1:B:227:LEU:HD23	1:B:227:LEU:HA	1.86	0.43
1:D:47:CYS:HA	1:D:106:ILE:O	2.19	0.43
1:A:37:SER:O	1:A:125:GLY:HA2	2.18	0.43
1:C:98:LEU:HD23	1:C:98:LEU:N	2.34	0.43
1:A:225:ASP:O	1:A:229:LYS:HG2	2.19	0.43
1:B:74:GLU:HA	1:B:82:VAL:O	2.17	0.43
1:C:91:VAL:HG21	1:C:100:TYR:HB3	2.01	0.43
1:C:206:LEU:HD22	1:C:209:ALA:HB2	2.01	0.43
1:D:142:GLU:O	1:D:146:GLN:HG3	2.18	0.43
1:D:178:TYR:HB3	1:D:310:PHE:HB3	2.00	0.43
1:B:287:LEU:N	1:B:287:LEU:HD12	2.34	0.43
1:C:87:SER:N	1:C:102:PHE:O	2.51	0.43
1:D:98:LEU:HD12	1:D:146:GLN:CD	2.43	0.43
1:C:131:LEU:HD22	1:C:192:TYR:CE1	2.54	0.43
1:D:157:TYR:HB2	1:D:310:PHE:CE1	2.54	0.43
1:B:222:VAL:CG2	1:B:223:PRO:HD2	2.49	0.42
1:C:16:PHE:HE2	1:C:33:LEU:HD12	1.83	0.42
1:C:234:LEU:HB3	1:C:236:VAL:HG12	2.01	0.42
1:D:201:VAL:HG11	1:D:230:MET:SD	2.59	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:GLN:HE22	1:A:59:ASP:HB3	1.84	0.42
1:A:128:TRP:CE3	1:A:315:TYR:CD1	3.05	0.42
1:D:8:LEU:CD2	1:D:33:LEU:HD11	2.50	0.42
1:D:153:LEU:N	1:D:153:LEU:HD23	2.34	0.42
1:A:179:GLU:CD	1:A:326:LYS:HA	2.43	0.42
1:B:98:LEU:HD21	1:B:139:ILE:CG2	2.49	0.42
1:B:277:ILE:HG12	1:B:286:MET:HG3	2.01	0.42
1:A:223:PRO:HD2	1:A:299:PHE:CE2	2.54	0.42
1:B:98:LEU:HD11	1:B:139:ILE:HG23	2.02	0.42
1:B:157:TYR:O	1:B:166:GLY:HA2	2.19	0.42
1:D:179:GLU:H	1:D:179:GLU:HG2	1.73	0.42
1:D:217:THR:HG22	1:D:219:ALA:H	1.84	0.42
1:A:25:ASN:CG	1:A:27:GLN:HG3	2.44	0.42
1:A:157:TYR:CE2	1:A:164:HIS:CE1	3.07	0.42
1:B:235:ASP:O	1:B:236:VAL:C	2.62	0.42
1:B:269:GLU:HB2	1:B:272:TYR:HE2	1.84	0.42
1:D:230:MET:C	1:D:232:GLN:N	2.77	0.42
1:D:235:ASP:HB2	1:D:253:LYS:NZ	2.34	0.42
1:A:296:VAL:HB	1:A:297:PRO:HD2	2.01	0.42
1:C:73:VAL:HG11	1:C:133:ILE:HG21	2.02	0.42
1:C:129:LYS:HB2	1:C:138:PRO:HD3	2.02	0.42
1:C:158:LEU:HD12	1:C:307:ARG:HG3	2.02	0.42
1:D:3:ASN:CG	1:D:172:GLY:HA3	2.45	0.42
1:D:184:TYR:CA	1:D:321:GLY:HA2	2.44	0.42
1:C:51:GLY:H	1:C:112:GLU:CD	2.27	0.42
1:C:296:VAL:HB	1:C:297:PRO:HD2	2.02	0.42
1:A:33:LEU:HD22	1:A:124:LEU:HD23	2.02	0.42
1:B:214:ASP:HB3	1:B:217:THR:HG22	2.01	0.42
1:C:131:LEU:HB2	1:C:191:LEU:CD2	2.50	0.42
1:B:303:ASP:O	1:B:304:PRO:C	2.63	0.42
1:A:68:LYS:HE3	1:A:85:PHE:CE2	2.55	0.42
1:A:130:ASP:CG	1:A:191:LEU:HA	2.45	0.42
1:A:168:LEU:O	1:A:170:ILE:HG13	2.20	0.42
1:B:222:VAL:HG23	1:B:299:PHE:CD2	2.55	0.42
1:C:79:SER:HB3	1:C:111:PHE:CE1	2.54	0.42
1:A:308:LYS:HD3	1:A:309:TYR:CE2	2.55	0.41
1:C:139:ILE:O	1:C:140:VAL:C	2.63	0.41
1:D:17:TYR:CD2	1:D:32:ILE:HG12	2.54	0.41
1:D:33:LEU:HD23	1:D:124:LEU:HD12	2.01	0.41
1:D:151:ASN:CG	1:D:153:LEU:HD22	2.44	0.41
1:A:163:LYS:CD	1:A:329:LEU:HD11	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:ILE:HG12	1:A:304:PRO:HB2	2.02	0.41
1:B:157:TYR:O	1:B:158:LEU:C	2.63	0.41
1:C:25:ASN:HD22	1:C:62:LYS:HB2	1.86	0.41
1:D:33:LEU:CD2	1:D:124:LEU:HD12	2.49	0.41
1:B:36:GLY:O	1:B:131:LEU:HD21	2.20	0.41
1:B:164:HIS:HB2	1:B:327:LYS:HG3	2.02	0.41
1:B:206:LEU:HD23	1:B:207:GLU:H	1.86	0.41
1:B:229:LYS:NZ	1:B:229:LYS:HB2	2.34	0.41
1:C:51:GLY:N	1:C:112:GLU:OE2	2.53	0.41
1:C:274:LEU:HD22	1:C:285:CYS:HB3	2.02	0.41
1:D:38:ALA:HB3	1:D:132:SER:HA	2.01	0.41
1:D:44:SER:O	1:D:57:LEU:HD22	2.20	0.41
1:D:55:LYS:HD3	1:D:115:TYR:CZ	2.55	0.41
1:D:173:ILE:HD13	1:D:173:ILE:HA	1.88	0.41
1:D:182:LEU:HD11	1:D:312:VAL:CG1	2.50	0.41
1:A:42:VAL:HG21	1:A:102:PHE:CD1	2.55	0.41
1:B:156:PHE:CD2	1:B:168:LEU:HD13	2.55	0.41
1:C:87:SER:C	1:C:101:LYS:HA	2.46	0.41
1:A:267:THR:O	1:A:308:LYS:NZ	2.46	0.41
1:B:131:LEU:HB2	1:B:191:LEU:CD2	2.50	0.41
1:B:201:VAL:HB	1:B:204:ILE:HB	2.02	0.41
1:C:40:LEU:CB	1:C:100:TYR:HE2	2.33	0.41
1:C:133:ILE:HG12	1:D:74:GLU:HG2	2.01	0.41
1:D:86:PHE:CG	1:D:101:LYS:HE2	2.55	0.41
1:D:129:LYS:HA	1:D:136:VAL:O	2.20	0.41
1:D:174:GLU:O	1:D:175:GLU:HB2	2.19	0.41
1:D:187:LEU:HD13	1:D:190:ASP:HA	2.01	0.41
1:A:41:TRP:CH2	1:A:103:ILE:HD13	2.56	0.41
1:A:128:TRP:CZ3	1:A:315:TYR:HD1	2.38	0.41
1:A:191:LEU:C	1:A:191:LEU:CD2	2.94	0.41
1:A:221:THR:HB	1:A:300:ILE:HB	2.01	0.41
1:B:220:ILE:HD13	1:B:302:GLY:H	1.84	0.41
1:C:128:TRP:CE3	1:C:315:TYR:HD2	2.39	0.41
1:C:179:GLU:H	1:C:179:GLU:HG2	1.71	0.41
1:D:32:ILE:HD11	1:D:120:PHE:CG	2.55	0.41
1:D:77:TYR:CE2	1:D:78:VAL:HG12	2.55	0.41
1:D:92:THR:HA	1:D:97:SER:HA	2.02	0.41
1:D:155:THR:O	1:D:168:LEU:HD12	2.21	0.41
1:A:46:LYS:NZ	1:A:85:PHE:HZ	2.19	0.41
1:A:189:HIS:ND1	1:A:191:LEU:HB2	2.35	0.41
1:B:200:HIS:CE1	1:B:207:GLU:HG2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:294:PHE:CZ	1:B:300:ILE:HD13	2.55	0.41
1:D:184:TYR:HA	1:D:320:VAL:O	2.21	0.41
1:A:89:ASP:HB3	1:A:102:PHE:HE2	1.86	0.41
1:B:10:ASP:HB2	1:B:159:PRO:CG	2.40	0.41
1:B:223:PRO:HD2	1:B:299:PHE:CE2	2.56	0.41
1:B:223:PRO:HB2	1:B:226:PHE:CD1	2.56	0.41
1:B:232:GLN:HA	1:B:236:VAL:HG12	2.01	0.41
1:C:100:TYR:HE1	1:C:136:VAL:CG1	2.34	0.41
1:D:3:ASN:HB3	1:D:169:THR:CG2	2.50	0.41
1:D:108:THR:O	1:D:112:GLU:N	2.53	0.41
1:D:157:TYR:CE2	1:D:164:HIS:CD2	3.09	0.41
1:D:247:THR:HG21	1:D:254:LEU:HD21	2.03	0.41
1:B:17:TYR:CD2	1:B:32:ILE:HD11	2.55	0.41
1:C:153:LEU:HA	1:C:313:PHE:O	2.20	0.41
1:D:220:ILE:HD13	1:D:302:GLY:H	1.86	0.41
1:A:29:PHE:HE2	1:A:57:LEU:N	2.19	0.40
1:B:37:SER:HB2	1:B:41:TRP:HZ3	1.86	0.40
1:C:158:LEU:HD11	1:C:306:MET:HE3	2.02	0.40
1:D:129:LYS:HB2	1:D:137:ASP:HA	2.02	0.40
1:A:40:LEU:HD12	1:A:41:TRP:N	2.37	0.40
1:B:75:MET:HB3	1:B:82:VAL:HG12	2.03	0.40
1:A:327:LYS:O	1:A:328:ASN:C	2.64	0.40
1:D:173:ILE:C	1:D:175:GLU:H	2.29	0.40
1:A:86:PHE:HB3	1:A:101:LYS:HE2	2.04	0.40
1:A:292:LEU:CD2	1:A:294:PHE:CE1	3.04	0.40
1:B:45:VAL:HA	1:B:57:LEU:HD13	2.03	0.40
1:A:90:LEU:HD11	1:A:97:SER:HB3	2.04	0.40
1:C:224:THR:HG23	1:C:293:ASP:OD1	2.21	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	327/329 (99%)	307 (94%)	20 (6%)	0	100	100
1	B	327/329 (99%)	306 (94%)	21 (6%)	0	100	100
1	C	327/329 (99%)	309 (94%)	18 (6%)	0	100	100
1	D	327/329 (99%)	301 (92%)	26 (8%)	0	100	100
All	All	1308/1316 (99%)	1223 (94%)	85 (6%)	0	100	100

There are no Ramachandran outliers to report.

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	294/294 (100%)	292 (99%)	2 (1%)	76	79
1	B	294/294 (100%)	288 (98%)	6 (2%)	48	67
1	C	294/294 (100%)	290 (99%)	4 (1%)	59	72
1	D	294/294 (100%)	290 (99%)	4 (1%)	59	72
All	All	1176/1176 (100%)	1160 (99%)	16 (1%)	59	72

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	114	THR
1	A	153	LEU
1	B	6	ILE
1	B	9	VAL
1	B	123	ILE
1	B	212	ILE
1	B	225	ASP
1	B	248	LEU
1	C	82	VAL
1	C	136	VAL
1	C	137	ASP
1	C	188	ASN
1	D	5	ASN

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Mol	Chain	Res	Type
1	D	82	VAL
1	D	280	VAL
1	D	312	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	27	GLN
1	A	147	ASN
1	A	210	ASN
1	A	232	GLN
1	B	3	ASN
1	B	26	GLN
1	B	76	ASN
1	B	109	ASN
1	B	189	HIS
1	B	200	HIS
1	B	251	ASN
1	B	318	HIS
1	C	76	ASN
1	C	151	ASN
1	C	161	HIS
1	C	164	HIS
1	C	251	ASN
1	D	5	ASN
1	D	27	GLN
1	D	39	ASN
1	D	147	ASN
1	D	188	ASN
1	D	194	GLN
1	D	318	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

1 ligand 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	I0J	A	401	-	32,32,32	0.34	0	41,43,43	0.37	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	I0J	A	401	-	-	7/12/19/19	0/4/4/4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

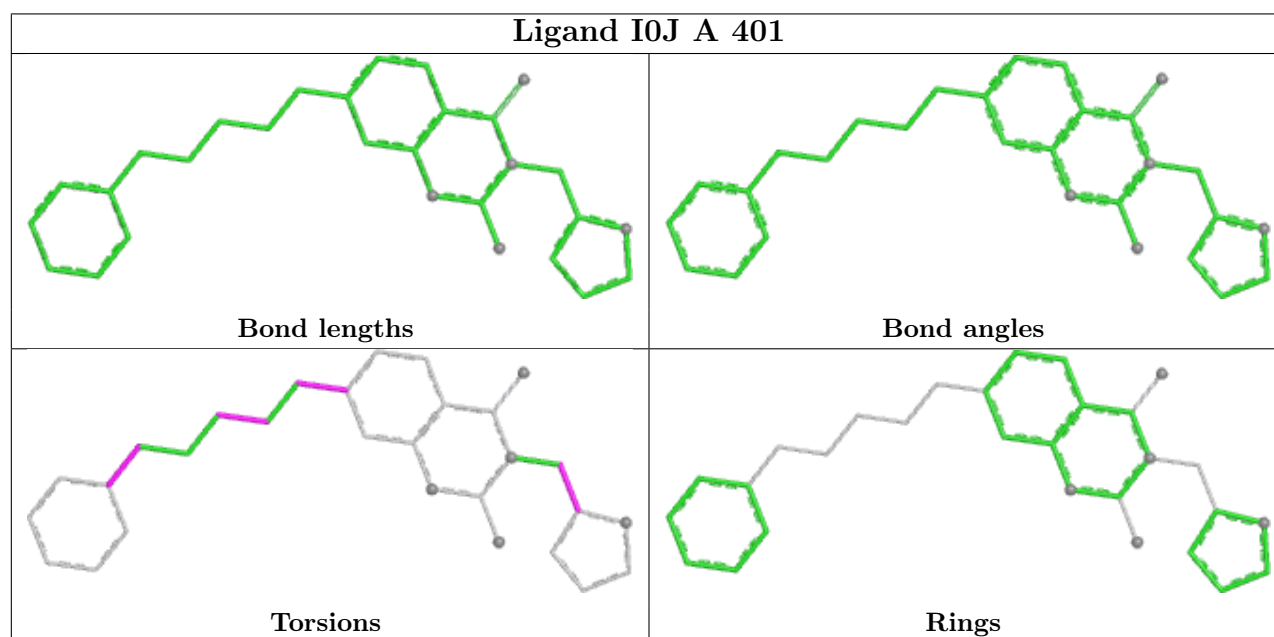
Mol	Chain	Res	Type	Atoms
2	A	401	I0J	C4-C5-C6-N7
2	A	401	I0J	O2-C5-C6-N7
2	A	401	I0J	C14-C18-C19-C20
2	A	401	I0J	C17-C18-C19-C20
2	A	401	I0J	C22-C23-C24-C25
2	A	401	I0J	C19-C20-C21-C22
2	A	401	I0J	C22-C23-C24-C29

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	I0J	1	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 [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	329/329 (100%)	0.83	36 (10%)	10 10	37, 69, 93, 111	0
1	B	329/329 (100%)	0.88	42 (12%)	7 7	39, 68, 106, 125	0
1	C	329/329 (100%)	0.94	46 (13%)	6 6	51, 89, 117, 142	0
1	D	329/329 (100%)	1.10	53 (16%)	4 5	61, 91, 128, 163	0
All	All	1316/1316 (100%)	0.93	177 (13%)	7 7	37, 79, 115, 163	0

All (177) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	171	GLY	7.5
1	D	257	PHE	5.4
1	D	259	PHE	5.4
1	D	192	TYR	5.3
1	D	100	TYR	5.2
1	D	230	MET	5.1
1	D	75	MET	5.0
1	C	102	PHE	4.7
1	B	237	ILE	4.6
1	C	148	LYS	4.5
1	C	100	TYR	4.5
1	D	101	LYS	4.4
1	C	257	PHE	4.4
1	B	177	PHE	4.4
1	C	177	PHE	4.4
1	B	109	ASN	4.2
1	C	170	ILE	4.2
1	D	289	ILE	4.1
1	B	108	THR	4.1
1	A	33	LEU	4.0
1	B	234	LEU	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	193	TRP	3.9
1	A	257	PHE	3.8
1	B	42	VAL	3.7
1	C	292	LEU	3.7
1	A	76	ASN	3.7
1	D	305	PHE	3.7
1	D	168	LEU	3.6
1	A	298	THR	3.6
1	B	33	LEU	3.6
1	A	124	LEU	3.5
1	C	301	LEU	3.5
1	C	105	VAL	3.5
1	B	299	PHE	3.5
1	D	33	LEU	3.5
1	D	156	PHE	3.5
1	D	184	TYR	3.5
1	D	189	HIS	3.4
1	A	102	PHE	3.4
1	C	33	LEU	3.4
1	B	284	LEU	3.4
1	D	76	ASN	3.4
1	D	143	LEU	3.4
1	A	6	ILE	3.3
1	D	16	PHE	3.3
1	D	193	TRP	3.3
1	B	86	PHE	3.2
1	C	241	PHE	3.2
1	B	93	VAL	3.2
1	A	123	ILE	3.2
1	B	239	VAL	3.2
1	D	110	GLY	3.1
1	B	52	CYS	3.1
1	C	76	ASN	3.1
1	C	124	LEU	3.1
1	C	320	VAL	3.1
1	A	230	MET	3.1
1	A	141	VAL	3.0
1	C	86	PHE	3.0
1	D	190	ASP	3.0
1	A	156	PHE	3.0
1	D	149	ILE	3.0
1	B	320	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	31	PHE	3.0
1	C	277	ILE	3.0
1	D	297	PRO	2.9
1	A	22	VAL	2.9
1	A	41	TRP	2.9
1	B	241	PHE	2.9
1	B	257	PHE	2.9
1	D	96	LEU	2.9
1	C	329	LEU	2.9
1	C	226	PHE	2.9
1	A	168	LEU	2.9
1	A	195	ILE	2.9
1	A	16	PHE	2.8
1	D	104	GLU	2.8
1	B	158	LEU	2.8
1	C	305	PHE	2.8
1	B	75	MET	2.8
1	B	124	LEU	2.8
1	C	230	MET	2.8
1	C	123	ILE	2.8
1	B	58	TYR	2.7
1	C	176	ARG	2.7
1	B	79	SER	2.7
1	B	78	VAL	2.7
1	B	6	ILE	2.7
1	D	41	TRP	2.7
1	A	74	GLU	2.7
1	A	178	TYR	2.7
1	B	100	TYR	2.7
1	C	19	ASP	2.7
1	A	140	VAL	2.7
1	C	167	PHE	2.7
1	D	232	GLN	2.6
1	D	218	SER	2.6
1	D	284	LEU	2.6
1	B	157	TYR	2.6
1	D	241	PHE	2.6
1	A	248	LEU	2.6
1	B	126	LEU	2.6
1	C	192	TYR	2.6
1	C	126	LEU	2.6
1	B	101	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	16	PHE	2.6
1	D	169	THR	2.6
1	D	247	THR	2.6
1	B	16	PHE	2.5
1	C	156	PHE	2.5
1	D	273	TYR	2.5
1	A	235	ASP	2.5
1	B	40	LEU	2.5
1	D	10	ASP	2.5
1	D	31	PHE	2.5
1	D	37	SER	2.5
1	D	177	PHE	2.5
1	C	18	GLY	2.5
1	D	172	GLY	2.4
1	B	120	PHE	2.4
1	D	86	PHE	2.4
1	D	226	PHE	2.4
1	D	299	PHE	2.4
1	A	192	TYR	2.4
1	D	287	LEU	2.4
1	C	280	VAL	2.4
1	B	178	TYR	2.4
1	D	77	TYR	2.4
1	D	283	GLY	2.4
1	D	195	ILE	2.4
1	A	241	PHE	2.4
1	C	42	VAL	2.4
1	D	91	VAL	2.3
1	D	238	LYS	2.3
1	B	69	ASP	2.3
1	D	148	LYS	2.3
1	C	234	LEU	2.3
1	C	259	PHE	2.3
1	C	283	GLY	2.3
1	B	230	MET	2.3
1	B	192	TYR	2.3
1	B	259	PHE	2.3
1	A	206	LEU	2.2
1	A	296	VAL	2.2
1	A	8	LEU	2.2
1	A	182	LEU	2.2
1	C	200	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	246	VAL	2.2
1	C	74	GLU	2.2
1	A	324	LEU	2.2
1	C	285	CYS	2.2
1	A	177	PHE	2.2
1	C	294	PHE	2.2
1	C	41	TRP	2.2
1	A	81	THR	2.1
1	C	46	LYS	2.1
1	B	240	PRO	2.1
1	A	143	LEU	2.1
1	A	320	VAL	2.1
1	C	279	ASP	2.1
1	C	131	LEU	2.1
1	C	194	GLN	2.1
1	B	51	GLY	2.1
1	D	39	ASN	2.1
1	D	320	VAL	2.1
1	C	191	LEU	2.1
1	C	184	TYR	2.1
1	D	313	PHE	2.1
1	A	42	VAL	2.1
1	D	327	LYS	2.1
1	A	170	ILE	2.1
1	A	301	LEU	2.1
1	D	102	PHE	2.1
1	B	231	LEU	2.0
1	B	123	ILE	2.0
1	B	266	TYR	2.0
1	B	238	LYS	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](#)

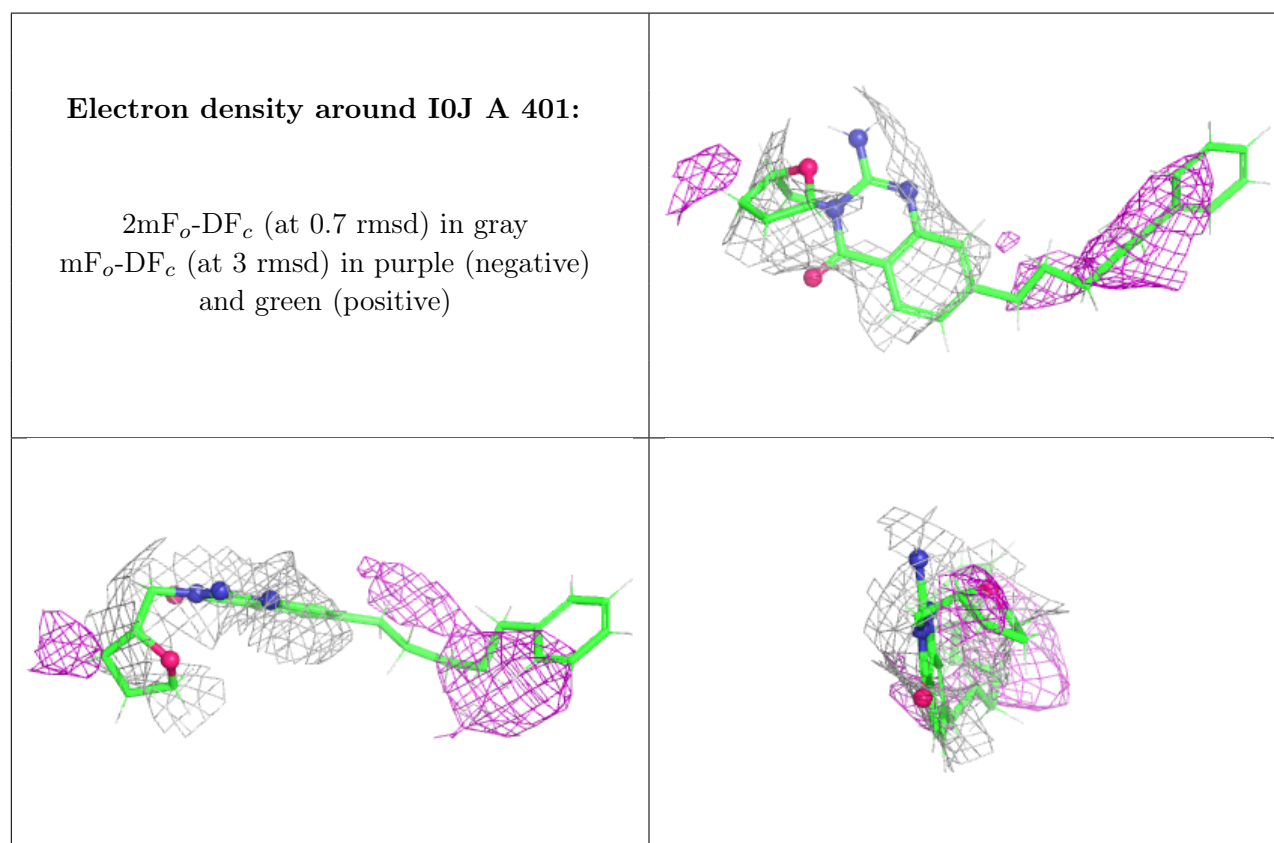
There are no oligosaccharides in this entry.

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($\text{\AA}^2$)	Q<0.9
2	I0J	A	401	29/29	0.84	0.15	69,91,113,117	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.