



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:26 PM UTC

PDB ID : 4UJ4 / pdb_00004uj4
Title : Crystal structure of human Rab11-Rabin8-FIP3
Authors : Vetter, M.; Lorentzen, E.
Deposited on : 2015-04-08
Resolution : 4.20 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

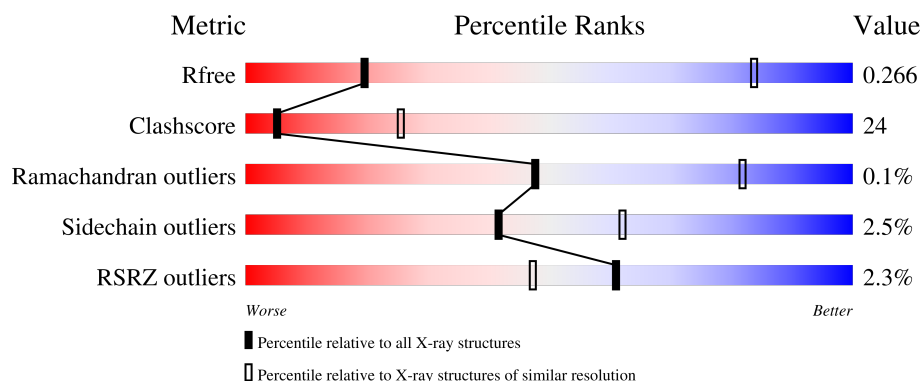
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.20 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	1004 (4.52-3.88)
Clashscore	190562	1019 (4.50-3.90)
Ramachandran outliers	187476	1020 (4.56-3.84)
Sidechain outliers	187428	1006 (4.56-3.84)
RSRZ outliers	180081	1001 (4.52-3.88)

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	185	4% 67% 26% • 5%
1	D	185	% 64% 30% • 5%
1	G	185	% 62% 29% • 5%
1	J	185	% 58% 35% • 6%
2	B	195	3% 48% 33% • 17%

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Mol	Chain	Length	Quality of chain
2	E	195	2% 54% 27% • 17%
2	H	195	% 42% 35% • 21%
2	K	195	2% 44% 35% • 18%
3	C	66	32% 23% 45%
3	F	66	2% 27% 29% 44%
3	I	66	9% 35% 26% • 38%
3	L	66	2% 42% 15% 42%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11717 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 Ras-related protein Rab-11A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	176	Total	C	N	O	S	0	0	0
			1365	862	232	270	1			
1	D	175	Total	C	N	O	S	0	2	0
			1368	864	233	270	1			
1	G	175	Total	C	N	O	S	0	2	0
			1370	865	233	271	1			
1	J	174	Total	C	N	O	S	0	0	0
			1336	847	225	263	1			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	GLY	-	expression tag	UNP P62491
A	3	ALA	-	expression tag	UNP P62491
A	70	LEU	GLN	engineered mutation	UNP P62491
D	2	GLY	-	expression tag	UNP P62491
D	3	ALA	-	expression tag	UNP P62491
D	70	LEU	GLN	engineered mutation	UNP P62491
G	2	GLY	-	expression tag	UNP P62491
G	3	ALA	-	expression tag	UNP P62491
G	70	LEU	GLN	engineered mutation	UNP P62491
J	2	GLY	-	expression tag	UNP P62491
J	3	ALA	-	expression tag	UNP P62491
J	70	LEU	GLN	engineered mutation	UNP P62491

- Molecule 2 is a protein called Rab-3A-interacting protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	161	Total	C	N	O	S	0	0	0
			1288	822	211	243	12			
2	E	162	Total	C	N	O	S	0	0	0
			1268	811	207	238	12			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	155	Total	C	N	O	S	0	2	0
			1263	810	208	233	12			
2	K	159	Total	C	N	O	S	0	1	0
			1259	804	204	239	12			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	266	GLY	-	expression tag	UNP Q96QF0
B	267	ALA	-	expression tag	UNP Q96QF0
B	268	ALA	-	expression tag	UNP Q96QF0
B	269	SER	-	expression tag	UNP Q96QF0
E	266	GLY	-	expression tag	UNP Q96QF0
E	267	ALA	-	expression tag	UNP Q96QF0
E	268	ALA	-	expression tag	UNP Q96QF0
E	269	SER	-	expression tag	UNP Q96QF0
H	266	GLY	-	expression tag	UNP Q96QF0
H	267	ALA	-	expression tag	UNP Q96QF0
H	268	ALA	-	expression tag	UNP Q96QF0
H	269	SER	-	expression tag	UNP Q96QF0
K	266	GLY	-	expression tag	UNP Q96QF0
K	267	ALA	-	expression tag	UNP Q96QF0
K	268	ALA	-	expression tag	UNP Q96QF0
K	269	SER	-	expression tag	UNP Q96QF0

- Molecule 3 is a protein called Rab11 family-interacting protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	36	Total	C	N	O	S	0	0	0
			265	169	46	48	2			
3	F	37	Total	C	N	O	S	0	0	0
			251	163	40	46	2			
3	I	41	Total	C	N	O	S	0	0	0
			287	184	47	54	2			
3	L	38	Total	C	N	O	S	0	0	0
			265	172	44	47	2			

There are 16 discrepancies between the modelled and reference sequences:

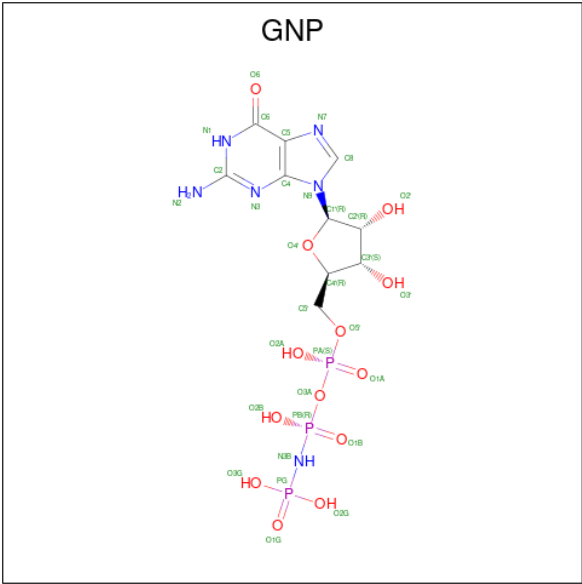
Chain	Residue	Modelled	Actual	Comment	Reference
C	691	GLY	-	expression tag	UNP O75154
C	692	ALA	-	expression tag	UNP O75154

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Chain	Residue	Modelled	Actual	Comment	Reference
C	693	ALA	-	expression tag	UNP O75154
C	694	SER	-	expression tag	UNP O75154
F	691	GLY	-	expression tag	UNP O75154
F	692	ALA	-	expression tag	UNP O75154
F	693	ALA	-	expression tag	UNP O75154
F	694	SER	-	expression tag	UNP O75154
I	691	GLY	-	expression tag	UNP O75154
I	692	ALA	-	expression tag	UNP O75154
I	693	ALA	-	expression tag	UNP O75154
I	694	SER	-	expression tag	UNP O75154
L	691	GLY	-	expression tag	UNP O75154
L	692	ALA	-	expression tag	UNP O75154
L	693	ALA	-	expression tag	UNP O75154
L	694	SER	-	expression tag	UNP O75154

- Molecule 4 is PHOSPHOAMINOPHOSPHONIC ACID-GUANYLATE ESTER (CCD ID: GNP) (formula: C₁₀H₁₇N₆O₁₃P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
4	D	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
4	G	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
4	J	1	Total	C	N	O	P	0	0
			32	10	6	13	3		

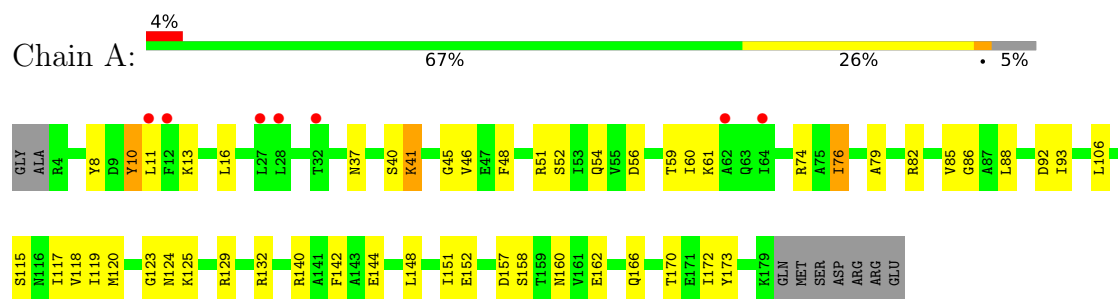
- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total 1	Mg 1	0	0
5	D	1	Total 1	Mg 1	0	0
5	G	1	Total 1	Mg 1	0	0
5	J	1	Total 1	Mg 1	0	0

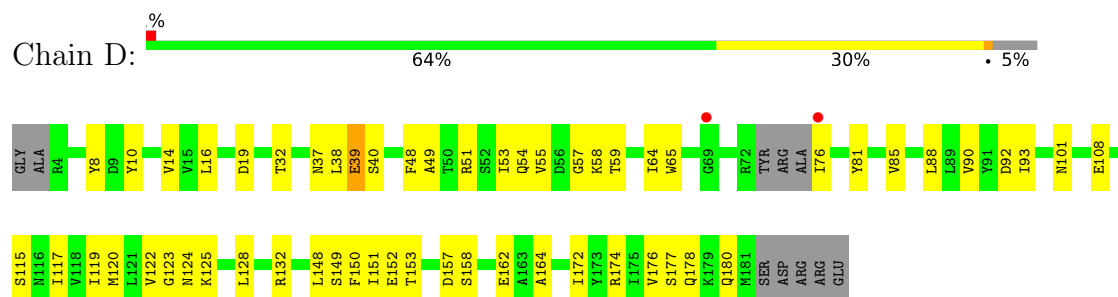
3 Residue-property plots [i](#)

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.

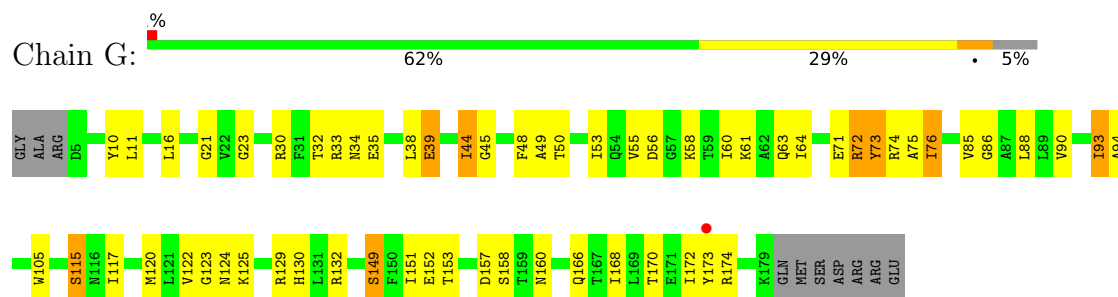
• Molecule 1: Ras-related protein Rab-11A



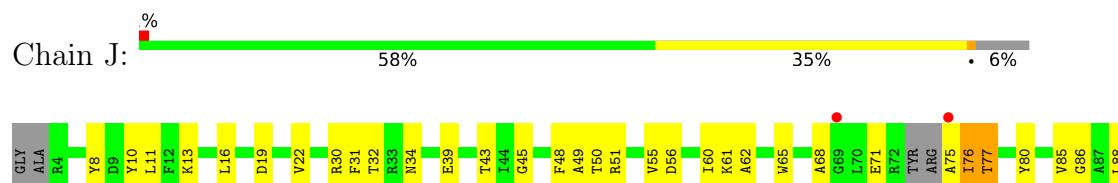
• Molecule 1: Ras-related protein Rab-11A

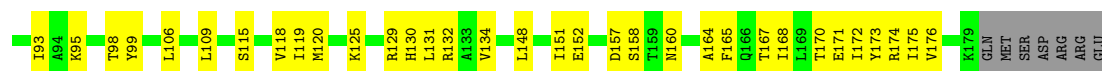


• Molecule 1: Ras-related protein Rab-11A

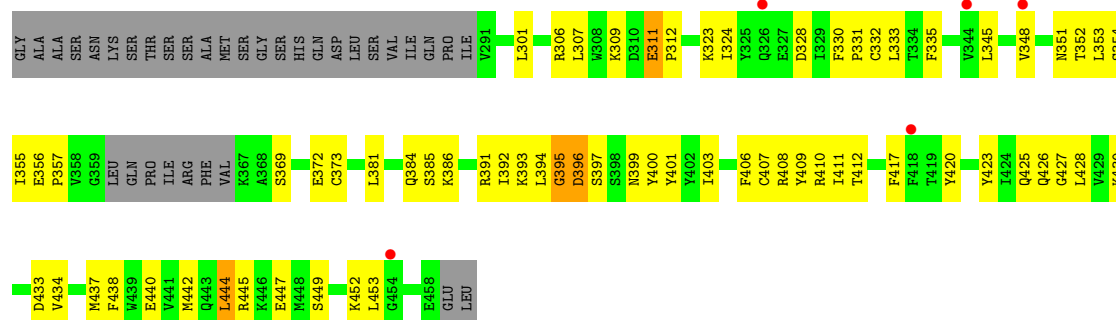


• Molecule 1: Ras-related protein Rab-11A

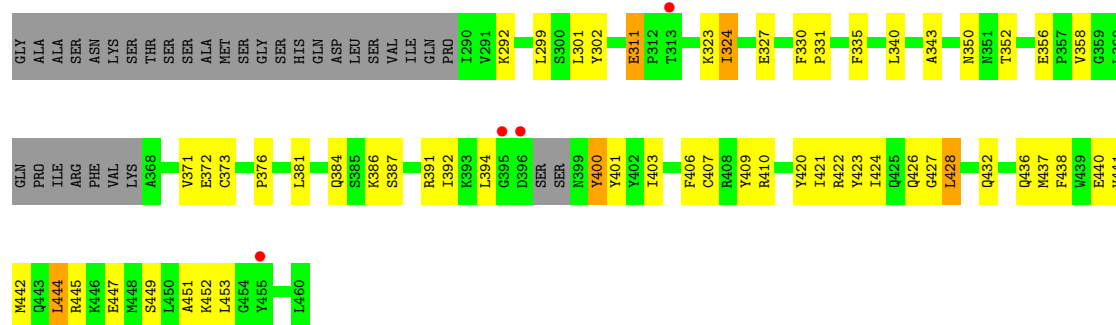




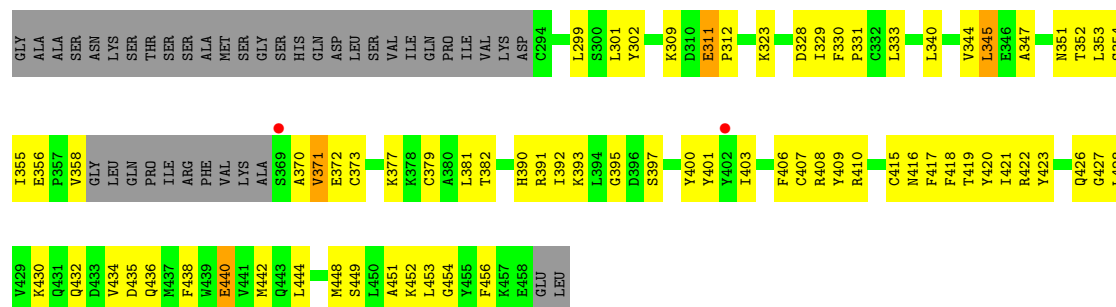
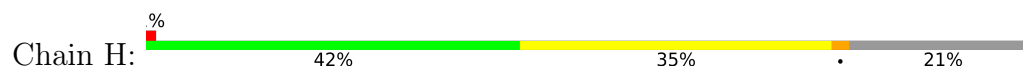
• Molecule 2: Rab-3A-interacting protein



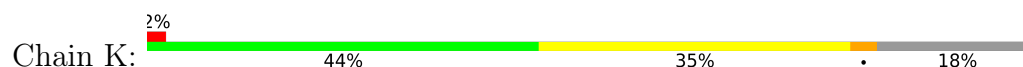
• Molecule 2: Rab-3A-interacting protein

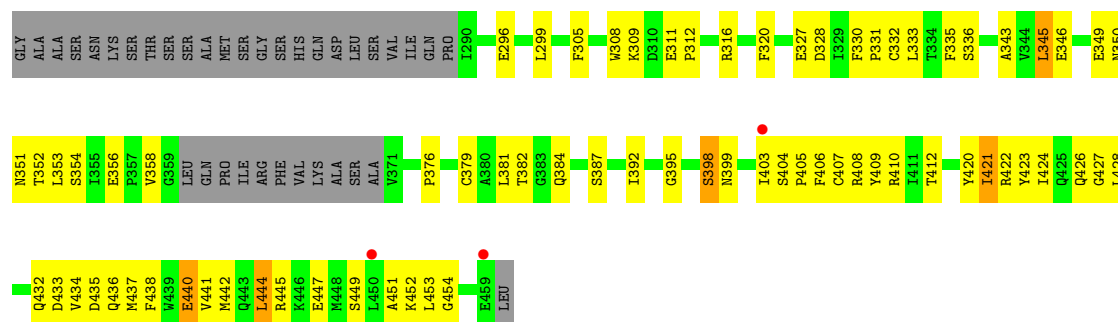


• Molecule 2: Rab-3A-interacting protein



• Molecule 2: Rab-3A-interacting protein





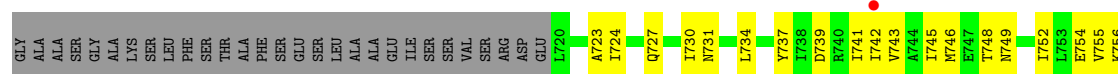
• Molecule 3: Rab11 family-interacting protein 3

Chain C: 32% 23% 45%



• Molecule 3: Rab11 family-interacting protein 3

Chain F: 2% 27% 29% 44%



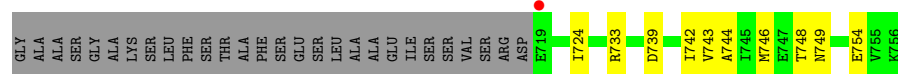
• Molecule 3: Rab11 family-interacting protein 3

Chain I: 9% 35% 26% 38%



• Molecule 3: Rab11 family-interacting protein 3

Chain L: 2% 42% 15% 42%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	66.12Å 161.70Å 192.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.47 – 4.20 49.47 – 4.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.47-4.20) 90.3 (49.47-4.20)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.68 (at 4.00Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.252 , 0.266 0.256 , 0.266	Depositor DCC
R_{free} test set	933 reflections (5.18%)	wwPDB-VP
Wilson B-factor (Å ²)	129.2	Xtriage
Anisotropy	0.299	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 172.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	11717	wwPDB-VP
Average B, all atoms (Å ²)	161.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.70% 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: MG, GNP

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	0.46	0/1386	0.83	1/1878 (0.1%)
1	D	0.46	0/1394	0.85	2/1889 (0.1%)
1	G	0.52	0/1397	0.92	2/1893 (0.1%)
1	J	0.46	0/1356	0.82	1/1839 (0.1%)
2	B	0.44	0/1315	0.88	3/1772 (0.2%)
2	E	0.39	0/1294	0.89	4/1748 (0.2%)
2	H	0.50	0/1296	0.94	4/1746 (0.2%)
2	K	0.39	0/1289	0.90	1/1742 (0.1%)
3	C	0.50	0/266	0.84	0/362
3	F	0.49	0/252	0.91	0/346
3	I	0.45	0/288	0.92	2/394 (0.5%)
3	L	0.46	0/266	0.88	0/364
All	All	0.46	0/11799	0.88	20/15973 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	G	0	3
2	B	0	1
2	K	0	1
All	All	0	5

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	371	VAL	N-CA-C	7.39	119.06	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	311	GLU	CA-C-N	7.15	127.12	119.76
2	B	311	GLU	C-N-CA	7.15	127.12	119.76
1	A	76	ILE	N-CA-C	-6.48	106.14	111.91
2	E	311	GLU	CA-C-N	6.16	126.11	119.76
2	E	311	GLU	C-N-CA	6.16	126.11	119.76
2	E	400	TYR	N-CA-C	6.07	118.53	109.25
2	B	395	GLY	N-CA-C	5.81	119.76	112.79
1	G	75	ALA	N-CA-C	5.75	123.04	110.80
3	I	717	ARG	N-CA-C	5.75	116.50	108.00
2	H	311	GLU	CA-C-N	5.62	125.54	119.76
2	H	311	GLU	C-N-CA	5.62	125.54	119.76
1	G	124	ASN	N-CA-C	5.59	118.80	110.14
3	I	717	ARG	CB-CA-C	-5.48	108.67	116.34
1	J	77	THR	N-CA-C	-5.39	105.44	112.23
2	H	371	VAL	CB-CA-C	-5.28	105.16	112.24
2	K	395	GLY	N-CA-C	5.19	119.06	110.97
1	D	8	TYR	N-CA-C	5.15	117.02	109.24
1	D	149	SER	N-CA-C	-5.14	104.09	110.41
2	E	371	VAL	N-CA-C	-5.11	104.60	111.44

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	396	ASP	Peptide
1	G	72	ARG	Peptide
1	G	73	TYR	Peptide
1	G	74	ARG	Peptide
2	K	398	SER	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1365	0	1322	68	1
1	D	1368	0	1322	56	2
1	G	1370	0	1328	58	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	J	1336	0	1290	80	3
2	B	1288	0	1249	76	0
2	E	1268	0	1197	65	1
2	H	1263	0	1244	77	3
2	K	1259	0	1185	72	1
3	C	265	0	255	33	0
3	F	251	0	227	46	0
3	I	287	0	261	36	0
3	L	265	0	249	22	1
4	A	32	0	13	1	0
4	D	32	0	13	1	0
4	G	32	0	13	6	0
4	J	32	0	13	2	0
5	A	1	0	0	0	0
5	D	1	0	0	0	0
5	G	1	0	0	0	0
5	J	1	0	0	0	0
All	All	11717	0	11181	561	6

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 24.

All (561) 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:J:76:ILE:CG2	3:L:743:VAL:HG22	1.60	1.29
1:J:76:ILE:CG2	3:L:743:VAL:CG2	2.25	1.15
1:J:76:ILE:HG21	3:L:743:VAL:CG2	1.78	1.12
1:A:76:ILE:HD11	3:C:742:ILE:HG22	1.31	1.12
1:J:48:PHE:CZ	3:L:754:GLU:HB2	1.87	1.10
1:A:76:ILE:HG13	3:C:743:VAL:CG2	1.82	1.09
1:A:76:ILE:HG13	3:C:743:VAL:HG22	1.30	1.08
3:F:727:GLN:NE2	3:I:728:GLU:HA	1.67	1.08
1:A:129:ARG:HD2	2:B:351:ASN:O	1.55	1.07
1:D:76:ILE:HD11	3:F:739:ASP:HA	1.33	1.07
2:B:391:ARG:HD3	2:B:400:TYR:HB3	1.11	1.06
1:J:76:ILE:HG22	3:L:743:VAL:HG22	1.39	1.03
1:D:76:ILE:HD11	3:F:739:ASP:CA	1.93	0.98
2:K:309:LYS:HE2	2:K:434:VAL:HG22	1.45	0.98
2:E:299:LEU:HD11	2:K:384:GLN:OE1	1.63	0.97
2:B:393:LYS:HE2	2:B:397:SER:HB3	1.45	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:ILE:HD11	3:C:742:ILE:CG2	1.95	0.95
1:D:65:TRP:HZ3	3:F:746:MET:HE2	1.32	0.94
2:B:391:ARG:HD3	2:B:400:TYR:CB	1.98	0.93
1:D:65:TRP:CZ3	3:F:746:MET:HE2	2.05	0.91
1:G:76:ILE:HG22	3:I:743:VAL:HG22	1.51	0.91
1:J:76:ILE:HG22	3:L:743:VAL:CG2	1.95	0.91
1:J:76:ILE:HG21	3:L:743:VAL:HG22	1.36	0.90
1:G:129[A]:ARG:HB3	2:H:354:SER:HB3	1.54	0.90
1:D:54:GLN:HE22	1:D:59:THR:HG22	1.34	0.89
2:E:391:ARG:HD2	2:E:400:TYR:HB3	1.56	0.88
2:B:384:GLN:NE2	2:H:299:LEU:HD21	1.89	0.87
2:E:292:LYS:NZ	3:I:718:ASP:OD2	2.08	0.86
1:G:129[B]:ARG:HB3	2:H:354:SER:HB3	1.55	0.86
1:D:85:VAL:HG23	1:D:172:ILE:HD13	1.59	0.85
2:B:391:ARG:CD	2:B:400:TYR:HB3	2.03	0.85
2:K:398:SER:OG	2:K:399:ASN:HA	1.76	0.85
1:J:16:LEU:HD23	1:J:88:LEU:HB2	1.59	0.84
1:A:41:LYS:H	1:A:41:LYS:HD2	1.42	0.83
1:J:129:ARG:HD2	2:K:351:ASN:O	1.78	0.83
2:B:384:GLN:HE22	2:H:299:LEU:HD21	1.41	0.82
2:B:384:GLN:CD	2:H:299:LEU:HD11	2.05	0.81
1:A:54:GLN:HE22	1:A:59:THR:HG22	1.47	0.79
2:K:333:LEU:HA	2:K:335:PHE:HE1	1.47	0.79
2:K:309:LYS:CE	2:K:434:VAL:HG22	2.12	0.79
2:K:438:PHE:CZ	2:K:442:MET:HE2	2.18	0.79
2:K:427:GLY:HA2	2:K:428:LEU:HB2	1.65	0.79
2:K:427:GLY:HA2	2:K:428:LEU:CB	2.13	0.78
1:D:16:LEU:HD23	1:D:88:LEU:HB2	1.63	0.78
1:A:76:ILE:CG1	3:C:743:VAL:HG22	2.14	0.77
2:H:452:LYS:HE3	2:H:453:LEU:HD11	1.67	0.77
3:F:745:ILE:HD11	3:I:745:ILE:HD11	1.67	0.76
3:F:727:GLN:NE2	3:I:728:GLU:CA	2.48	0.76
2:B:453:LEU:O	2:H:442:MET:HE3	1.86	0.76
2:B:403:ILE:HG23	2:B:407:CYS:HB2	1.67	0.75
1:A:79:ALA:CB	3:C:746:MET:HE3	2.17	0.75
2:E:384:GLN:HE21	2:K:299:LEU:HD11	1.51	0.75
1:G:129[A]:ARG:HB3	2:H:354:SER:CB	2.17	0.75
2:E:427:GLY:HA2	2:E:428:LEU:CB	2.18	0.74
3:I:716:SER:N	3:I:720:LEU:HD11	2.03	0.74
3:F:731:ASN:OD1	3:I:731:ASN:ND2	2.20	0.73
2:B:409:TYR:OH	2:B:447:GLU:OE1	2.06	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:88:LEU:CD2	1:D:120:MET:HE2	2.18	0.73
1:A:11:LEU:HD13	1:A:61:LYS:HD3	1.69	0.73
2:B:395:GLY:HA3	2:B:396:ASP:C	2.12	0.73
1:D:54:GLN:NE2	1:D:59:THR:HG22	2.04	0.72
2:E:391:ARG:HD2	2:E:400:TYR:CB	2.20	0.72
1:G:129[B]:ARG:HB3	2:H:354:SER:CB	2.18	0.72
1:J:11:LEU:HD13	1:J:61:LYS:HD3	1.70	0.72
1:A:85:VAL:HG23	1:A:172:ILE:HD13	1.71	0.72
2:K:336:SER:HB2	2:K:382:THR:HA	1.70	0.72
2:H:301:LEU:CD1	2:H:442:MET:HE1	2.20	0.71
1:D:76:ILE:CD1	3:F:739:ASP:HA	2.14	0.71
1:J:48:PHE:CZ	3:L:754:GLU:CB	2.71	0.71
1:A:76:ILE:HG13	3:C:743:VAL:HG23	1.70	0.71
3:F:737:TYR:CB	1:G:44:ILE:HD12	2.20	0.71
2:K:343:ALA:O	2:K:346:GLU:HG2	1.90	0.71
3:F:724:ILE:HG12	3:I:724:ILE:CG1	2.21	0.70
2:K:420:TYR:CZ	2:K:440:GLU:HG2	2.26	0.70
3:F:724:ILE:HG12	3:I:724:ILE:HG13	1.73	0.70
2:B:372:GLU:HG2	2:B:373:CYS:H	1.55	0.70
2:B:438:PHE:CZ	2:B:442:MET:HE2	2.26	0.70
2:E:420:TYR:CE2	2:E:440:GLU:HG2	2.26	0.69
2:K:406:PHE:CZ	2:K:410:ARG:HD2	2.27	0.69
2:E:324:ILE:CD1	2:E:445:ARG:HD2	2.23	0.69
3:F:737:TYR:HB2	1:G:44:ILE:HD12	1.74	0.69
3:F:741:ILE:O	3:F:745:ILE:HG13	1.93	0.69
3:F:730:ILE:HG22	3:I:731:ASN:OD1	1.92	0.69
1:G:16:LEU:HD23	1:G:88:LEU:HB2	1.74	0.69
2:E:301:LEU:CD1	2:E:442:MET:HE1	2.22	0.69
2:B:393:LYS:CE	2:B:397:SER:HB3	2.22	0.68
1:G:38:LEU:HD22	2:H:428:LEU:HD22	1.76	0.67
2:E:301:LEU:HD13	2:E:442:MET:HE1	1.76	0.67
1:D:90:VAL:HA	1:D:122:VAL:O	1.94	0.67
1:G:90:VAL:HA	1:G:122:VAL:O	1.95	0.67
2:E:427:GLY:HA2	2:E:428:LEU:HB3	1.76	0.67
3:F:727:GLN:HE21	3:I:728:GLU:HA	1.58	0.67
3:F:727:GLN:HE22	3:I:728:GLU:HA	1.57	0.67
1:D:76:ILE:HG12	3:F:743:VAL:CG2	2.24	0.67
3:L:739:ASP:O	3:L:743:VAL:HG23	1.94	0.66
1:J:120:MET:HE1	1:J:164:ALA:HB1	1.76	0.66
2:E:453:LEU:O	2:K:442:MET:HE3	1.95	0.66
2:E:384:GLN:NE2	2:K:299:LEU:HD11	2.11	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:409:TYR:OH	2:E:447:GLU:OE1	2.11	0.66
2:H:351:ASN:ND2	2:H:351:ASN:O	2.27	0.66
2:H:333:LEU:HD12	2:H:345:LEU:HD23	1.79	0.65
1:A:54:GLN:NE2	1:A:59:THR:HG22	2.11	0.65
1:J:56:ASP:HB3	1:J:173:TYR:HE2	1.61	0.65
1:A:79:ALA:HB1	3:C:746:MET:HE3	1.78	0.65
3:F:737:TYR:CG	1:G:44:ILE:HD12	2.31	0.65
2:H:427:GLY:HA2	2:H:428:LEU:HB2	1.78	0.65
2:K:427:GLY:CA	2:K:428:LEU:HB2	2.26	0.65
2:H:427:GLY:HA2	2:H:428:LEU:CB	2.26	0.65
1:A:45:GLY:HA2	3:C:742:ILE:HD12	1.79	0.65
2:H:301:LEU:HD12	2:H:442:MET:HE1	1.79	0.65
2:K:433[B]:ASP:OD1	2:K:435:ASP:N	2.28	0.65
1:J:109:LEU:HD23	1:J:119:ILE:CD1	2.27	0.65
3:F:737:TYR:HB2	1:G:44:ILE:CD1	2.27	0.64
1:D:88:LEU:HD21	1:D:120:MET:HE2	1.78	0.64
4:G:200:GNP:H5'2	4:G:200:GNP:H8	1.79	0.64
1:J:43:THR:O	1:J:68:ALA:HB2	1.97	0.64
1:J:151:ILE:HD11	1:J:160:ASN:HB3	1.78	0.64
2:B:406:PHE:CZ	2:B:410:ARG:HD2	2.33	0.64
2:E:426:GLN:O	2:E:428:LEU:HB2	1.98	0.64
1:A:88:LEU:CD2	1:A:120:MET:HE2	2.28	0.64
1:A:129:ARG:CD	2:B:351:ASN:O	2.41	0.64
2:K:350:ASN:HB2	2:K:352:THR:HG23	1.79	0.63
2:K:420:TYR:CE2	2:K:440:GLU:HG2	2.32	0.63
1:D:53:ILE:CD1	1:D:55:VAL:HG23	2.27	0.63
1:D:76:ILE:HG22	1:D:76:ILE:O	1.98	0.63
2:B:427:GLY:HA2	2:B:428:LEU:CB	2.28	0.63
1:G:76:ILE:HG22	3:I:743:VAL:CG2	2.26	0.63
2:K:328:ASP:OD2	2:K:445:ARG:NE	2.30	0.63
2:H:328:ASP:HB3	2:H:448:MET:HE3	1.81	0.63
2:K:438:PHE:HZ	2:K:442:MET:HE2	1.64	0.62
1:G:125:LYS:HG2	4:G:200:GNP:C6	2.29	0.62
1:G:132:ARG:HD2	2:H:356:GLU:HG2	1.79	0.62
2:K:420:TYR:CE2	2:K:424:ILE:HD11	2.34	0.62
1:J:55:VAL:HG12	1:J:173:TYR:CD2	2.34	0.62
1:A:41:LYS:H	1:A:41:LYS:CD	2.03	0.62
3:I:745:ILE:HG23	3:I:752:ILE:HD12	1.82	0.61
2:B:372:GLU:HG2	2:B:373:CYS:N	2.16	0.61
1:D:174:ARG:O	1:D:178:GLN:HG3	2.00	0.61
1:J:109:LEU:HD23	1:J:119:ILE:HD13	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:ILE:CG1	3:C:743:VAL:CG2	2.69	0.61
2:B:391:ARG:CD	2:B:400:TYR:HD2	2.13	0.61
1:G:11:LEU:HD13	1:G:61:LYS:HD3	1.83	0.61
3:I:741:ILE:O	3:I:745:ILE:HG13	2.00	0.61
1:A:79:ALA:HB3	3:C:746:MET:HE3	1.82	0.61
2:B:442:MET:HE3	2:H:453:LEU:O	2.01	0.61
2:H:347:ALA:HA	2:H:352:THR:HG22	1.82	0.61
1:J:10:TYR:HB2	1:J:60:ILE:CD1	2.31	0.61
1:A:8:TYR:HA	1:A:59:THR:OG1	2.01	0.60
1:J:118:VAL:HG21	1:J:172:ILE:HG12	1.82	0.60
2:K:336:SER:CB	2:K:382:THR:HA	2.30	0.60
1:G:85:VAL:HG23	1:G:172:ILE:HD13	1.84	0.60
1:G:88:LEU:CD2	1:G:120:MET:HE2	2.32	0.60
1:G:38:LEU:HB2	2:H:428:LEU:CD2	2.31	0.60
1:A:45:GLY:CA	3:C:742:ILE:CD1	2.79	0.60
1:D:132:ARG:HD2	2:E:356:GLU:HG2	1.82	0.60
1:J:95:LYS:HE2	1:J:98:THR:HG23	1.84	0.60
1:D:76:ILE:HD11	3:F:739:ASP:C	2.25	0.60
1:J:171:GLU:O	1:J:175:ILE:HG13	2.01	0.60
2:K:392:ILE:HG23	2:K:403:ILE:HG13	1.83	0.60
3:F:737:TYR:HD2	3:I:738:ILE:HD13	1.67	0.60
1:A:45:GLY:HA2	3:C:742:ILE:CD1	2.32	0.59
3:F:727:GLN:HE22	3:I:728:GLU:CB	2.15	0.59
2:E:391:ARG:HD2	2:E:400:TYR:CG	2.37	0.59
1:J:125:LYS:HG2	4:J:200:GNP:C6	2.32	0.59
1:J:75:ALA:O	1:J:76:ILE:HG13	2.02	0.59
2:B:427:GLY:HA2	2:B:428:LEU:HB2	1.83	0.59
2:E:442:MET:HE3	2:K:453:LEU:O	2.02	0.59
2:H:347:ALA:HB1	2:H:353:LEU:HB2	1.83	0.59
1:A:76:ILE:CD1	3:C:742:ILE:CG2	2.76	0.59
2:E:386:LYS:NZ	2:K:296:GLU:CD	2.61	0.59
2:B:393:LYS:HG3	2:B:397:SER:CB	2.32	0.59
2:H:426:GLN:HG3	2:H:428:LEU:HD12	1.85	0.59
1:A:56:ASP:HB3	1:A:173:TYR:HE2	1.68	0.58
1:J:76:ILE:HG22	3:L:743:VAL:HG21	1.82	0.58
3:L:748:THR:OG1	3:L:749:ASN:N	2.35	0.58
1:G:120:MET:HG3	1:G:149:SER:HB2	1.84	0.58
2:K:452:LYS:HE3	2:K:453:LEU:HD11	1.84	0.58
2:H:345:LEU:HD13	2:H:345:LEU:C	2.28	0.58
2:H:347:ALA:CB	2:H:353:LEU:HB2	2.33	0.58
2:E:335:PHE:CE2	2:E:381:LEU:HD13	2.39	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:392:ILE:HD13	2:B:403:ILE:HG13	1.84	0.58
2:B:427:GLY:CA	2:B:428:LEU:HB2	2.33	0.58
2:E:324:ILE:HD11	2:E:445:ARG:HD2	1.86	0.58
1:J:76:ILE:HG21	3:L:743:VAL:HG23	1.77	0.58
2:K:422:ARG:O	2:K:426:GLN:HG2	2.04	0.58
3:C:739:ASP:O	3:C:743:VAL:HG23	2.04	0.58
1:J:99:TYR:CD2	1:J:134:VAL:HG11	2.39	0.58
1:D:76:ILE:HG12	3:F:743:VAL:HG23	1.85	0.57
1:G:53:ILE:HD11	1:G:166:GLN:CD	2.29	0.57
1:J:76:ILE:CG2	3:L:743:VAL:HG21	2.28	0.57
3:F:724:ILE:HD13	3:I:723:ALA:HB3	1.86	0.57
1:A:166:GLN:O	1:A:170:THR:HG23	2.05	0.57
2:E:386:LYS:HZ2	2:K:296:GLU:CD	2.12	0.57
2:E:406:PHE:CZ	2:E:410:ARG:HD2	2.40	0.57
2:E:452:LYS:HD2	2:K:327:GLU:OE1	2.04	0.57
2:K:350:ASN:CB	2:K:352:THR:HG23	2.35	0.57
1:A:8:TYR:CD2	1:A:61:LYS:HD2	2.40	0.57
2:E:391:ARG:CD	2:E:400:TYR:HB3	2.31	0.57
1:D:53:ILE:HD13	1:D:55:VAL:HG23	1.87	0.56
1:G:86:GLY:HA3	1:G:172:ILE:HD11	1.85	0.56
1:J:132:ARG:HD2	2:K:356:GLU:HG2	1.87	0.56
1:G:88:LEU:HD22	1:G:120:MET:HE2	1.87	0.56
2:B:384:GLN:NE2	2:H:299:LEU:HD11	2.20	0.56
2:B:420:TYR:CE2	2:B:440:GLU:HG2	2.40	0.56
1:G:21:GLY:H	4:G:200:GNP:HNB3	1.53	0.56
1:J:48:PHE:CE1	3:L:754:GLU:HB2	2.39	0.56
2:E:386:LYS:NZ	2:K:296:GLU:OE1	2.37	0.56
2:H:427:GLY:CA	2:H:428:LEU:HB2	2.36	0.56
2:K:437:MET:O	2:K:441:VAL:HG23	2.06	0.56
2:B:309:LYS:HE2	2:B:434:VAL:HG22	1.87	0.55
2:B:391:ARG:HD3	2:B:400:TYR:CD2	2.41	0.55
2:H:403:ILE:HG23	2:H:407:CYS:HB2	1.88	0.55
1:D:92:ASP:CG	1:D:125:LYS:HD2	2.31	0.55
2:H:351:ASN:C	2:H:351:ASN:HD22	2.14	0.55
1:D:32:THR:HG21	1:D:49:ALA:HB3	1.88	0.55
1:J:55:VAL:CG1	1:J:170:THR:HG22	2.36	0.55
1:G:55:VAL:HG12	1:G:173:TYR:CD2	2.41	0.55
1:J:55:VAL:HG12	1:J:173:TYR:HD2	1.69	0.55
2:E:403:ILE:HG23	2:E:407:CYS:HB2	1.88	0.55
1:G:56:ASP:HB3	1:G:173:TYR:HE2	1.71	0.55
2:H:301:LEU:HD13	2:H:442:MET:HE1	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:717:ARG:O	3:I:720:LEU:HD13	2.06	0.55
1:J:157:ASP:O	1:J:158:SER:HB2	2.06	0.55
2:B:373:CYS:SG	2:B:399:ASN:HB3	2.46	0.55
1:G:30:ARG:O	1:G:34:ASN:HA	2.07	0.55
2:H:417:PHE:HA	2:H:444:LEU:HD22	1.89	0.55
2:H:353:LEU:C	2:H:353:LEU:HD23	2.31	0.54
3:C:741:ILE:O	3:C:745:ILE:HG13	2.07	0.54
2:E:438:PHE:CZ	2:E:442:MET:HE2	2.42	0.54
2:K:308:TRP:CH2	2:K:421:ILE:HD12	2.42	0.54
2:E:424:ILE:HD13	2:E:437:MET:CE	2.37	0.54
1:A:45:GLY:CA	3:C:742:ILE:HD11	2.37	0.54
2:B:311:GLU:N	2:B:312:PRO:HD3	2.23	0.54
1:G:157:ASP:O	1:G:158:SER:HB2	2.08	0.54
2:H:452:LYS:HG2	2:H:453:LEU:HD12	1.88	0.54
1:J:10:TYR:HB2	1:J:60:ILE:HD13	1.90	0.54
2:K:433[B]:ASP:OD2	2:K:436:GLN:HG3	2.08	0.54
2:B:381:LEU:HD12	2:B:403:ILE:HD12	1.89	0.54
2:E:437:MET:O	2:E:441:VAL:HG23	2.07	0.54
1:J:76:ILE:HD12	1:J:76:ILE:O	2.08	0.54
2:B:353:LEU:HD12	2:B:394:LEU:HD23	1.89	0.54
1:J:39:GLU:OE2	2:K:432:GLN:HB2	2.08	0.54
2:H:392:ILE:HD13	2:H:403:ILE:HG13	1.90	0.54
1:D:54:GLN:HE22	1:D:59:THR:CG2	2.15	0.53
3:F:748:THR:OG1	3:F:749:ASN:N	2.41	0.53
2:B:369:SER:OG	2:B:372:GLU:OE2	2.22	0.53
2:E:330:PHE:N	2:E:331:PRO:HD2	2.23	0.53
1:A:74:ARG:C	1:A:76:ILE:H	2.16	0.53
1:D:172:ILE:O	1:D:176:VAL:HG22	2.07	0.53
2:K:423:TYR:O	2:K:427:GLY:HA2	2.09	0.53
3:I:748:THR:OG1	3:I:749:ASN:N	2.42	0.53
3:C:748:THR:OG1	3:C:749:ASN:N	2.42	0.53
1:D:76:ILE:HD11	3:F:739:ASP:O	2.09	0.53
3:F:724:ILE:CG1	3:I:724:ILE:HG13	2.38	0.53
1:D:53:ILE:C	1:D:53:ILE:HD12	2.34	0.53
3:F:727:GLN:HE22	3:I:728:GLU:CA	2.18	0.53
3:F:755:VAL:HG12	3:F:756:LYS:N	2.24	0.53
1:G:55:VAL:HG13	1:G:170:THR:HG22	1.91	0.53
2:H:452:LYS:HE3	2:H:453:LEU:CD1	2.37	0.53
1:J:55:VAL:HG13	1:J:170:THR:HG22	1.91	0.53
2:B:438:PHE:HZ	2:B:442:MET:HE2	1.74	0.53
1:D:119:ILE:CG2	1:D:148:LEU:HD22	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:55:VAL:CG1	1:J:173:TYR:HD2	2.21	0.52
1:D:88:LEU:HD22	1:D:120:MET:HE2	1.90	0.52
2:E:422:ARG:O	2:E:426:GLN:HG2	2.10	0.52
3:F:724:ILE:CD1	3:I:723:ALA:HB3	2.40	0.52
2:K:333:LEU:HA	2:K:335:PHE:CE1	2.36	0.52
2:B:301:LEU:CD1	2:B:442:MET:HE1	2.40	0.52
2:E:392:ILE:HD13	2:E:403:ILE:HG13	1.90	0.52
3:C:745:ILE:HG23	3:C:752:ILE:HD12	1.91	0.52
1:A:76:ILE:HD11	3:C:742:ILE:CB	2.39	0.52
1:A:46:VAL:N	3:C:742:ILE:HD13	2.25	0.52
2:K:316:ARG:NH2	2:K:349:GLU:OE2	2.43	0.52
2:K:409:TYR:OH	2:K:447:GLU:OE1	2.21	0.52
2:K:433[B]:ASP:OD1	2:K:435:ASP:HB2	2.09	0.52
1:D:92:ASP:OD1	1:D:125:LYS:HD2	2.10	0.52
2:H:406:PHE:CE1	2:H:454:GLY:HA2	2.45	0.52
1:A:86:GLY:HA3	1:A:172:ILE:HD11	1.91	0.51
2:E:350:ASN:HB2	2:E:352:THR:HG23	1.91	0.51
2:E:350:ASN:CB	2:E:352:THR:HG23	2.39	0.51
1:G:76:ILE:CG2	3:I:743:VAL:HG22	2.35	0.51
2:H:358:VAL:HB	2:H:391:ARG:HD2	1.92	0.51
3:L:744:ALA:O	3:L:748:THR:HG23	2.09	0.51
2:K:403:ILE:HG23	2:K:407:CYS:HB2	1.91	0.51
1:J:85:VAL:HG23	1:J:172:ILE:HD13	1.91	0.51
1:D:157:ASP:O	1:D:158:SER:HB2	2.10	0.51
1:G:48:PHE:O	3:I:754:GLU:HA	2.11	0.51
1:A:13:LYS:NZ	1:A:82:ARG:O	2.33	0.51
2:B:355:ILE:HG12	2:B:392:ILE:HD11	1.93	0.51
1:G:55:VAL:CG1	1:G:170:THR:HG22	2.41	0.51
1:A:88:LEU:HD22	1:A:120:MET:HE2	1.91	0.51
1:J:119:ILE:CG2	1:J:148:LEU:HD22	2.40	0.51
1:A:16:LEU:HD23	1:A:88:LEU:HB2	1.93	0.51
2:K:449:SER:O	2:K:453:LEU:HD13	2.11	0.51
1:A:119:ILE:CG2	1:A:148:LEU:HD22	2.40	0.51
3:F:724:ILE:HG12	3:I:724:ILE:HG12	1.92	0.51
1:D:76:ILE:HD13	3:F:742:ILE:HB	1.93	0.51
2:E:324:ILE:HD13	2:E:445:ARG:HD2	1.92	0.51
2:E:327:GLU:OE1	2:K:452:LYS:HD2	2.11	0.51
2:E:420:TYR:CZ	2:E:440:GLU:HG2	2.45	0.51
2:K:345:LEU:C	2:K:345:LEU:HD13	2.36	0.50
1:A:157:ASP:O	1:A:158:SER:HB2	2.11	0.50
2:E:424:ILE:HD13	2:E:437:MET:HE2	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:167:THR:O	1:J:171:GLU:HG3	2.11	0.50
1:A:115:SER:C	1:A:117:ILE:H	2.19	0.50
2:H:330:PHE:N	2:H:331:PRO:HD2	2.26	0.50
2:B:393:LYS:HG3	2:B:397:SER:HB3	1.92	0.50
3:C:736:ASP:O	3:C:740:ARG:HG3	2.11	0.50
1:D:120:MET:HE1	1:D:164:ALA:O	2.11	0.50
1:J:170:THR:O	1:J:174:ARG:HG2	2.11	0.50
1:J:16:LEU:HD21	1:J:88:LEU:HD12	1.94	0.50
1:J:86:GLY:HA3	1:J:172:ILE:HD11	1.94	0.50
2:E:301:LEU:HD12	2:E:442:MET:HE1	1.93	0.50
2:H:329:ILE:HD13	2:H:418:PHE:CE2	2.47	0.50
2:H:379:CYS:SG	2:H:382:THR:HG22	2.52	0.50
2:K:420:TYR:O	2:K:424:ILE:HG13	2.12	0.50
1:J:71:GLU:OE1	1:J:77:THR:HG22	2.11	0.49
1:A:93:ILE:HD11	1:A:123:GLY:HA3	1.94	0.49
3:C:724:ILE:HD12	3:C:725:GLN:N	2.27	0.49
1:D:10:TYR:HE1	1:D:58:LYS:HD3	1.77	0.49
1:G:125:LYS:HG2	4:G:200:GNP:C5	2.41	0.49
2:H:370:ALA:HB3	2:H:373:CYS:SG	2.51	0.49
2:K:376:PRO:O	2:K:387:SER:HA	2.11	0.49
1:A:54:GLN:HE22	1:A:59:THR:CG2	2.21	0.49
1:A:129:ARG:O	2:B:354:SER:HA	2.11	0.49
2:B:423:TYR:O	2:B:427:GLY:HA2	2.12	0.49
2:K:392:ILE:CG2	2:K:403:ILE:HG13	2.42	0.49
1:J:88:LEU:HD22	1:J:120:MET:HE2	1.94	0.49
3:F:724:ILE:CG1	3:I:724:ILE:CG1	2.89	0.49
1:A:51:ARG:NH2	1:A:162:GLU:OE2	2.46	0.49
1:J:130:HIS:CD2	1:J:131:LEU:HG	2.48	0.49
2:K:452:LYS:HG2	2:K:453:LEU:HD12	1.95	0.49
1:D:88:LEU:HD23	1:D:120:MET:HB3	1.94	0.49
2:H:438:PHE:CZ	2:H:442:MET:HE2	2.48	0.49
1:J:76:ILE:CB	3:L:743:VAL:HG22	2.38	0.49
2:H:328:ASP:CB	2:H:448:MET:HE3	2.43	0.48
2:K:330:PHE:N	2:K:331:PRO:HD2	2.28	0.48
2:E:427:GLY:CA	2:E:428:LEU:CB	2.91	0.48
2:E:428:LEU:HD23	2:E:428:LEU:O	2.13	0.48
2:H:353:LEU:HD23	2:H:354:SER:N	2.28	0.48
2:B:309:LYS:CE	2:B:434:VAL:HG22	2.43	0.48
2:E:423:TYR:O	2:E:427:GLY:HA2	2.14	0.48
1:G:56:ASP:HB3	1:G:173:TYR:CE2	2.48	0.48
1:G:56:ASP:CB	1:G:173:TYR:HE2	2.25	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:SER:HA	1:A:60:ILE:O	2.13	0.48
1:J:55:VAL:CG2	1:J:170:THR:HG22	2.44	0.48
1:J:118:VAL:CG2	1:J:172:ILE:HG12	2.42	0.48
2:K:353:LEU:HG	2:K:354:SER:N	2.28	0.48
2:E:301:LEU:HD21	2:E:323:LYS:HE2	1.96	0.48
2:B:333:LEU:HD12	2:B:345:LEU:HD13	1.96	0.48
1:G:58:LYS:HD2	1:G:173:TYR:OH	2.14	0.48
2:H:415:CYS:O	2:H:419[A]:THR:HG22	2.13	0.48
2:H:449:SER:O	2:H:453:LEU:HD13	2.14	0.48
1:J:151:ILE:HG12	1:J:152:GLU:N	2.29	0.48
2:E:340:LEU:HD23	2:E:381:LEU:HD23	1.96	0.48
4:G:200:GNP:H5'2	4:G:200:GNP:C8	2.44	0.48
1:A:79:ALA:HB3	3:C:746:MET:CE	2.44	0.48
2:B:348:VAL:HG22	2:B:353:LEU:HD22	1.96	0.48
1:D:38:LEU:HD22	2:E:428:LEU:HD22	1.96	0.48
2:H:331:PRO:HB2	2:H:452:LYS:HG3	1.95	0.48
1:A:118:VAL:CG2	1:A:172:ILE:HG12	2.44	0.47
2:B:425:GLN:OE1	2:B:426:GLN:HG3	2.14	0.47
1:D:123:GLY:O	1:D:153:THR:N	2.35	0.47
1:G:151:ILE:HG12	1:G:152:GLU:N	2.29	0.47
1:G:33:ARG:O	1:G:35:GLU:HG3	2.15	0.47
2:H:420:TYR:CE2	2:H:440:GLU:HG2	2.49	0.47
2:K:335:PHE:CD2	2:K:381:LEU:HB3	2.50	0.47
2:B:331:PRO:HB2	2:B:452:LYS:HG3	1.96	0.47
2:E:391:ARG:HD2	2:E:400:TYR:CD2	2.48	0.47
2:H:406:PHE:HE1	2:H:454:GLY:HA2	1.78	0.47
2:B:301:LEU:HD13	2:B:442:MET:HE1	1.95	0.47
2:K:305:PHE:HD1	2:K:320:PHE:CE1	2.32	0.47
1:A:132:ARG:HD2	2:B:356:GLU:HG2	1.97	0.47
2:E:331:PRO:HB2	2:E:452:LYS:HG3	1.97	0.47
3:C:722:GLU:OE1	3:C:722:GLU:N	2.48	0.47
1:D:120:MET:HE1	1:D:164:ALA:HB1	1.97	0.47
1:G:48:PHE:CZ	3:I:754:GLU:HB2	2.50	0.47
1:A:46:VAL:N	3:C:742:ILE:CD1	2.77	0.46
1:A:151:ILE:HD11	1:A:160:ASN:HB3	1.96	0.46
2:B:391:ARG:CD	2:B:400:TYR:CD2	2.94	0.46
2:B:408:ARG:O	2:B:412:THR:HG23	2.14	0.46
2:E:384:GLN:NE2	2:K:299:LEU:HD21	2.30	0.46
2:H:416:ASN:HA	2:H:419[A]:THR:HG22	1.97	0.46
2:K:406:PHE:CE1	2:K:454:GLY:HA2	2.50	0.46
1:A:92:ASP:CG	1:A:125:LYS:HD2	2.40	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:727:GLN:NE2	3:I:728:GLU:CB	2.78	0.46
2:H:311:GLU:N	2:H:312:PRO:HD3	2.30	0.46
1:A:151:ILE:HG12	1:A:152:GLU:N	2.31	0.46
2:E:392:ILE:HD13	2:E:403:ILE:CG1	2.45	0.46
1:J:172:ILE:O	1:J:176:VAL:HG23	2.15	0.46
2:B:373:CYS:HA	2:B:401:TYR:CD2	2.51	0.46
2:B:393:LYS:HG3	2:B:397:SER:HB2	1.97	0.46
2:E:432:GLN:HG2	2:E:436:GLN:HB2	1.97	0.46
1:A:85:VAL:C	1:A:117:ILE:HG13	2.40	0.46
1:A:106:LEU:HD11	1:A:148:LEU:HD11	1.98	0.46
2:B:393:LYS:HD2	2:B:400:TYR:CZ	2.51	0.46
1:G:23:GLY:HA2	4:G:200:GNP:O1A	2.15	0.46
1:J:8:TYR:CD2	1:J:61:LYS:HD2	2.50	0.46
1:D:37:ASN:HB3	1:D:40:SER:HB3	1.98	0.46
2:H:418:PHE:O	2:H:422[B]:ARG:HG3	2.15	0.46
2:H:423:TYR:O	2:H:427:GLY:HA2	2.16	0.46
1:J:129:ARG:CD	2:K:351:ASN:O	2.57	0.46
2:E:340:LEU:O	2:E:343:ALA:HB3	2.15	0.46
2:E:372:GLU:O	2:E:373:CYS:HB2	2.15	0.46
2:H:371:VAL:HG13	2:H:372:GLU:HG3	1.98	0.46
2:H:390:HIS:ND1	2:H:408:ARG:HD2	2.30	0.46
1:J:50:THR:HA	1:J:62:ALA:O	2.16	0.46
1:J:88:LEU:CD2	1:J:120:MET:HE2	2.46	0.46
2:B:391:ARG:HD3	2:B:400:TYR:CG	2.51	0.46
2:E:421:ILE:HD13	2:E:421:ILE:HA	1.73	0.46
3:F:745:ILE:HG23	3:F:752:ILE:HD12	1.97	0.46
1:D:124:ASN:HA	1:D:153:THR:O	2.16	0.45
2:B:381:LEU:HG	2:B:403:ILE:HD11	1.97	0.45
2:H:377:LYS:HD3	2:H:377:LYS:HA	1.80	0.45
1:J:55:VAL:CG1	1:J:173:TYR:CD2	2.97	0.45
2:B:353:LEU:HD21	2:B:411:ILE:HG21	1.97	0.45
1:D:39:GLU:OE2	2:E:432:GLN:HB2	2.17	0.45
1:D:51:ARG:NH1	1:D:162:GLU:OE2	2.50	0.45
1:J:48:PHE:CE1	3:L:754:GLU:CA	2.99	0.45
1:D:48:PHE:CZ	3:F:754:GLU:HB2	2.51	0.45
3:F:724:ILE:HD13	3:I:723:ALA:CB	2.45	0.45
1:D:65:TRP:HZ3	3:F:746:MET:CE	2.17	0.45
2:H:381:LEU:HD12	2:H:403:ILE:HD12	1.98	0.45
2:K:305:PHE:CE1	2:K:441:VAL:HG21	2.52	0.45
1:A:45:GLY:HA3	3:C:742:ILE:HD11	1.98	0.45
1:A:48:PHE:O	3:C:754:GLU:HA	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:395:GLY:HA3	2:B:397:SER:N	2.31	0.45
2:E:410:ARG:HG2	2:E:451:ALA:HB1	1.99	0.45
1:A:46:VAL:HG23	3:C:742:ILE:HG21	1.99	0.45
1:G:123:GLY:O	1:G:153:THR:N	2.42	0.45
1:J:31:PHE:HB2	1:J:165:PHE:CE2	2.52	0.45
2:E:444:LEU:HD13	2:E:444:LEU:HA	1.81	0.44
2:K:410:ARG:HG2	2:K:451:ALA:HB1	1.99	0.44
2:K:452:LYS:HE3	2:K:453:LEU:CD1	2.47	0.44
2:B:335:PHE:CE2	2:B:381:LEU:HD13	2.52	0.44
2:E:438:PHE:HZ	2:E:442:MET:HE2	1.83	0.44
1:A:76:ILE:CD1	3:C:742:ILE:HB	2.48	0.44
2:E:373:CYS:HA	2:E:401:TYR:HA	2.00	0.44
2:H:309:LYS:HE2	2:H:434:VAL:HG22	2.00	0.44
1:A:76:ILE:CG1	3:C:743:VAL:HG23	2.41	0.44
2:B:355:ILE:HG12	2:B:392:ILE:CD1	2.47	0.44
2:B:393:LYS:CG	2:B:397:SER:HB3	2.46	0.44
2:H:403:ILE:HG23	2:H:407:CYS:CB	2.48	0.44
1:J:32:THR:HG21	1:J:49:ALA:HB3	2.00	0.44
1:J:131:LEU:HD23	1:J:131:LEU:HA	1.74	0.44
2:B:403:ILE:HG23	2:B:407:CYS:CB	2.43	0.44
1:G:115:SER:C	1:G:117:ILE:H	2.25	0.44
2:H:409:TYR:CG	2:H:456:PHE:HB2	2.53	0.44
1:J:16:LEU:CD2	1:J:88:LEU:HB2	2.41	0.44
1:A:46:VAL:CG2	3:C:742:ILE:HG21	2.48	0.44
1:A:140:ARG:O	1:A:144:GLU:HG3	2.18	0.44
1:D:115:SER:C	1:D:117:ILE:H	2.26	0.43
2:H:403:ILE:HD13	2:H:403:ILE:HA	1.88	0.43
2:H:393:LYS:HB2	2:H:400:TYR:CE1	2.53	0.43
2:H:395:GLY:C	2:H:397:SER:H	2.25	0.43
1:A:37:ASN:HB3	1:A:40:SER:HB3	2.00	0.43
2:B:440:GLU:HG3	2:B:444:LEU:HD23	1.99	0.43
2:E:376:PRO:O	2:E:387:SER:HA	2.19	0.43
1:J:106:LEU:HD11	1:J:148:LEU:HD11	2.00	0.43
2:B:433:ASP:O	2:B:437:MET:HG3	2.18	0.43
1:G:50:THR:HG22	1:G:63:GLN:HA	2.00	0.43
2:B:306:ARG:HG3	2:B:434:VAL:HG11	2.00	0.43
1:J:55:VAL:HG22	1:J:170:THR:HG22	2.00	0.43
2:E:302:TYR:CD1	2:E:302:TYR:C	2.97	0.43
1:G:130:HIS:O	2:H:355:ILE:HB	2.18	0.43
1:J:45:GLY:HA2	3:L:742:ILE:CD1	2.49	0.43
2:K:421:ILE:HA	2:K:421:ILE:HD13	1.77	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:723:ALA:HB3	3:I:724:ILE:HD11	2.00	0.43
1:A:93:ILE:CD1	1:A:123:GLY:C	2.92	0.43
2:B:306:ARG:NH1	2:B:434:VAL:HG11	2.33	0.43
1:G:76:ILE:O	1:G:76:ILE:HG13	2.18	0.43
2:K:440:GLU:O	2:K:444:LEU:HD23	2.19	0.43
3:F:737:TYR:OH	1:G:45:GLY:HA3	2.18	0.43
2:E:449:SER:O	2:E:452:LYS:HB3	2.19	0.42
2:H:409:TYR:CE2	2:H:451:ALA:HB2	2.53	0.42
1:J:99:TYR:CD1	1:J:99:TYR:C	2.97	0.42
2:K:379:CYS:HB3	2:K:382:THR:O	2.19	0.42
1:A:132:ARG:O	2:B:357:PRO:HD3	2.19	0.42
2:B:330:PHE:N	2:B:331:PRO:HD2	2.34	0.42
3:I:734:LEU:HD23	3:I:734:LEU:HA	1.75	0.42
2:B:437:MET:O	2:B:440:GLU:HB3	2.20	0.42
1:D:180:GLN:O	1:D:180:GLN:HG2	2.18	0.42
2:B:324:ILE:HD13	2:B:417:PHE:CE1	2.53	0.42
1:G:32:THR:HG21	1:G:49:ALA:HB3	2.01	0.42
1:G:170:THR:O	1:G:174:ARG:HG3	2.19	0.42
2:H:421:ILE:HD13	2:H:421:ILE:HA	1.81	0.42
2:K:408:ARG:O	2:K:412:THR:HG23	2.19	0.42
3:F:731:ASN:ND2	3:I:727:GLN:NE2	2.68	0.42
2:B:395:GLY:HA3	2:B:397:SER:HA	2.01	0.42
1:D:14:VAL:HB	1:D:64:ILE:HD13	2.02	0.42
2:H:340:LEU:HD13	2:H:401:TYR:CE2	2.55	0.42
2:H:410:ARG:HG2	2:H:451:ALA:HB1	2.01	0.42
1:J:31:PHE:CZ	1:J:51:ARG:HD3	2.55	0.42
1:D:19:ASP:OD2	1:D:101:ASN:ND2	2.46	0.42
3:F:734:LEU:HD22	3:I:738:ILE:HD12	2.02	0.42
1:G:48:PHE:HA	1:G:64:ILE:O	2.18	0.42
2:H:432:GLN:HG2	2:H:436:GLN:HB2	2.00	0.42
1:J:125:LYS:HG2	4:J:200:GNP:C5	2.50	0.42
1:J:80:TYR:HA	3:L:746:MET:HE1	2.02	0.42
2:B:449:SER:O	2:B:452:LYS:HB3	2.20	0.42
1:D:38:LEU:HD21	2:E:423:TYR:CD2	2.55	0.42
3:F:739:ASP:O	3:F:743:VAL:HG23	2.19	0.42
1:G:168:ILE:O	1:G:172:ILE:HG13	2.20	0.42
2:K:404:SER:HB2	2:K:405:PRO:HD2	2.02	0.42
2:K:358:VAL:O	2:K:358:VAL:HG13	2.20	0.41
2:K:398:SER:CB	2:K:399:ASN:HA	2.48	0.41
1:A:142:PHE:CD1	1:A:142:PHE:C	2.98	0.41
2:H:340:LEU:O	2:H:344:VAL:HG23	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:48:PHE:HZ	3:L:754:GLU:HB2	1.67	0.41
1:A:118:VAL:HG21	1:A:172:ILE:HG12	2.02	0.41
2:E:324:ILE:HD13	2:E:445:ARG:CD	2.50	0.41
2:E:358:VAL:HG13	2:E:358:VAL:O	2.20	0.41
1:J:30:ARG:O	1:J:34:ASN:HA	2.19	0.41
2:B:306:ARG:HH12	2:B:434:VAL:HB	1.85	0.41
1:G:50:THR:CG2	1:G:63:GLN:HG3	2.50	0.41
1:J:45:GLY:HA2	3:L:742:ILE:HD12	2.01	0.41
1:J:48:PHE:CE1	3:L:754:GLU:N	2.88	0.41
2:B:393:LYS:CD	2:B:397:SER:HB3	2.49	0.41
1:D:81:TYR:OH	1:D:108:GLU:OE2	2.33	0.41
1:A:10:TYR:HB2	1:A:60:ILE:CD1	2.51	0.41
1:G:71:GLU:HB2	1:G:105:TRP:HH2	1.84	0.41
2:K:311:GLU:N	2:K:312:PRO:HD3	2.35	0.41
1:D:151:ILE:HG12	1:D:152:GLU:N	2.35	0.41
2:B:385:SER:C	2:B:386:LYS:HG2	2.45	0.41
1:D:88:LEU:HD23	1:D:88:LEU:HA	1.88	0.41
1:D:119:ILE:HG21	1:D:148:LEU:HD22	2.03	0.41
2:B:428:LEU:H	2:B:430:LYS:HE3	1.86	0.41
2:H:352:THR:OG1	2:H:395:GLY:HA3	2.20	0.41
1:J:95:LYS:HE2	1:J:95:LYS:HB3	1.54	0.41
1:J:168:ILE:O	1:J:172:ILE:HG13	2.20	0.41
1:A:124:ASN:O	1:A:125:LYS:HB2	2.21	0.41
2:B:355:ILE:HD12	2:B:412:THR:HG22	2.03	0.41
3:F:731:ASN:OD1	3:I:731:ASN:CG	2.64	0.41
1:J:10:TYR:HB2	1:J:60:ILE:HD12	2.03	0.41
1:J:13:LYS:HG3	1:J:65:TRP:CD1	2.57	0.41
2:H:302:TYR:OH	2:H:435:ASP:OD1	2.34	0.40
1:A:115:SER:C	1:A:117:ILE:N	2.79	0.40
1:D:120:MET:CE	1:D:164:ALA:HB1	2.51	0.40
1:D:128:LEU:HD21	4:D:200:GNP:N2	2.35	0.40
1:D:177:SER:O	1:D:180:GLN:HB3	2.19	0.40
1:G:151:ILE:HD11	1:G:160:ASN:HB3	2.03	0.40
2:H:420:TYR:CZ	2:H:440:GLU:HG2	2.55	0.40
1:G:10:TYR:HB2	1:G:60:ILE:HD12	2.03	0.40
1:G:71:GLU:HB3	1:G:72:ARG:H	1.70	0.40
1:J:19:ASP:O	1:J:22:VAL:HG13	2.21	0.40
2:K:426:GLN:O	2:K:428:LEU:HB2	2.21	0.40
1:A:125:LYS:HG2	4:A:200:GNP:C6	2.51	0.40
2:B:328:ASP:OD2	2:B:445:ARG:NE	2.47	0.40
1:D:122:VAL:HA	1:D:123:GLY:HA3	1.89	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:55:VAL:CG2	1:G:170:THR:HG22	2.51	0.40
1:J:56:ASP:HB3	1:J:173:TYR:CE2	2.48	0.40
1:J:56:ASP:CB	1:J:173:TYR:HE2	2.30	0.40
2:K:392:ILE:HG23	2:K:403:ILE:CG1	2.50	0.40
3:C:724:ILE:HD12	3:C:724:ILE:C	2.46	0.40
3:F:723:ALA:HB3	3:I:724:ILE:CD1	2.52	0.40
1:G:39:GLU:CD	2:H:430:LYS:H	2.30	0.40
1:G:93:ILE:HG13	1:G:94:ALA:N	2.35	0.40
2:H:340:LEU:CD1	2:H:401:TYR:CE2	3.04	0.40

All (6) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:371:VAL:CG1	1:J:160:ASN:OD1[3_554]	1.06	1.14
1:D:150:PHE:O	2:E:311:GLU:OE2[4_545]	1.66	0.54
1:D:57:GLY:CA	2:K:336:SER:O[4_545]	1.81	0.39
2:H:371:VAL:CG1	1:J:160:ASN:CG[3_554]	2.07	0.13
2:H:371:VAL:CB	1:J:160:ASN:OD1[3_554]	2.15	0.05
1:A:41:LYS:CB	3:L:733:ARG:NH1[1_455]	2.17	0.03

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	174/185 (94%)	169 (97%)	5 (3%)	0	100	100
1	D	173/185 (94%)	169 (98%)	4 (2%)	0	100	100
1	G	175/185 (95%)	168 (96%)	5 (3%)	2 (1%)	11	45
1	J	170/185 (92%)	165 (97%)	5 (3%)	0	100	100
2	B	157/195 (80%)	152 (97%)	5 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	E	156/195 (80%)	152 (97%)	4 (3%)	0	100	100
2	H	153/195 (78%)	150 (98%)	3 (2%)	0	100	100
2	K	156/195 (80%)	149 (96%)	7 (4%)	0	100	100
3	C	34/66 (52%)	33 (97%)	1 (3%)	0	100	100
3	F	35/66 (53%)	33 (94%)	2 (6%)	0	100	100
3	I	39/66 (59%)	36 (92%)	3 (8%)	0	100	100
3	L	36/66 (54%)	35 (97%)	1 (3%)	0	100	100
All	All	1458/1784 (82%)	1411 (97%)	45 (3%)	2 (0%)	48	82

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	76	ILE
1	G	73	TYR

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	142/157 (90%)	140 (99%)	2 (1%)	59	70
1	D	143/157 (91%)	141 (99%)	2 (1%)	59	70
1	G	143/157 (91%)	138 (96%)	5 (4%)	32	53
1	J	138/157 (88%)	135 (98%)	3 (2%)	45	64
2	B	141/175 (81%)	136 (96%)	5 (4%)	32	53
2	E	133/175 (76%)	129 (97%)	4 (3%)	36	57
2	H	140/175 (80%)	137 (98%)	3 (2%)	47	65
2	K	134/175 (77%)	129 (96%)	5 (4%)	30	51
3	C	25/56 (45%)	25 (100%)	0	100	100
3	F	21/56 (38%)	21 (100%)	0	100	100
3	I	25/56 (45%)	25 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	L	23/56 (41%)	22 (96%)	1 (4%)	26 48
All	All	1208/1552 (78%)	1178 (98%)	30 (2%)	42 62

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	TYR
1	A	41	LYS
2	B	307	LEU
2	B	323	LYS
2	B	332	CYS
2	B	352	THR
2	B	444	LEU
1	D	39	GLU
1	D	93	ILE
2	E	324	ILE
2	E	394	LEU
2	E	428	LEU
2	E	444	LEU
1	G	39	GLU
1	G	44	ILE
1	G	93	ILE
1	G	115	SER
1	G	149	SER
2	H	323	LYS
2	H	345	LEU
2	H	440	GLU
1	J	76	ILE
1	J	93	ILE
1	J	115	SER
2	K	332	CYS
2	K	345	LEU
2	K	421	ILE
2	K	440	GLU
2	K	444	LEU
3	L	724	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	436	GLN

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Mol	Chain	Res	Type
1	D	54	GLN
2	E	384	GLN
2	E	436	GLN
3	F	727	GLN
1	G	54	GLN
2	H	351	ASN
3	I	727	GLN
1	J	54	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

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	$\# Z > 2$
4	GNP	G	200	5	34,34,34	1.68	5 (14%)	47,54,54	1.24	4 (8%)
4	GNP	D	200	5	34,34,34	1.74	5 (14%)	47,54,54	1.07	3 (6%)
4	GNP	A	200	5	34,34,34	1.59	5 (14%)	47,54,54	0.98	3 (6%)
4	GNP	J	200	5	34,34,34	1.64	4 (11%)	47,54,54	1.20	4 (8%)

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
4	GNP	G	200	5	-	4/18/38/38	0/3/3/3
4	GNP	D	200	5	-	4/18/38/38	0/3/3/3
4	GNP	A	200	5	-	2/18/38/38	0/3/3/3
4	GNP	J	200	5	-	2/18/38/38	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	200	GNP	PA-O3A	-5.73	1.53	1.59
4	J	200	GNP	PA-O3A	-5.70	1.53	1.59
4	G	200	GNP	PA-O3A	-5.67	1.53	1.59
4	A	200	GNP	PA-O3A	-5.04	1.54	1.59
4	D	200	GNP	PB-O3A	-4.73	1.53	1.59
4	G	200	GNP	PB-O3A	-4.46	1.53	1.59
4	J	200	GNP	PB-O3A	-4.21	1.53	1.59
4	D	200	GNP	PG-O1G	4.00	1.52	1.46
4	A	200	GNP	PB-O3A	-3.75	1.54	1.59
4	A	200	GNP	PG-O1G	3.63	1.51	1.46
4	J	200	GNP	PG-O1G	3.53	1.51	1.46
4	G	200	GNP	PG-O1G	3.10	1.50	1.46
4	A	200	GNP	PB-O2B	-3.09	1.48	1.56
4	D	200	GNP	PB-O2B	-3.07	1.48	1.56
4	J	200	GNP	PB-O2B	-3.01	1.48	1.56
4	G	200	GNP	PB-O2B	-3.00	1.48	1.56
4	G	200	GNP	PG-O2G	-2.22	1.50	1.56
4	D	200	GNP	PG-O2G	-2.14	1.51	1.56
4	A	200	GNP	PG-O2G	-2.13	1.51	1.56

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	J	200	GNP	O2B-PB-O1B	4.59	119.72	109.87
4	G	200	GNP	O2G-PG-O1G	-4.43	102.34	113.45
4	D	200	GNP	O2B-PB-O1B	3.90	118.24	109.87
4	G	200	GNP	O2G-PG-O3G	3.78	117.76	107.59
4	D	200	GNP	O2G-PG-O1G	-3.75	104.05	113.45
4	J	200	GNP	O2G-PG-O1G	-3.62	104.36	113.45
4	J	200	GNP	O2G-PG-O3G	3.52	117.06	107.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	200	GNP	O2G-PG-O1G	-3.17	105.50	113.45
4	D	200	GNP	O2G-PG-O3G	3.09	115.89	107.59
4	G	200	GNP	O2B-PB-O1B	3.08	116.48	109.87
4	A	200	GNP	O2G-PG-O3G	2.92	115.44	107.59
4	A	200	GNP	O2B-PB-O1B	2.54	115.32	109.87
4	G	200	GNP	O2A-PA-O3A	2.28	113.44	107.27
4	J	200	GNP	O1B-PB-N3B	-2.15	108.60	111.77

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	200	GNP	PG-N3B-PB-O1B
4	A	200	GNP	PG-N3B-PB-O3A
4	D	200	GNP	PG-N3B-PB-O1B
4	D	200	GNP	PA-O3A-PB-O2B
4	G	200	GNP	PG-N3B-PB-O1B
4	G	200	GNP	PA-O3A-PB-O2B
4	J	200	GNP	PG-N3B-PB-O1B
4	G	200	GNP	O4'-C4'-C5'-O5'
4	J	200	GNP	PG-N3B-PB-O3A
4	D	200	GNP	PA-O3A-PB-O1B
4	G	200	GNP	PA-O3A-PB-O1B
4	D	200	GNP	PG-N3B-PB-O3A

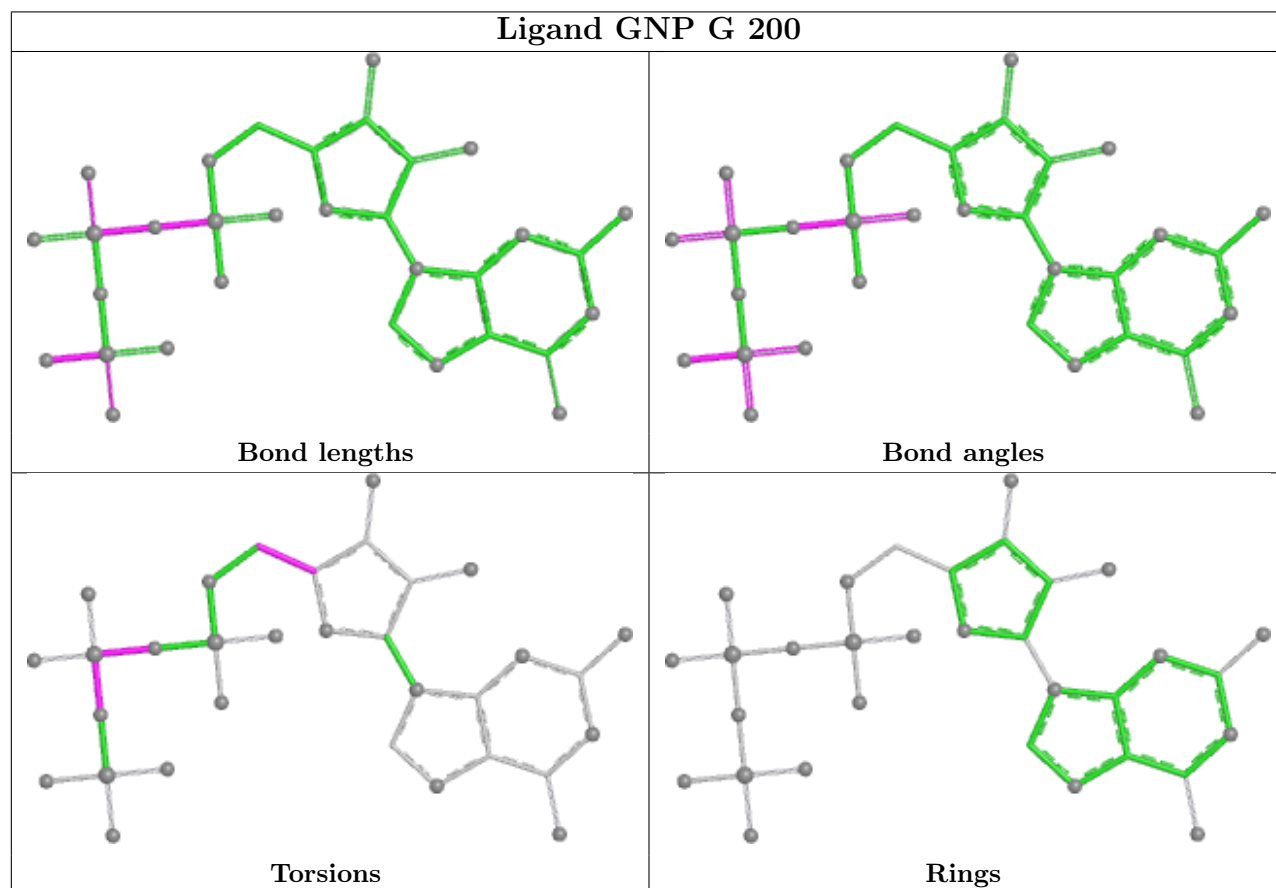
There are no ring outliers.

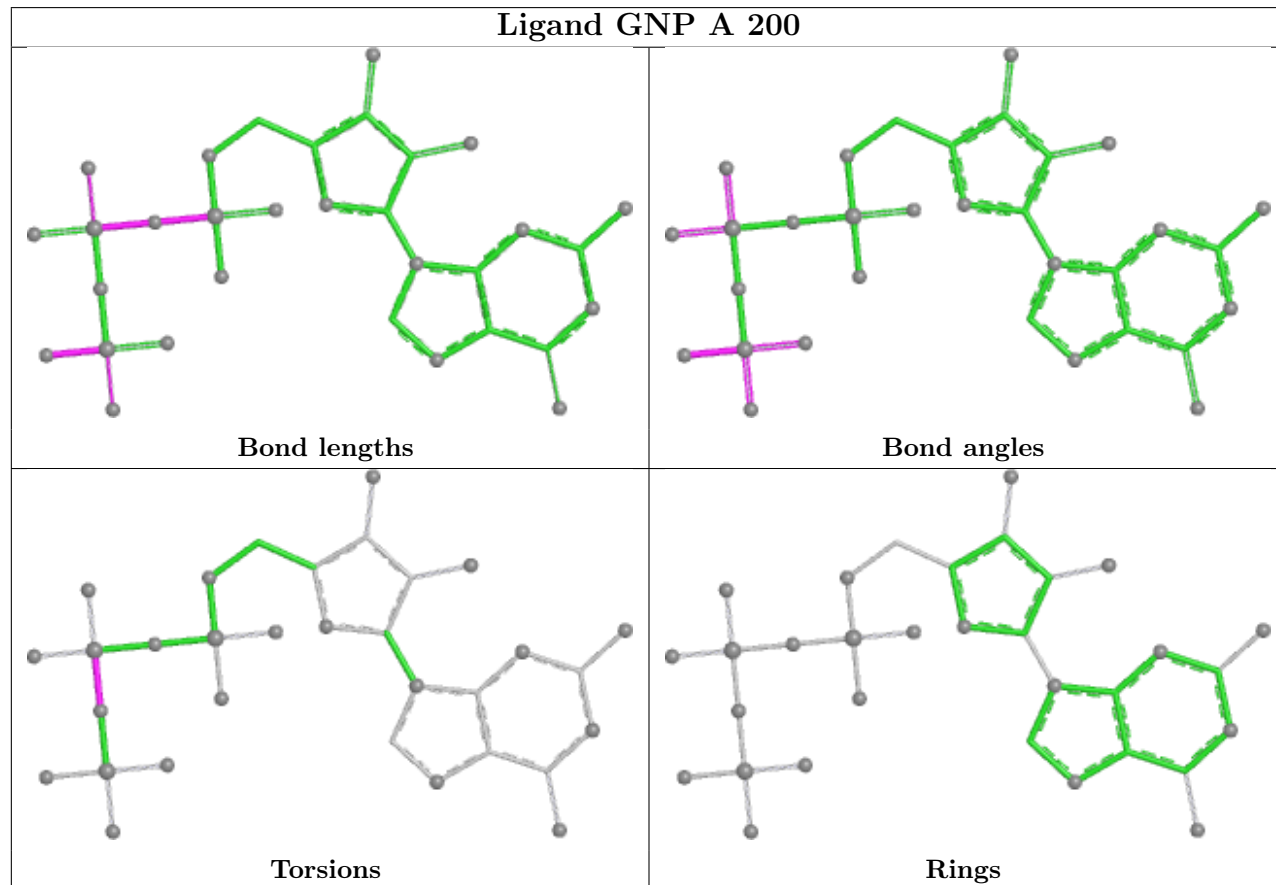
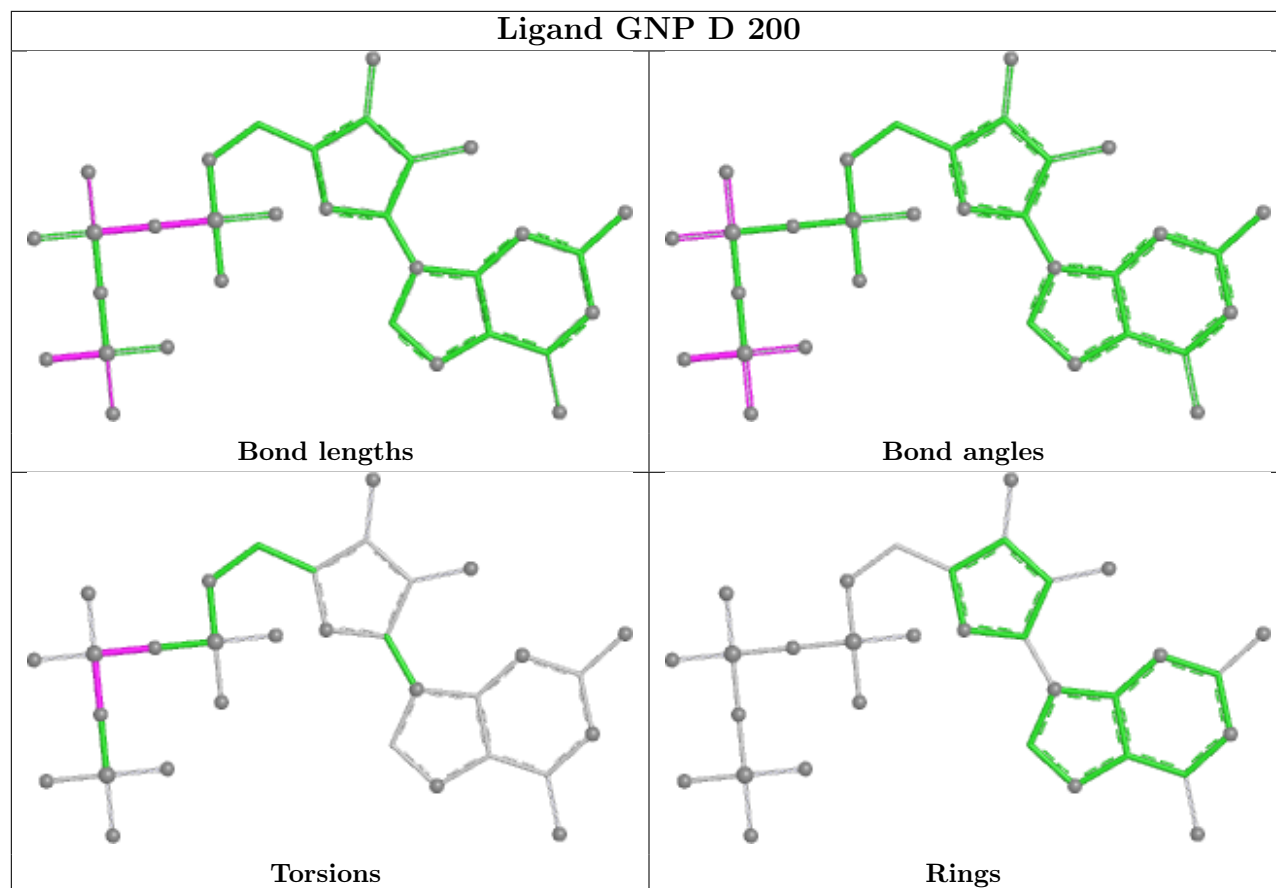
4 monomers are involved in 10 short contacts:

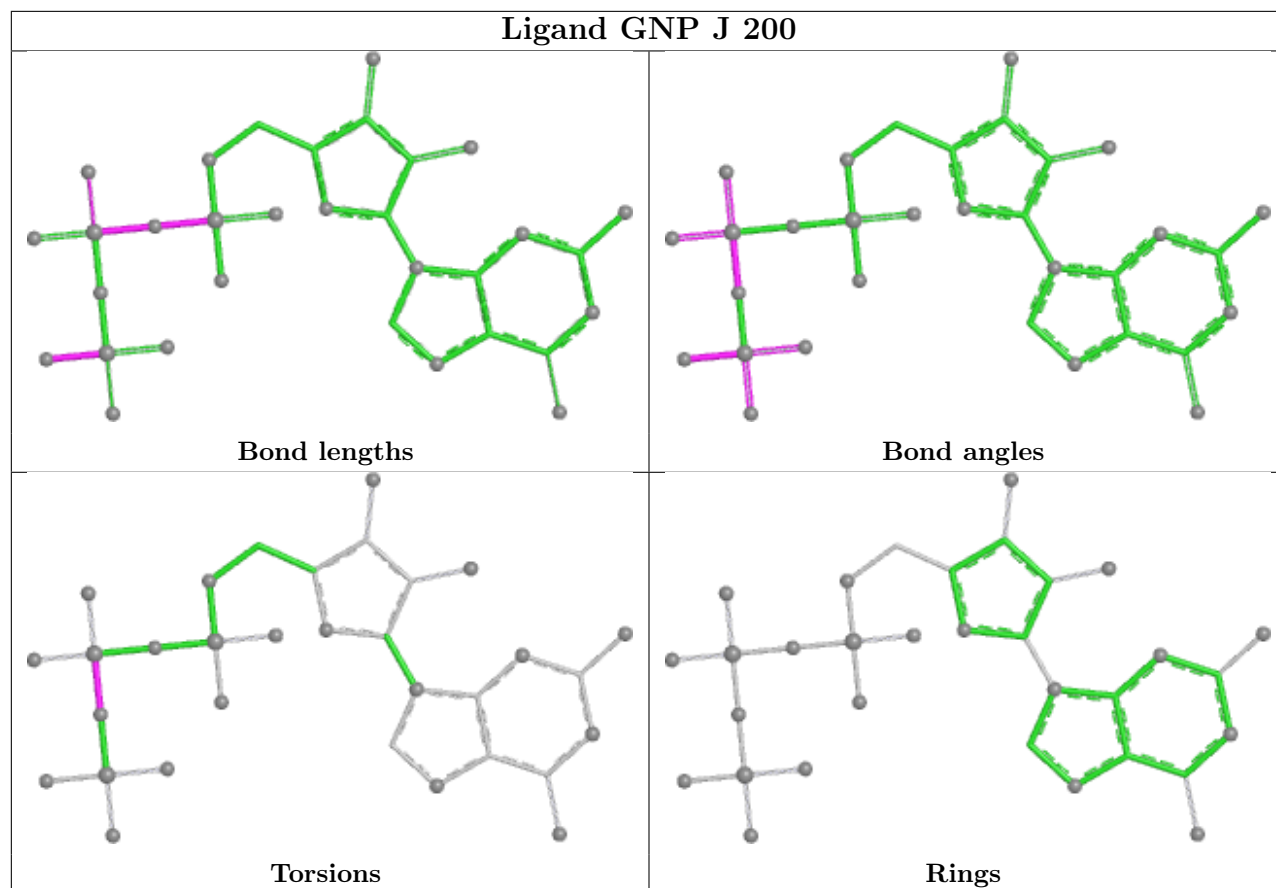
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	G	200	GNP	6	0
4	D	200	GNP	1	0
4	A	200	GNP	1	0
4	J	200	GNP	2	0

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	176/185 (95%)	0.28	7 (3%) 42 35	144, 166, 205, 258	0
1	D	175/185 (94%)	0.15	2 (1%) 78 62	95, 158, 214, 253	2 (1%)
1	G	175/185 (94%)	0.14	1 (0%) 85 72	81, 141, 188, 226	2 (1%)
1	J	174/185 (94%)	0.14	2 (1%) 78 62	121, 149, 200, 235	0
2	B	161/195 (82%)	0.30	5 (3%) 51 39	131, 166, 225, 273	0
2	E	162/195 (83%)	0.17	4 (2%) 58 45	120, 169, 205, 238	0
2	H	155/195 (79%)	0.23	2 (1%) 75 59	82, 149, 174, 201	2 (1%)
2	K	159/195 (81%)	0.29	3 (1%) 66 51	94, 162, 208, 218	1 (0%)
3	C	36/66 (54%)	0.26	0 100 100	139, 164, 221, 224	0
3	F	37/66 (56%)	0.20	1 (2%) 56 43	122, 143, 197, 209	0
3	I	41/66 (62%)	0.72	6 (14%) 6 9	133, 153, 227, 236	0
3	L	38/66 (57%)	0.30	1 (2%) 57 43	141, 168, 219, 231	0
All	All	1489/1784 (83%)	0.23	34 (2%) 61 47	81, 159, 210, 273	7 (0%)

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	I	720	LEU	5.5
1	A	28	LEU	4.2
1	A	11	LEU	3.9
2	E	396	ASP	3.9
1	J	75	ALA	3.3
2	B	418	PHE	3.3
2	K	450	LEU	3.0
1	A	62	ALA	2.8
1	A	12	PHE	2.7
3	L	719	GLU	2.6
2	H	369	SER	2.6

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Mol	Chain	Res	Type	RSRZ
2	E	455	TYR	2.6
1	A	32	THR	2.6
2	E	313	THR	2.5
2	H	402	TYR	2.5
3	I	719	GLU	2.4
1	A	27	LEU	2.4
2	B	454	GLY	2.4
1	A	64	ILE	2.4
3	I	716	SER	2.4
3	I	751	SER	2.3
2	K	403	ILE	2.3
2	B	326	GLN	2.3
2	B	344	VAL	2.2
2	E	395	GLY	2.2
3	I	723	ALA	2.2
1	D	69	GLY	2.2
1	G	173	TYR	2.2
1	D	76	ILE	2.1
2	K	459	GLU	2.1
2	B	348	VAL	2.1
3	F	742	ILE	2.1
3	I	721	MET	2.0
1	J	69	GLY	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](#)

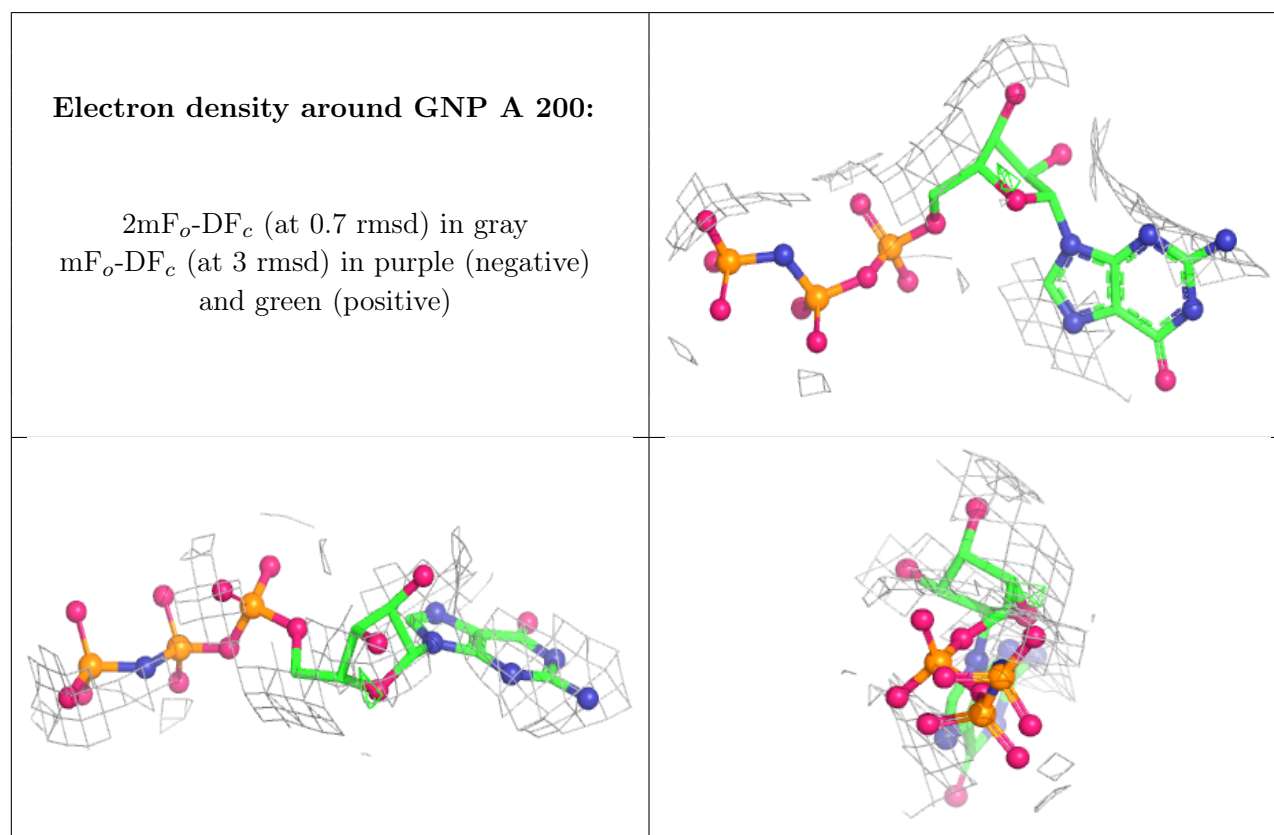
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists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

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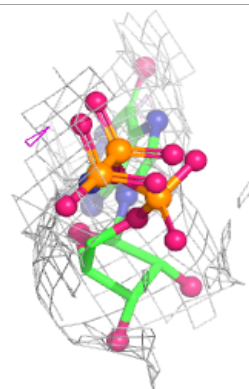
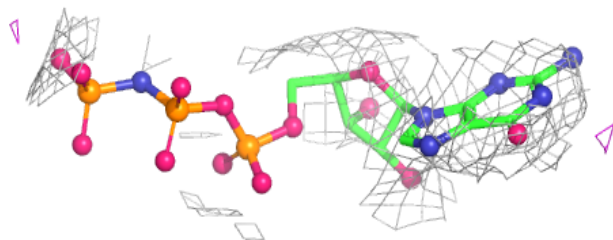
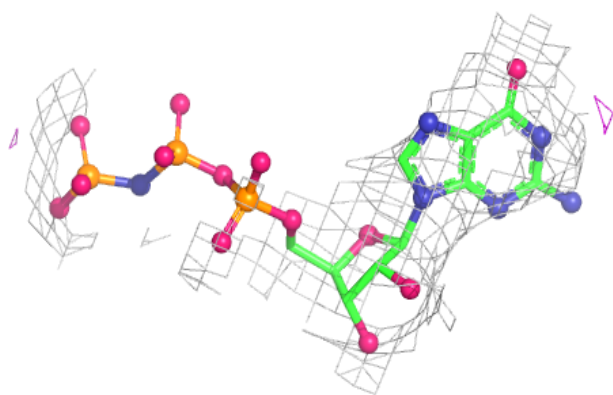
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	GNP	A	200	32/32	0.90	0.08	140,150,162,164	0
5	MG	J	201	1/1	0.91	0.07	131,131,131,131	0
4	GNP	D	200	32/32	0.92	0.09	130,143,153,161	0
4	GNP	J	200	32/32	0.93	0.08	117,129,146,148	0
5	MG	A	201	1/1	0.94	0.08	145,145,145,145	0
4	GNP	G	200	32/32	0.96	0.06	113,129,142,149	0
5	MG	G	201	1/1	0.97	0.09	111,111,111,111	0
5	MG	D	201	1/1	0.98	0.05	132,132,132,132	0

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

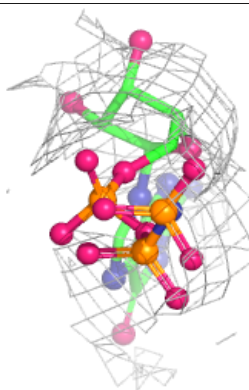
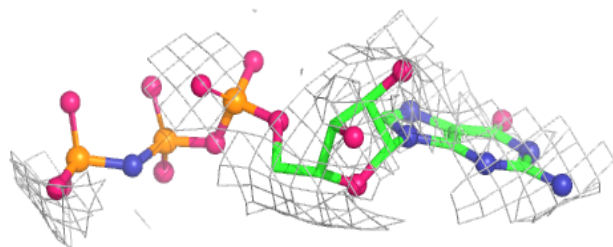
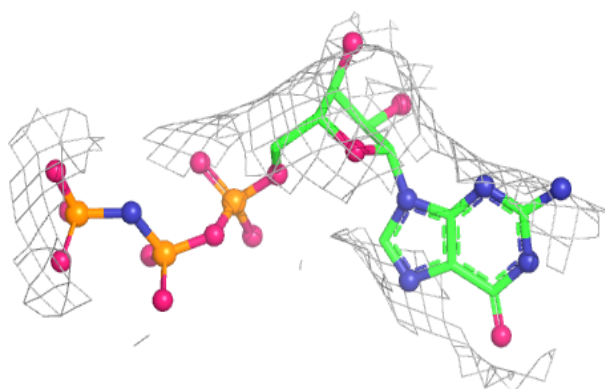


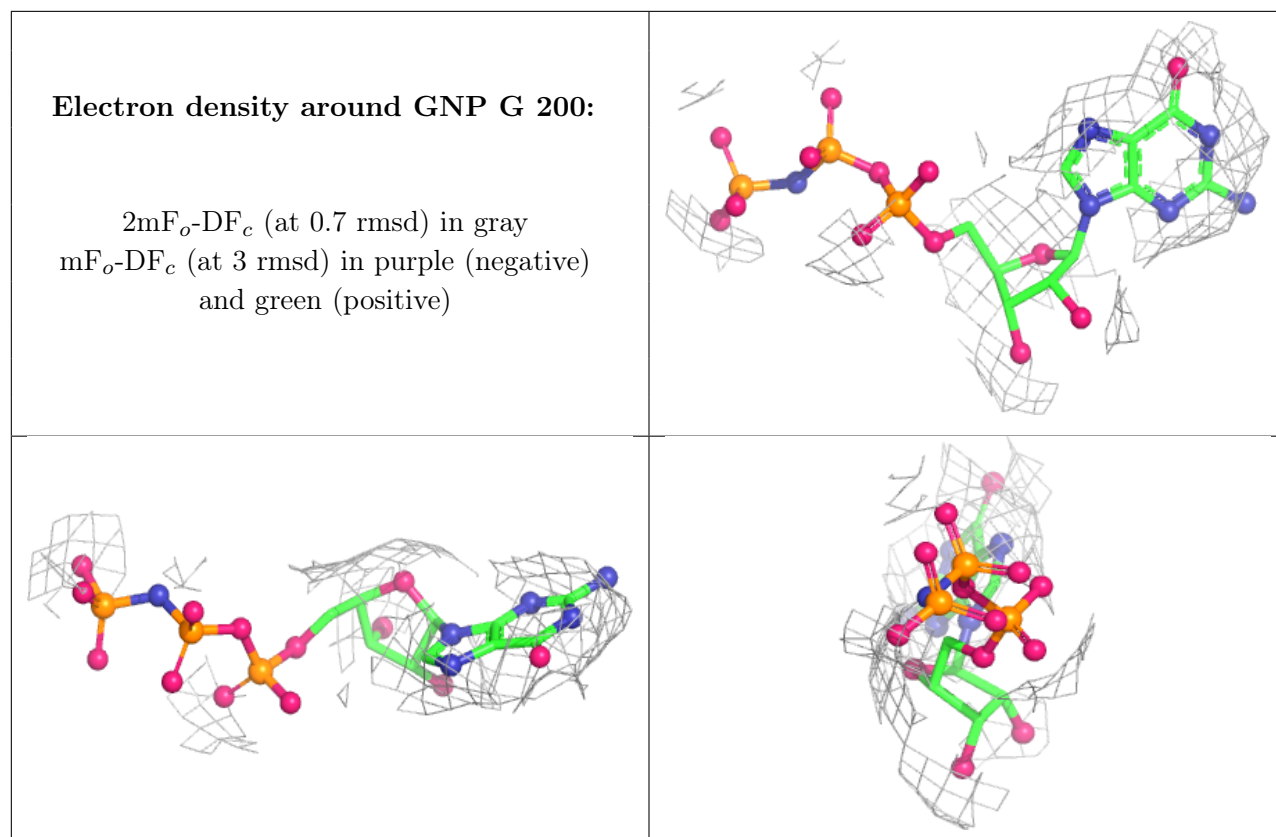
Electron density around GNP D 200:

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

**Electron density around GNP J 200:**

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





6.5 Other polymers [i](#)

There are no such residues in this entry.