



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 02:12 PM UTC

PDB ID : 3QOF / pdb_00003qof
Title : Crystal structure of the cytosolic domain of human atlastin-1 in complex with GDP, orthorhombic form
Authors : Liu, X.
Deposited on : 2011-02-09
Resolution : 2.80 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

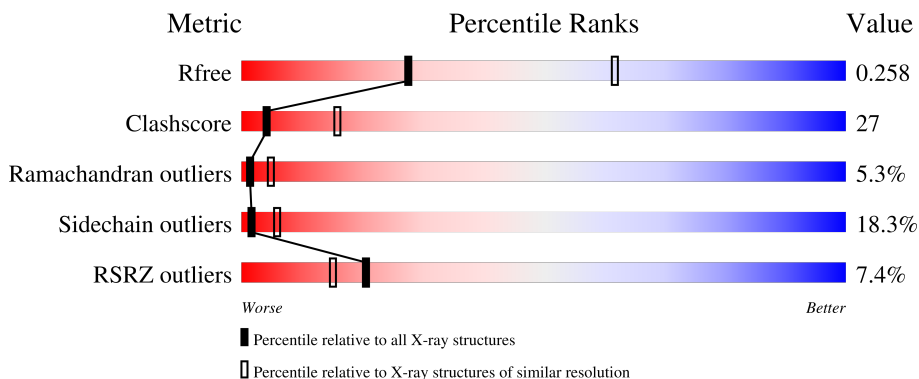
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	459	
1	B	459	
1	C	459	
1	D	459	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12574 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 Atlastin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	404	Total	C	N	O	S	0	0	0
			3108	1994	518	583	13			
1	B	397	Total	C	N	O	S	0	0	0
			3079	1979	509	578	13			
1	C	405	Total	C	N	O	S	0	0	0
			3116	2003	518	582	13			
1	D	398	Total	C	N	O	S	0	0	0
			3029	1948	508	561	12			

There are 116 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	expression tag	UNP Q8WXF7
A	-10	HIS	-	expression tag	UNP Q8WXF7
A	-9	HIS	-	expression tag	UNP Q8WXF7
A	-8	HIS	-	expression tag	UNP Q8WXF7
A	-7	HIS	-	expression tag	UNP Q8WXF7
A	-6	HIS	-	expression tag	UNP Q8WXF7
A	-5	HIS	-	expression tag	UNP Q8WXF7
A	-4	GLY	-	expression tag	UNP Q8WXF7
A	-3	SER	-	expression tag	UNP Q8WXF7
A	-2	HIS	-	expression tag	UNP Q8WXF7
A	-1	MET	-	expression tag	UNP Q8WXF7
A	0	ALA	-	expression tag	UNP Q8WXF7
A	1	SER	-	expression tag	UNP Q8WXF7
A	2	ALA	-	expression tag	UNP Q8WXF7
A	3	LYS	-	expression tag	UNP Q8WXF7
A	4	ASN	-	expression tag	UNP Q8WXF7
A	5	ARG	-	expression tag	UNP Q8WXF7
A	6	ARG	-	expression tag	UNP Q8WXF7
A	7	ASP	-	expression tag	UNP Q8WXF7
A	8	ARG	-	expression tag	UNP Q8WXF7
A	9	ASN	-	expression tag	UNP Q8WXF7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	10	SER	-	expression tag	UNP Q8WXF7
A	11	TRP	-	expression tag	UNP Q8WXF7
A	12	GLY	-	expression tag	UNP Q8WXF7
A	13	GLY	-	expression tag	UNP Q8WXF7
A	14	PHE	-	expression tag	UNP Q8WXF7
A	15	SER	-	expression tag	UNP Q8WXF7
A	16	GLU	-	expression tag	UNP Q8WXF7
A	17	LYS	-	expression tag	UNP Q8WXF7
B	-11	MET	-	expression tag	UNP Q8WXF7
B	-10	HIS	-	expression tag	UNP Q8WXF7
B	-9	HIS	-	expression tag	UNP Q8WXF7
B	-8	HIS	-	expression tag	UNP Q8WXF7
B	-7	HIS	-	expression tag	UNP Q8WXF7
B	-6	HIS	-	expression tag	UNP Q8WXF7
B	-5	HIS	-	expression tag	UNP Q8WXF7
B	-4	GLY	-	expression tag	UNP Q8WXF7
B	-3	SER	-	expression tag	UNP Q8WXF7
B	-2	HIS	-	expression tag	UNP Q8WXF7
B	-1	MET	-	expression tag	UNP Q8WXF7
B	0	ALA	-	expression tag	UNP Q8WXF7
B	1	SER	-	expression tag	UNP Q8WXF7
B	2	ALA	-	expression tag	UNP Q8WXF7
B	3	LYS	-	expression tag	UNP Q8WXF7
B	4	ASN	-	expression tag	UNP Q8WXF7
B	5	ARG	-	expression tag	UNP Q8WXF7
B	6	ARG	-	expression tag	UNP Q8WXF7
B	7	ASP	-	expression tag	UNP Q8WXF7
B	8	ARG	-	expression tag	UNP Q8WXF7
B	9	ASN	-	expression tag	UNP Q8WXF7
B	10	SER	-	expression tag	UNP Q8WXF7
B	11	TRP	-	expression tag	UNP Q8WXF7
B	12	GLY	-	expression tag	UNP Q8WXF7
B	13	GLY	-	expression tag	UNP Q8WXF7
B	14	PHE	-	expression tag	UNP Q8WXF7
B	15	SER	-	expression tag	UNP Q8WXF7
B	16	GLU	-	expression tag	UNP Q8WXF7
B	17	LYS	-	expression tag	UNP Q8WXF7
C	-11	MET	-	expression tag	UNP Q8WXF7
C	-10	HIS	-	expression tag	UNP Q8WXF7
C	-9	HIS	-	expression tag	UNP Q8WXF7
C	-8	HIS	-	expression tag	UNP Q8WXF7
C	-7	HIS	-	expression tag	UNP Q8WXF7

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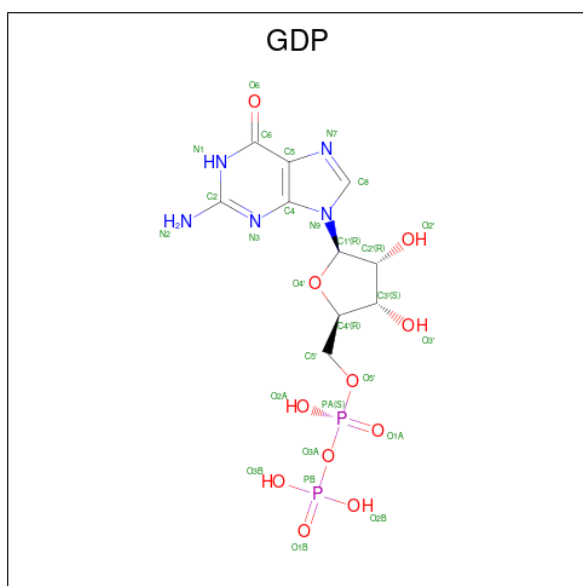
Chain	Residue	Modelled	Actual	Comment	Reference
C	-6	HIS	-	expression tag	UNP Q8WXF7
C	-5	HIS	-	expression tag	UNP Q8WXF7
C	-4	GLY	-	expression tag	UNP Q8WXF7
C	-3	SER	-	expression tag	UNP Q8WXF7
C	-2	HIS	-	expression tag	UNP Q8WXF7
C	-1	MET	-	expression tag	UNP Q8WXF7
C	0	ALA	-	expression tag	UNP Q8WXF7
C	1	SER	-	expression tag	UNP Q8WXF7
C	2	ALA	-	expression tag	UNP Q8WXF7
C	3	LYS	-	expression tag	UNP Q8WXF7
C	4	ASN	-	expression tag	UNP Q8WXF7
C	5	ARG	-	expression tag	UNP Q8WXF7
C	6	ARG	-	expression tag	UNP Q8WXF7
C	7	ASP	-	expression tag	UNP Q8WXF7
C	8	ARG	-	expression tag	UNP Q8WXF7
C	9	ASN	-	expression tag	UNP Q8WXF7
C	10	SER	-	expression tag	UNP Q8WXF7
C	11	TRP	-	expression tag	UNP Q8WXF7
C	12	GLY	-	expression tag	UNP Q8WXF7
C	13	GLY	-	expression tag	UNP Q8WXF7
C	14	PHE	-	expression tag	UNP Q8WXF7
C	15	SER	-	expression tag	UNP Q8WXF7
C	16	GLU	-	expression tag	UNP Q8WXF7
C	17	LYS	-	expression tag	UNP Q8WXF7
D	-11	MET	-	expression tag	UNP Q8WXF7
D	-10	HIS	-	expression tag	UNP Q8WXF7
D	-9	HIS	-	expression tag	UNP Q8WXF7
D	-8	HIS	-	expression tag	UNP Q8WXF7
D	-7	HIS	-	expression tag	UNP Q8WXF7
D	-6	HIS	-	expression tag	UNP Q8WXF7
D	-5	HIS	-	expression tag	UNP Q8WXF7
D	-4	GLY	-	expression tag	UNP Q8WXF7
D	-3	SER	-	expression tag	UNP Q8WXF7
D	-2	HIS	-	expression tag	UNP Q8WXF7
D	-1	MET	-	expression tag	UNP Q8WXF7
D	0	ALA	-	expression tag	UNP Q8WXF7
D	1	SER	-	expression tag	UNP Q8WXF7
D	2	ALA	-	expression tag	UNP Q8WXF7
D	3	LYS	-	expression tag	UNP Q8WXF7
D	4	ASN	-	expression tag	UNP Q8WXF7
D	5	ARG	-	expression tag	UNP Q8WXF7
D	6	ARG	-	expression tag	UNP Q8WXF7

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Chain	Residue	Modelled	Actual	Comment	Reference
D	7	ASP	-	expression tag	UNP Q8WXF7
D	8	ARG	-	expression tag	UNP Q8WXF7
D	9	ASN	-	expression tag	UNP Q8WXF7
D	10	SER	-	expression tag	UNP Q8WXF7
D	11	TRP	-	expression tag	UNP Q8WXF7
D	12	GLY	-	expression tag	UNP Q8WXF7
D	13	GLY	-	expression tag	UNP Q8WXF7
D	14	PHE	-	expression tag	UNP Q8WXF7
D	15	SER	-	expression tag	UNP Q8WXF7
D	16	GLU	-	expression tag	UNP Q8WXF7
D	17	LYS	-	expression tag	UNP Q8WXF7

- Molecule 2 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	B	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	C	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	D	1	Total	C	N	O	P	0	0
			28	10	5	11	2		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mg 1 1	0	0
3	B	1	Total Mg 1 1	0	0
3	C	1	Total Mg 1 1	0	0
3	D	1	Total Mg 1 1	0	0

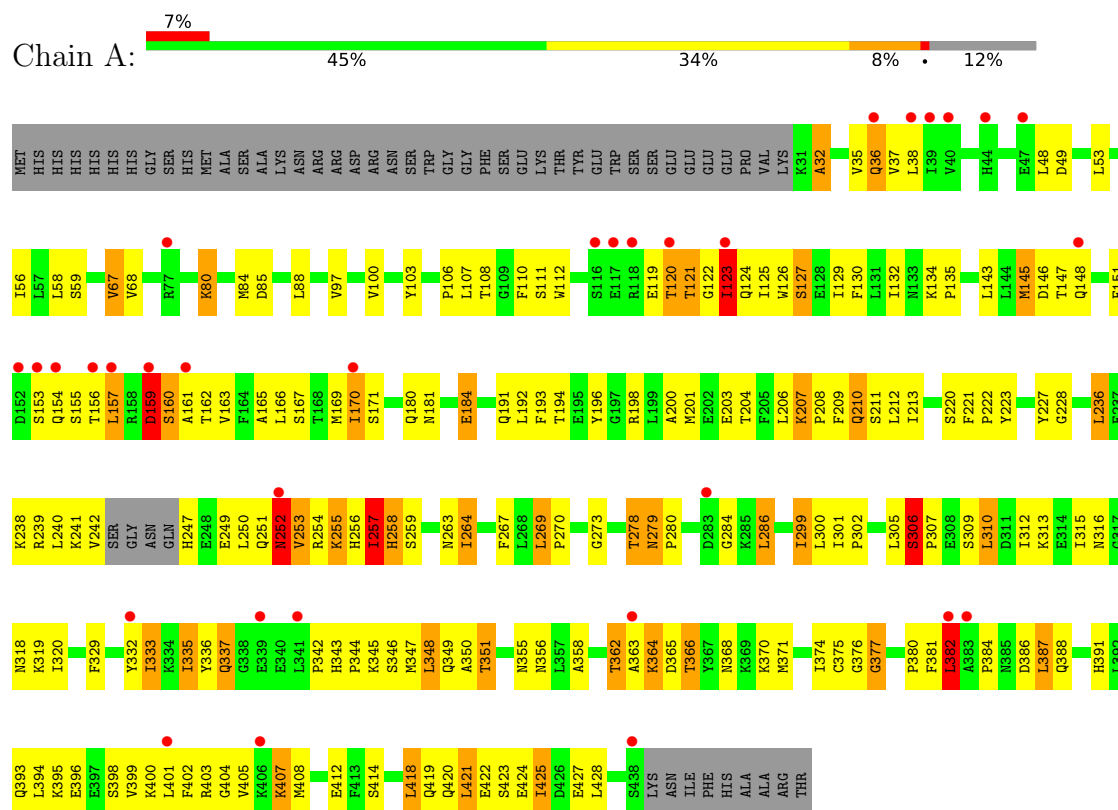
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	37	Total O 37 37	0	0
4	B	29	Total O 29 29	0	0
4	C	37	Total O 37 37	0	0
4	D	23	Total O 23 23	0	0

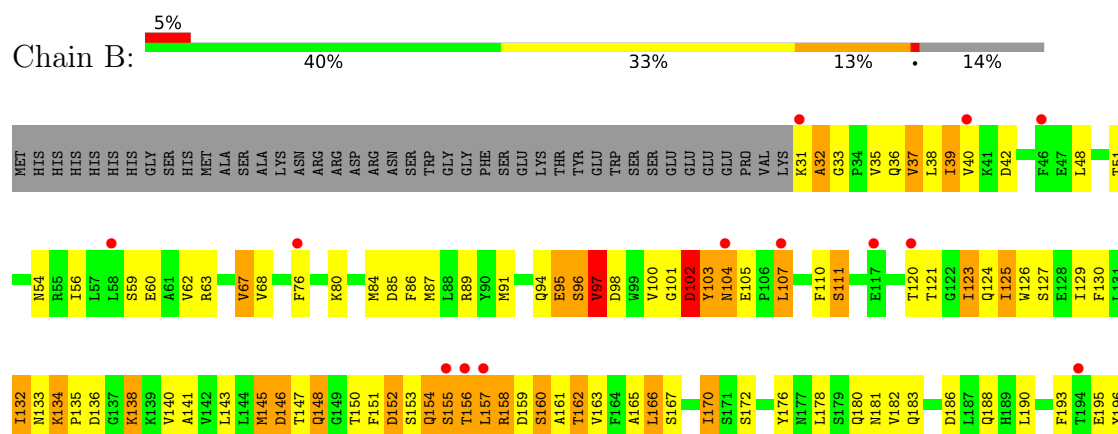
3 Residue-property plots

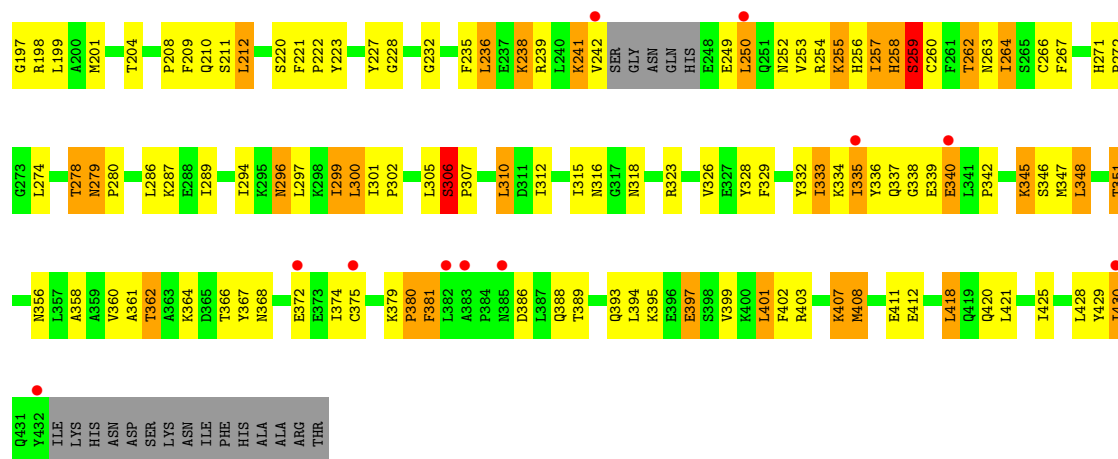
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Atlastin-1

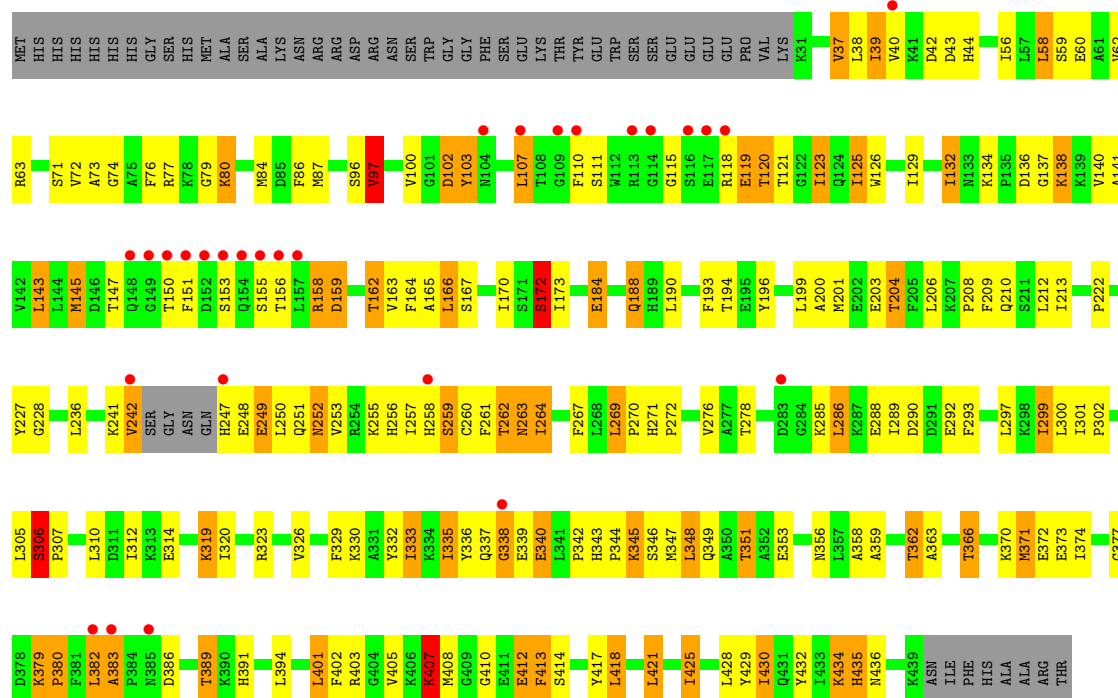
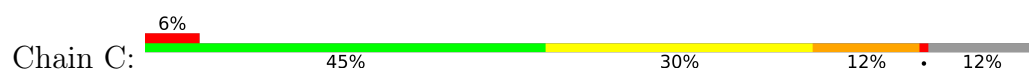


• Molecule 1: Atlastin-1

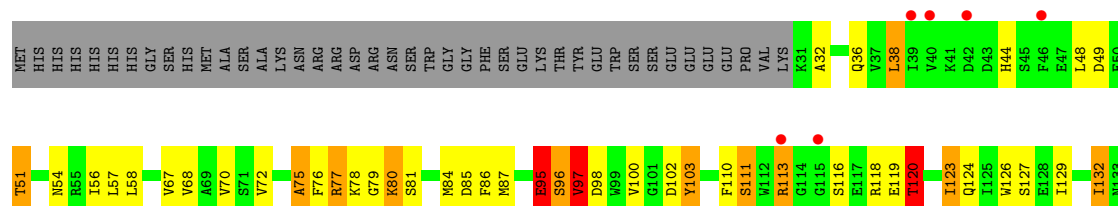


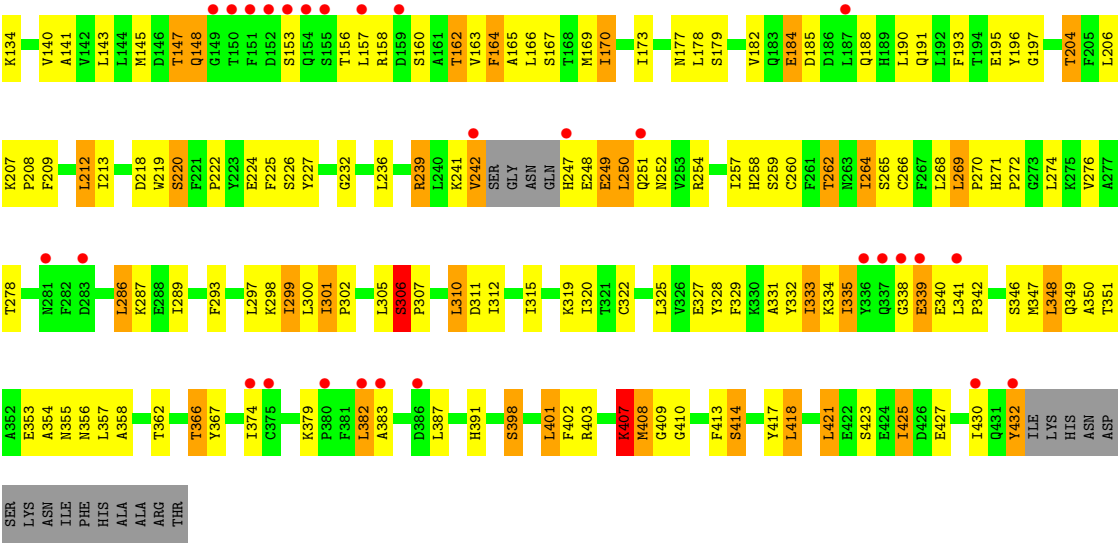


• Molecule 1: Atlastin-1



• Molecule 1: Atlastin-1





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	104.74Å 133.44Å 176.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.80 40.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.8 (40.00-2.80) 98.8 (40.00-2.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.25 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.225 , 0.267 0.217 , 0.258	Depositor DCC
R_{free} test set	3079 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	56.5	Xtriage
Anisotropy	0.688	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 50.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12574	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.73% 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: GDP, MG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.52	0/3173	0.97	11/4301 (0.3%)
1	B	0.52	0/3142	0.96	13/4252 (0.3%)
1	C	0.58	0/3180	0.99	13/4307 (0.3%)
1	D	0.49	0/3092	0.97	6/4195 (0.1%)
All	All	0.53	0/12587	0.98	43/17055 (0.3%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	D	0	1
All	All	0	2

There are no bond length outliers.

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	207	LYS	CA-C-N	12.00	131.72	119.24
1	A	207	LYS	C-N-CA	12.00	131.72	119.24
1	A	306	SER	CA-C-N	9.38	131.57	119.84
1	A	306	SER	C-N-CA	9.38	131.57	119.84
1	B	259	SER	N-CA-C	-9.15	101.83	113.72
1	D	306	SER	CA-C-N	8.04	129.89	119.84
1	D	306	SER	C-N-CA	8.04	129.89	119.84
1	D	259	SER	N-CA-C	-7.94	101.34	113.89
1	C	306	SER	CA-C-N	7.82	129.62	119.84
1	C	306	SER	C-N-CA	7.82	129.62	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	407	LYS	N-CA-C	7.38	123.93	114.56
1	C	259	SER	N-CA-C	-7.12	104.47	113.72
1	C	172	SER	N-CA-C	-6.37	106.73	114.75
1	C	372	GLU	N-CA-C	-6.20	106.08	113.21
1	C	353	GLU	N-CA-C	-5.97	104.68	111.07
1	B	338	GLY	N-CA-C	-5.90	105.84	112.33
1	D	407	LYS	N-CA-C	5.81	121.31	112.54
1	A	279	ASN	N-CA-C	5.78	119.61	108.45
1	B	271	HIS	CA-C-N	5.73	125.34	119.56
1	B	271	HIS	C-N-CA	5.73	125.34	119.56
1	C	413	PHE	N-CA-C	-5.72	101.57	110.10
1	D	379	LYS	CA-C-N	5.71	126.51	120.11
1	D	379	LYS	C-N-CA	5.71	126.51	120.11
1	A	67	VAL	N-CA-C	5.68	117.52	108.89
1	C	263	ASN	N-CA-C	5.62	116.59	107.32
1	B	279	ASN	CA-C-N	5.62	125.14	119.19
1	B	279	ASN	C-N-CA	5.62	125.14	119.19
1	B	306	SER	CA-C-N	5.57	126.81	119.84
1	B	306	SER	C-N-CA	5.57	126.81	119.84
1	C	253	VAL	N-CA-C	-5.57	104.94	110.62
1	C	97	VAL	N-CA-C	5.53	120.85	109.34
1	A	279	ASN	CA-C-N	5.50	124.69	118.97
1	A	279	ASN	C-N-CA	5.50	124.69	118.97
1	B	97	VAL	N-CA-C	5.40	120.56	109.34
1	B	102	ASP	N-CA-C	5.39	122.28	110.80
1	C	252	ASN	N-CA-C	-5.38	106.59	112.72
1	B	211	SER	N-CA-C	5.34	117.61	108.90
1	A	257	ILE	N-CA-C	-5.33	101.95	109.63
1	B	67	VAL	N-CA-C	5.26	116.06	108.48
1	A	366	THR	N-CA-C	-5.17	105.55	111.14
1	B	372	GLU	N-CA-C	-5.17	106.89	114.39
1	C	339	GLU	N-CA-C	-5.15	106.10	112.38
1	A	253	VAL	N-CA-C	-5.04	104.21	111.17

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	252	ASN	Peptide
1	D	95	GLU	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	3108	0	2980	156	0
1	B	3079	0	2997	189	0
1	C	3116	0	3007	171	0
1	D	3029	0	2899	156	0
2	A	28	0	12	0	0
2	B	28	0	12	2	0
2	C	28	0	12	3	0
2	D	28	0	12	4	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	37	0	0	2	0
4	B	29	0	0	1	0
4	C	37	0	0	3	0
4	D	23	0	0	1	0
All	All	12574	0	11931	667	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 27.

All (667) 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:196:TYR:HB2	1:B:347:MET:HE3	1.26	1.11
1:C:196:TYR:HB2	1:C:347:MET:HE3	1.39	1.01
1:C:84:MET:HG2	1:C:87:MET:HE2	1.43	0.99
1:B:403:ARG:HH21	1:B:418:LEU:HD11	1.21	0.98
1:B:402:PHE:O	1:B:407:LYS:HE3	1.65	0.95
1:A:165:ALA:HA	1:A:193:PHE:HE2	1.33	0.92
1:D:102:ASP:O	1:D:103:TYR:HB2	1.69	0.92
1:A:400:LYS:HZ3	1:A:403:ARG:HH11	1.16	0.92
1:A:165:ALA:HA	1:A:193:PHE:CE2	2.05	0.91
1:B:48:LEU:HB2	1:B:333:ILE:HG13	1.56	0.88
1:C:356:ASN:ND2	1:C:414:SER:HB2	1.88	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:250:LEU:H	1:D:250:LEU:HD23	1.39	0.87
1:D:48:LEU:HB2	1:D:333:ILE:HG13	1.56	0.87
1:A:400:LYS:NZ	1:A:403:ARG:HH11	1.73	0.86
1:C:305:LEU:O	1:C:306:SER:HB3	1.73	0.86
1:A:344:PRO:HA	1:A:347:MET:HE3	1.58	0.85
1:C:252:ASN:HB3	1:C:255:LYS:HE2	1.60	0.83
1:B:38:LEU:HB3	1:B:123:ILE:HG23	1.62	0.81
1:C:335:ILE:HD11	1:C:346:SER:O	1.80	0.81
1:D:191:GLN:HE22	1:D:250:LEU:HD12	1.45	0.81
1:A:200:ALA:O	1:A:204:THR:HG22	1.81	0.81
1:C:199:LEU:HD13	1:C:345:LYS:HE3	1.61	0.81
1:D:191:GLN:NE2	1:D:250:LEU:HD12	1.95	0.81
1:A:335:ILE:HD12	1:A:350:ALA:HB2	1.63	0.81
1:C:209:PHE:O	1:C:262:THR:HG22	1.82	0.80
1:D:95:GLU:O	1:D:96:SER:HB2	1.80	0.79
1:A:257:ILE:O	1:A:258:HIS:HB2	1.82	0.79
1:D:209:PHE:O	1:D:262:THR:HG22	1.81	0.79
1:D:407:LYS:HB2	1:D:407:LYS:NZ	1.97	0.79
1:B:102:ASP:HA	1:B:286:LEU:HD12	1.66	0.78
1:C:414:SER:HA	1:C:417:TYR:HD1	1.50	0.77
1:D:204:THR:HG22	1:D:206:LEU:H	1.48	0.76
1:C:103:TYR:HA	1:C:286:LEU:HB3	1.66	0.76
1:A:194:THR:O	1:A:198:ARG:HG3	1.86	0.76
1:C:402:PHE:O	1:C:407:LYS:HE3	1.85	0.75
1:B:335:ILE:HD11	1:B:346:SER:O	1.87	0.74
1:B:196:TYR:CB	1:B:347:MET:HE3	2.12	0.74
1:C:60:GLU:HA	1:C:63:ARG:HH21	1.53	0.74
1:C:343:HIS:ND1	1:C:344:PRO:HD2	2.02	0.74
1:A:387:LEU:HG	1:A:387:LEU:O	1.88	0.74
1:B:157:LEU:HD22	1:B:159:ASP:HB3	1.69	0.74
1:C:210:GLN:HA	1:C:262:THR:HG23	1.69	0.74
1:A:38:LEU:HB3	1:A:123:ILE:HG23	1.70	0.73
1:A:332:TYR:OH	1:A:351:THR:HB	1.87	0.73
1:A:264:ILE:HD13	1:A:264:ILE:O	1.87	0.73
1:B:305:LEU:O	1:B:306:SER:HB3	1.88	0.73
1:C:258:HIS:HB3	1:C:264:ILE:HG23	1.69	0.73
1:B:210:GLN:HA	1:B:262:THR:HG23	1.71	0.72
1:D:264:ILE:HD13	1:D:264:ILE:O	1.88	0.72
1:C:407:LYS:O	1:C:408:MET:HB2	1.87	0.72
1:A:204:THR:HG23	1:A:206:LEU:H	1.54	0.72
1:D:178:LEU:HD12	1:D:182:VAL:HG22	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:78:LYS:HD2	1:D:179:SER:HB2	1.71	0.71
1:C:358:ALA:O	1:C:362:THR:HG22	1.90	0.71
1:B:210:GLN:HA	1:B:262:THR:CG2	2.21	0.71
1:C:71:SER:HB2	1:C:145:MET:HE2	1.73	0.71
1:A:165:ALA:O	1:A:169:MET:HG2	1.91	0.71
1:B:364:LYS:HE3	1:B:428:LEU:HD11	1.72	0.70
1:C:72:VAL:HG12	1:C:80:LYS:HG2	1.71	0.70
1:C:103:TYR:HA	1:C:286:LEU:HD12	1.72	0.70
1:C:258:HIS:HB3	1:C:264:ILE:CG2	2.21	0.70
1:D:44:HIS:HB3	1:D:339:GLU:HB3	1.74	0.70
1:D:80:LYS:HG3	1:D:177:ASN:HD22	1.57	0.70
1:C:412:GLU:CD	1:C:412:GLU:H	2.00	0.70
1:D:347:MET:C	1:D:349:GLN:H	2.00	0.70
1:B:250:LEU:O	1:B:250:LEU:HG	1.92	0.69
1:C:312:ILE:HG21	1:C:319:LYS:HG3	1.73	0.69
1:A:337:GLN:HE21	1:A:337:GLN:HA	1.58	0.69
1:D:56:ILE:HD13	1:D:127:SER:HA	1.73	0.69
1:C:356:ASN:ND2	1:C:414:SER:CB	2.56	0.69
1:A:85:ASP:OD1	1:A:111:SER:HA	1.92	0.68
1:A:401:LEU:O	1:A:405:VAL:HG23	1.93	0.68
1:C:371:MET:C	1:C:373:GLU:H	2.02	0.68
1:B:196:TYR:HB2	1:B:347:MET:CE	2.15	0.68
1:D:332:TYR:O	1:D:335:ILE:HG22	1.93	0.68
1:D:402:PHE:O	1:D:407:LYS:HE3	1.94	0.68
1:A:110:PHE:HB3	1:A:126:TRP:HH2	1.59	0.67
1:D:327:GLU:OE1	1:D:357:LEU:HB3	1.94	0.67
1:D:335:ILE:HD12	1:D:350:ALA:HB2	1.75	0.67
1:B:104:ASN:O	1:B:105:GLU:HG3	1.95	0.67
1:C:151:PHE:HA	1:C:156:THR:HB	1.77	0.67
1:D:315:ILE:HD13	1:D:328:TYR:CE2	2.30	0.67
1:D:409:GLY:O	1:D:413:PHE:HB3	1.94	0.67
1:B:198:ARG:HD2	1:B:256:HIS:CG	2.30	0.66
1:B:84:MET:HG2	1:B:87:MET:HE2	1.77	0.66
1:C:301:ILE:HB	1:C:302:PRO:HD3	1.77	0.66
1:A:160:SER:C	1:A:162:THR:H	2.03	0.66
1:B:59:SER:O	1:B:63:ARG:HB2	1.95	0.66
1:D:299:ILE:O	1:D:299:ILE:HG13	1.91	0.66
1:B:395:LYS:HG3	1:B:425:ILE:HD11	1.78	0.66
1:D:38:LEU:HB3	1:D:123:ILE:HG23	1.76	0.66
1:B:286:LEU:HD22	1:B:294:ILE:HD11	1.77	0.66
1:D:356:ASN:HD21	1:D:414:SER:HB3	1.59	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:286:LEU:HD23	1:D:289:ILE:HD12	1.78	0.65
1:D:70:VAL:HG22	1:D:173:ILE:HD12	1.79	0.65
1:C:332:TYR:OH	1:C:351:THR:HB	1.96	0.65
1:B:236:LEU:HG	1:B:266:CYS:HB2	1.77	0.65
1:A:250:LEU:HD23	1:A:250:LEU:H	1.60	0.65
1:A:156:THR:HA	1:A:160:SER:HB2	1.79	0.65
1:D:156:THR:OG1	1:D:160:SER:HB2	1.97	0.65
1:C:196:TYR:CB	1:C:347:MET:HE3	2.23	0.65
1:C:210:GLN:HA	1:C:262:THR:CG2	2.27	0.64
1:B:107:LEU:HD22	1:B:107:LEU:O	1.97	0.64
1:A:403:ARG:HH21	1:A:418:LEU:HD11	1.61	0.64
1:D:103:TYR:HA	1:D:286:LEU:HB3	1.78	0.64
1:D:301:ILE:HB	1:D:302:PRO:HD3	1.79	0.64
1:B:165:ALA:HA	1:B:193:PHE:CE2	2.33	0.64
1:D:222:PRO:HG3	1:D:227:TYR:CZ	2.33	0.64
1:A:316:ASN:ND2	1:A:408:MET:HG3	2.13	0.64
1:C:272:PRO:HG3	2:C:593:GDP:C6	2.32	0.64
1:D:123:ILE:HD11	1:D:145:MET:HE3	1.78	0.64
1:B:136:ASP:OD1	1:B:138:LYS:HG2	1.97	0.63
1:B:315:ILE:HD13	1:B:328:TYR:CE2	2.33	0.63
1:C:204:THR:HG22	1:C:206:LEU:H	1.64	0.63
1:D:96:SER:O	1:D:97:VAL:O	2.16	0.63
1:B:85:ASP:OD1	1:B:111:SER:HA	1.98	0.63
1:B:296:ASN:N	1:B:296:ASN:HD22	1.97	0.63
1:C:386:ASP:O	1:C:389:THR:HG22	1.99	0.63
1:D:119:GLU:O	1:D:120:THR:C	2.41	0.63
1:A:85:ASP:CG	1:A:111:SER:HA	2.25	0.62
1:C:407:LYS:HB2	1:C:407:LYS:NZ	2.14	0.62
1:A:155:SER:O	1:A:159:ASP:HB3	1.98	0.62
1:A:407:LYS:O	1:A:408:MET:HB2	1.98	0.62
1:B:386:ASP:HA	1:B:389:THR:HG22	1.81	0.62
1:B:255:LYS:HE3	1:B:256:HIS:CE1	2.34	0.62
1:C:348:LEU:O	1:C:348:LEU:HD22	1.98	0.62
1:B:195:GLU:OE2	1:B:195:GLU:HA	1.98	0.62
1:A:257:ILE:O	1:A:258:HIS:CB	2.48	0.62
1:D:335:ILE:HD11	1:D:346:SER:O	2.00	0.61
1:D:356:ASN:ND2	1:D:414:SER:HB3	2.15	0.61
1:A:48:LEU:HB2	1:A:333:ILE:HG13	1.82	0.61
1:C:170:ILE:HD12	1:C:329:PHE:CZ	2.35	0.61
1:C:251:GLN:OE1	1:C:251:GLN:HA	2.00	0.61
1:D:169:MET:HE1	1:D:196:TYR:CE1	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:305:LEU:O	1:D:306:SER:HB3	1.99	0.61
1:A:103:TYR:HA	1:A:286:LEU:HB3	1.81	0.61
1:A:241:LYS:O	1:A:242:VAL:HB	1.99	0.61
1:D:407:LYS:HB2	1:D:407:LYS:HZ3	1.64	0.61
1:D:48:LEU:HD12	1:D:49:ASP:N	2.15	0.61
1:C:172:SER:HA	4:C:456:HOH:O	2.00	0.61
1:C:356:ASN:HD21	1:C:414:SER:CB	2.14	0.61
1:D:252:ASN:C	1:D:254:ARG:H	2.09	0.61
1:C:38:LEU:HB3	1:C:123:ILE:HG23	1.82	0.60
1:C:343:HIS:CE1	1:C:344:PRO:HD2	2.36	0.60
1:D:36:GLN:HE22	1:D:124:GLN:NE2	1.99	0.60
1:D:209:PHE:HB2	1:D:260:CYS:O	2.00	0.60
1:C:118:ARG:C	1:C:120:THR:H	2.10	0.60
1:C:143:LEU:HD23	4:C:452:HOH:O	2.01	0.60
1:C:199:LEU:CD1	1:C:345:LYS:HE3	2.30	0.60
1:B:147:THR:HG21	1:B:167:SER:CB	2.32	0.60
1:B:35:VAL:HG21	1:B:56:ILE:HD11	1.83	0.60
1:B:182:VAL:HB	1:B:239:ARG:HG3	1.82	0.60
1:B:299:ILE:HD12	1:B:299:ILE:O	2.00	0.60
1:A:110:PHE:HB3	1:A:126:TRP:CH2	2.37	0.59
1:A:162:THR:HG23	1:A:163:VAL:H	1.66	0.59
1:D:353:GLU:HG3	1:D:417:TYR:OH	2.02	0.59
1:D:358:ALA:O	1:D:362:THR:HG23	2.02	0.59
1:D:147:THR:HG21	1:D:167:SER:HB2	1.84	0.59
1:A:192:LEU:HD11	1:A:347:MET:HE1	1.85	0.59
1:A:252:ASN:N	1:A:252:ASN:HD22	2.00	0.59
1:A:423:SER:O	1:A:427:GLU:HG2	2.02	0.59
1:D:272:PRO:HG3	2:D:593:GDP:C6	2.38	0.59
1:C:63:ARG:HD2	1:C:323:ARG:NH2	2.18	0.58
1:C:256:HIS:O	1:C:259:SER:HB3	2.03	0.58
1:B:162:THR:HG22	1:B:163:VAL:N	2.18	0.58
1:A:80:LYS:O	1:A:84:MET:HG3	2.04	0.58
1:A:236:LEU:HD22	1:A:240:LEU:HG	1.86	0.58
1:D:102:ASP:O	1:D:103:TYR:CB	2.49	0.58
1:B:201:MET:HE2	1:B:259:SER:O	2.03	0.58
1:C:119:GLU:OE1	1:C:150:THR:HG22	2.04	0.58
1:C:345:LYS:HB2	1:C:345:LYS:NZ	2.18	0.58
1:C:305:LEU:O	1:C:306:SER:CB	2.46	0.58
1:B:430:ILE:HG13	1:B:430:ILE:O	2.04	0.58
1:C:299:ILE:O	1:C:299:ILE:HG13	1.99	0.58
1:B:32:ALA:HB3	1:B:110:PHE:CE2	2.39	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:103:TYR:CA	1:C:286:LEU:HB3	2.35	0.57
1:A:162:THR:HG23	1:A:163:VAL:N	2.20	0.57
1:C:151:PHE:HZ	1:C:164:PHE:HB2	1.68	0.57
1:B:165:ALA:HA	1:B:193:PHE:HE2	1.67	0.57
1:B:196:TYR:OH	1:B:351:THR:HG21	2.05	0.57
1:D:72:VAL:HG12	1:D:80:LYS:HG2	1.85	0.57
1:D:204:THR:CG2	1:D:206:LEU:H	2.17	0.57
1:C:228:GLY:HA2	1:C:267:PHE:CZ	2.40	0.57
1:A:335:ILE:CD1	1:A:350:ALA:HB2	2.33	0.56
1:A:348:LEU:HD22	1:A:408:MET:HE1	1.86	0.56
1:A:355:ASN:HD22	1:A:408:MET:HB2	1.69	0.56
1:D:239:ARG:HG2	1:D:239:ARG:HH11	1.69	0.56
1:A:160:SER:C	1:A:162:THR:N	2.63	0.56
1:A:160:SER:N	1:A:162:THR:HG22	2.20	0.56
1:C:356:ASN:HD22	1:C:414:SER:HB2	1.71	0.56
1:D:103:TYR:O	1:D:287:LYS:HG2	2.06	0.56
1:D:347:MET:C	1:D:349:GLN:N	2.60	0.56
1:D:382:LEU:HD21	1:D:432:TYR:HB2	1.86	0.56
1:C:147:THR:HG21	1:C:167:SER:CB	2.35	0.56
1:C:102:ASP:O	1:C:103:TYR:CB	2.54	0.56
1:D:49:ASP:OD2	1:D:51:THR:HG23	2.06	0.56
1:B:102:ASP:O	1:B:103:TYR:O	2.23	0.56
1:C:40:VAL:HB	1:C:121:THR:HG23	1.88	0.56
1:C:107:LEU:O	1:C:107:LEU:HD22	2.06	0.56
1:B:305:LEU:O	1:B:306:SER:CB	2.54	0.56
1:B:86:PHE:CG	1:B:297:LEU:HD21	2.40	0.56
1:B:37:VAL:HG12	1:B:38:LEU:N	2.21	0.56
1:B:190:LEU:HD13	1:B:257:ILE:CD1	2.36	0.55
1:D:355:ASN:HD22	1:D:408:MET:HB2	1.69	0.55
1:B:358:ALA:O	1:B:362:THR:HG22	2.06	0.55
1:C:190:LEU:HD13	1:C:257:ILE:HD11	1.88	0.55
1:C:407:LYS:HB2	1:C:407:LYS:HZ3	1.71	0.55
1:D:305:LEU:O	1:D:306:SER:CB	2.54	0.55
1:A:278:THR:CG2	1:A:278:THR:O	2.54	0.55
1:B:241:LYS:O	1:B:254:ARG:HD3	2.06	0.55
1:C:173:ILE:HD13	4:C:448:HOH:O	2.06	0.55
1:C:242:VAL:HG11	1:C:247:HIS:CB	2.36	0.55
1:D:68:VAL:HG13	1:D:311:ASP:O	2.07	0.55
1:D:148:GLN:CA	1:D:148:GLN:HE21	2.20	0.55
1:B:96:SER:O	1:B:97:VAL:O	2.23	0.55
1:D:421:LEU:O	1:D:425:ILE:HG12	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:228:GLY:HA2	1:B:267:PHE:CZ	2.42	0.55
1:B:31:LYS:O	1:B:32:ALA:HB3	2.07	0.55
1:C:86:PHE:CG	1:C:297:LEU:HD21	2.41	0.55
1:A:32:ALA:HA	1:A:127:SER:OG	2.07	0.55
1:A:196:TYR:OH	1:A:351:THR:HG21	2.06	0.55
1:A:301:ILE:HB	1:A:302:PRO:HD3	1.89	0.55
1:A:364:LYS:O	1:A:364:LYS:HG2	2.06	0.54
1:B:154:GLN:HG2	1:B:155:SER:OG	2.06	0.54
1:B:86:PHE:CD2	1:B:297:LEU:HD21	2.42	0.54
1:B:335:ILE:HD11	1:B:346:SER:C	2.31	0.54
1:C:158:ARG:O	1:C:159:ASP:CB	2.56	0.54
1:C:258:HIS:CD2	1:C:264:ILE:HD12	2.43	0.54
1:A:400:LYS:HZ3	1:A:403:ARG:HD2	1.72	0.54
1:B:54:ASN:HD21	1:B:326:VAL:HG11	1.72	0.54
1:C:421:LEU:O	1:C:425:ILE:HG12	2.08	0.54
1:A:318:ASN:O	1:A:320:ILE:HD12	2.07	0.54
1:A:358:ALA:O	1:A:362:THR:HG22	2.08	0.54
1:B:123:ILE:CD1	1:B:145:MET:HE3	2.38	0.54
1:B:232:GLY:O	1:B:266:CYS:HB3	2.08	0.54
1:A:364:LYS:HD2	1:A:424:GLU:OE1	2.08	0.54
1:B:162:THR:CG2	1:B:163:VAL:N	2.70	0.54
1:C:344:PRO:HA	1:C:347:MET:HE2	1.90	0.54
1:C:413:PHE:O	1:C:414:SER:CB	2.55	0.54
1:A:363:ALA:C	1:A:365:ASP:H	2.15	0.53
1:C:356:ASN:HD21	1:C:414:SER:HB2	1.67	0.53
1:D:148:GLN:HE21	1:D:148:GLN:HA	1.73	0.53
1:B:335:ILE:HG22	1:B:336:TYR:CD1	2.42	0.53
1:B:339:GLU:O	1:B:340:GLU:HB2	2.08	0.53
1:C:264:ILE:O	1:C:264:ILE:HD13	2.08	0.53
1:C:335:ILE:HG23	1:C:342:PRO:HB3	1.90	0.53
1:B:238:LYS:HB2	1:B:238:LYS:NZ	2.23	0.53
1:A:108:THR:O	1:A:108:THR:HG23	2.09	0.53
1:B:267:PHE:CG	1:B:300:LEU:HD23	2.44	0.53
1:D:170:ILE:HD13	1:D:329:PHE:CZ	2.43	0.53
1:D:75:ALA:HB3	1:D:78:LYS:HG3	1.91	0.53
1:D:156:THR:OG1	1:D:157:LEU:N	2.40	0.53
1:D:165:ALA:O	1:D:169:MET:HG2	2.08	0.53
1:D:76:PHE:CG	1:D:77:ARG:N	2.77	0.53
1:A:169:MET:HE2	1:A:332:TYR:CE1	2.44	0.53
1:A:278:THR:O	1:A:278:THR:HG22	2.09	0.53
1:C:222:PRO:HG3	1:C:227:TYR:CZ	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:228:GLY:HA2	1:A:267:PHE:CZ	2.43	0.53
1:B:362:THR:O	1:B:366:THR:HG22	2.09	0.53
1:A:201:MET:HE1	1:A:259:SER:HB3	1.91	0.53
1:D:209:PHE:O	1:D:262:THR:CG2	2.54	0.53
1:B:150:THR:C	1:B:152:ASP:H	2.17	0.52
1:C:79:GLY:HA2	2:C:593:GDP:O1A	2.09	0.52
1:C:250:LEU:C	1:C:252:ASN:H	2.16	0.52
1:B:209:PHE:O	1:B:262:THR:HG22	2.10	0.52
1:D:407:LYS:HB2	1:D:407:LYS:HZ2	1.72	0.52
1:B:154:GLN:HG2	1:B:155:SER:N	2.24	0.52
1:C:269:LEU:HD23	1:C:270:PRO:HD2	1.90	0.52
1:D:97:VAL:HG12	1:D:98:ASP:OD2	2.08	0.52
1:B:222:PRO:HG3	1:B:227:TYR:CZ	2.44	0.52
1:B:274:LEU:O	1:B:278:THR:HB	2.09	0.52
1:A:156:THR:O	1:A:157:LEU:C	2.52	0.52
1:C:208:PRO:HD2	1:C:260:CYS:O	2.09	0.52
1:D:212:LEU:HD12	1:D:264:ILE:HB	1.92	0.52
1:B:38:LEU:HB3	1:B:123:ILE:CG2	2.36	0.52
1:C:379:LYS:CB	1:C:380:PRO:HD3	2.39	0.52
1:D:252:ASN:C	1:D:254:ARG:N	2.68	0.52
1:A:269:LEU:CD2	1:A:270:PRO:HD2	2.39	0.52
1:C:190:LEU:HD13	1:C:257:ILE:CD1	2.40	0.52
1:D:178:LEU:CD1	1:D:182:VAL:HG22	2.38	0.52
1:A:280:PRO:HG3	1:B:223:TYR:CE1	2.45	0.52
1:A:400:LYS:HZ3	1:A:403:ARG:NH1	1.98	0.52
1:A:335:ILE:HD11	1:A:346:SER:O	2.10	0.51
1:B:166:LEU:O	1:B:170:ILE:HG23	2.10	0.51
1:D:362:THR:O	1:D:366:THR:HG22	2.10	0.51
1:B:157:LEU:HD22	1:B:159:ASP:CB	2.38	0.51
1:C:241:LYS:O	1:C:242:VAL:HB	2.10	0.51
1:C:371:MET:C	1:C:373:GLU:N	2.69	0.51
1:D:251:GLN:HA	1:D:251:GLN:OE1	2.09	0.51
1:B:94:GLN:HE22	1:B:132:ILE:HD12	1.74	0.51
1:A:400:LYS:NZ	1:A:403:ARG:NH1	2.51	0.51
1:B:356:ASN:ND2	1:B:407:LYS:HD2	2.26	0.51
1:C:165:ALA:O	1:C:166:LEU:C	2.54	0.51
1:D:156:THR:C	1:D:158:ARG:H	2.18	0.51
1:D:333:ILE:HG22	1:D:334:LYS:N	2.26	0.51
1:A:134:LYS:HB3	1:A:135:PRO:CD	2.40	0.51
1:A:316:ASN:HD22	1:A:408:MET:HG3	1.75	0.51
1:B:129:ILE:HG21	1:B:141:ALA:HB1	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:PRO:O	1:A:305:LEU:O	2.28	0.51
1:B:59:SER:OG	1:B:62:VAL:HG22	2.11	0.51
1:C:76:PHE:CE2	1:C:77:ARG:HG2	2.46	0.51
1:A:134:LYS:HB3	1:A:135:PRO:HD2	1.93	0.51
1:B:54:ASN:ND2	1:B:326:VAL:HG11	2.26	0.51
1:D:147:THR:HG21	1:D:167:SER:CB	2.40	0.51
1:D:156:THR:CB	1:D:160:SER:HB2	2.41	0.51
1:A:68:VAL:HG11	1:A:310:LEU:HG	1.92	0.50
1:B:91:MET:HB3	1:B:130:PHE:CE2	2.46	0.50
1:B:170:ILE:HD13	1:B:329:PHE:CE2	2.47	0.50
1:C:151:PHE:CZ	1:C:164:PHE:HB2	2.46	0.50
1:A:422:GLU:HA	1:A:425:ILE:HG13	1.92	0.50
1:B:399:VAL:O	1:B:403:ARG:HB2	2.12	0.50
1:A:223:TYR:CD1	1:B:280:PRO:HG3	2.46	0.50
1:B:258:HIS:HD1	1:B:258:HIS:C	2.20	0.50
1:D:407:LYS:O	1:D:408:MET:O	2.29	0.50
1:A:421:LEU:O	1:A:425:ILE:HG12	2.10	0.50
1:B:147:THR:HG21	1:B:167:SER:HB2	1.92	0.50
1:B:286:LEU:O	1:B:289:ILE:HB	2.11	0.50
1:C:204:THR:HG22	1:C:206:LEU:N	2.25	0.50
1:C:37:VAL:HG21	1:C:125:ILE:HD13	1.92	0.50
1:C:269:LEU:CD2	1:C:270:PRO:HD2	2.41	0.50
1:D:226:SER:O	1:D:268:LEU:HD12	2.11	0.50
1:D:272:PRO:HB2	1:D:276:VAL:HB	1.93	0.50
1:C:252:ASN:CB	1:C:255:LYS:HE2	2.37	0.50
1:A:88:LEU:HD11	1:A:112:TRP:NE1	2.26	0.49
1:A:191:GLN:HE22	1:A:250:LEU:HD12	1.76	0.49
1:D:190:LEU:HD13	1:D:257:ILE:HD11	1.93	0.49
1:A:273:GLY:HA2	1:B:274:LEU:HD11	1.94	0.49
1:B:101:GLY:O	1:B:102:ASP:O	2.30	0.49
1:C:165:ALA:HA	1:C:193:PHE:HZ	1.76	0.49
1:C:222:PRO:HG3	1:C:227:TYR:CE1	2.47	0.49
1:B:274:LEU:HG	4:B:448:HOH:O	2.13	0.49
1:C:212:LEU:HB2	1:C:261:PHE:CD2	2.47	0.49
1:D:262:THR:HA	4:D:452:HOH:O	2.13	0.49
1:A:368:ASN:HB2	1:A:428:LEU:HD21	1.93	0.49
1:B:151:PHE:O	1:B:152:ASP:C	2.55	0.49
1:D:110:PHE:HB3	1:D:126:TRP:CZ2	2.48	0.49
1:A:343:HIS:O	1:A:344:PRO:C	2.54	0.49
1:B:257:ILE:C	1:B:259:SER:H	2.21	0.49
1:D:145:MET:SD	1:D:170:ILE:HD11	2.52	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:123:ILE:HD11	1:B:145:MET:HE3	1.95	0.49
1:C:118:ARG:O	1:C:120:THR:HG22	2.13	0.49
1:A:165:ALA:HA	1:A:193:PHE:CZ	2.47	0.48
1:B:241:LYS:HG2	1:B:242:VAL:H	1.77	0.48
1:B:36:GLN:HE22	1:B:124:GLN:NE2	2.11	0.48
1:B:190:LEU:HD13	1:B:257:ILE:HD12	1.95	0.48
1:C:269:LEU:HD13	1:C:293:PHE:HE2	1.76	0.48
1:C:196:TYR:OH	1:C:351:THR:HG21	2.13	0.48
1:D:197:GLY:HA2	1:D:208:PRO:HG2	1.94	0.48
1:A:184:GLU:HG3	1:A:251:GLN:HG3	1.95	0.48
1:B:335:ILE:HG23	1:B:342:PRO:CG	2.43	0.48
1:C:403:ARG:HH21	1:C:418:LEU:HD11	1.79	0.48
1:D:213:ILE:CD1	1:D:265:SER:OG	2.61	0.48
1:C:335:ILE:HG22	1:C:336:TYR:CD1	2.48	0.48
1:B:358:ALA:O	1:B:361:ALA:HB3	2.14	0.48
1:D:250:LEU:H	1:D:250:LEU:CD2	2.15	0.48
1:A:299:ILE:O	1:A:299:ILE:HG13	2.09	0.48
1:B:267:PHE:CD1	1:B:300:LEU:HD23	2.49	0.48
1:A:395:LYS:HG3	1:A:425:ILE:HD12	1.94	0.48
1:A:404:GLY:O	1:A:405:VAL:C	2.55	0.48
1:C:333:ILE:HD13	1:C:333:ILE:O	2.14	0.48
1:D:156:THR:HG23	1:D:157:LEU:N	2.28	0.48
1:D:347:MET:O	1:D:351:THR:HG22	2.14	0.48
1:A:337:GLN:HE21	1:A:337:GLN:CA	2.24	0.48
1:B:272:PRO:HG3	2:B:593:GDP:C6	2.49	0.48
1:C:347:MET:O	1:C:351:THR:HG22	2.13	0.48
1:C:363:ALA:HA	1:C:366:THR:CG2	2.43	0.48
1:D:67:VAL:HG11	1:D:325:LEU:HD22	1.95	0.48
1:D:271:HIS:ND1	1:D:272:PRO:HD2	2.29	0.48
1:B:332:TYR:OH	1:B:351:THR:HB	2.12	0.47
1:C:136:ASP:OD1	1:C:138:LYS:HG2	2.14	0.47
1:C:371:MET:HE2	1:C:371:MET:HB3	1.74	0.47
1:A:221:PHE:N	1:A:222:PRO:HD3	2.29	0.47
1:A:241:LYS:O	1:A:254:ARG:HD3	2.13	0.47
1:A:252:ASN:N	1:A:252:ASN:ND2	2.62	0.47
1:B:183:GLN:O	1:B:186:ASP:HB2	2.14	0.47
1:C:110:PHE:HB3	1:C:126:TRP:HH2	1.79	0.47
1:A:146:ASP:C	1:A:146:ASP:OD1	2.56	0.47
1:B:425:ILE:HA	1:B:428:LEU:HB2	1.97	0.47
1:C:302:PRO:O	1:C:306:SER:HB3	2.13	0.47
1:A:335:ILE:HG22	1:A:336:TYR:CD1	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:LEU:HD13	1:B:107:LEU:H	1.80	0.47
1:D:169:MET:HE1	1:D:196:TYR:CZ	2.49	0.47
1:A:269:LEU:HD23	1:A:270:PRO:HD2	1.97	0.47
1:B:147:THR:HG21	1:B:167:SER:HB3	1.95	0.47
1:B:379:LYS:CB	1:B:380:PRO:HD2	2.45	0.47
1:D:38:LEU:HD13	1:D:333:ILE:HD11	1.95	0.47
1:A:169:MET:HE3	1:A:169:MET:HB3	1.78	0.47
1:C:248:GLU:O	1:C:249:GLU:O	2.33	0.47
1:D:129:ILE:HG21	1:D:141:ALA:HB1	1.96	0.47
1:D:218:ASP:OD2	2:D:593:GDP:N2	2.41	0.47
1:D:355:ASN:ND2	1:D:408:MET:HB2	2.28	0.47
1:A:208:PRO:HB2	1:A:209:PHE:CD1	2.50	0.47
1:B:31:LYS:O	1:B:32:ALA:CB	2.63	0.47
1:B:60:GLU:HA	1:B:63:ARG:HH21	1.79	0.47
1:D:403:ARG:HH21	1:D:418:LEU:HD11	1.80	0.47
1:A:38:LEU:HB3	1:A:123:ILE:CG2	2.42	0.46
1:B:153:SER:O	1:B:154:GLN:C	2.57	0.46
1:C:166:LEU:HD22	1:C:336:TYR:CD1	2.50	0.46
1:D:398:SER:HA	1:D:401:LEU:HB2	1.97	0.46
1:A:250:LEU:C	1:A:252:ASN:H	2.23	0.46
1:C:434:LYS:O	1:C:436:ASN:N	2.49	0.46
1:D:224:GLU:HG2	1:D:225:PHE:CD1	2.51	0.46
1:A:35:VAL:O	1:A:37:VAL:HG23	2.15	0.46
1:B:316:ASN:ND2	1:B:408:MET:HG3	2.30	0.46
1:B:394:LEU:HA	1:B:397:GLU:CD	2.41	0.46
1:B:366:THR:C	1:B:368:ASN:H	2.22	0.46
1:C:194:THR:HA	1:C:260:CYS:SG	2.56	0.46
1:A:48:LEU:HD11	1:A:53:LEU:HD22	1.96	0.46
1:B:407:LYS:O	1:B:408:MET:O	2.33	0.46
1:A:400:LYS:HD3	1:A:400:LYS:HA	1.60	0.46
1:C:343:HIS:O	1:C:344:PRO:C	2.57	0.46
1:C:382:LEU:HA	1:C:436:ASN:OD1	2.15	0.46
1:D:196:TYR:HE2	1:D:348:LEU:HA	1.81	0.46
1:A:180:GLN:O	1:A:181:ASN:HB2	2.15	0.46
1:A:381:PHE:O	1:A:382:LEU:HB2	2.15	0.46
1:D:196:TYR:OH	1:D:351:THR:HG21	2.14	0.46
1:D:220:SER:C	1:D:222:PRO:HD3	2.40	0.46
1:D:232:GLY:O	1:D:266:CYS:HB3	2.16	0.46
1:C:42:ASP:C	1:C:44:HIS:H	2.23	0.46
1:D:341:LEU:HA	1:D:342:PRO:HD3	1.76	0.46
1:C:118:ARG:C	1:C:120:THR:N	2.74	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:285:LYS:O	1:C:288:GLU:HB2	2.16	0.46
1:D:423:SER:O	1:D:427:GLU:HG3	2.15	0.46
1:A:161:ALA:O	1:A:165:ALA:HB2	2.16	0.46
1:B:89:ARG:NH2	1:B:107:LEU:HB2	2.30	0.46
1:B:208:PRO:HD2	1:B:260:CYS:O	2.16	0.46
1:B:316:ASN:C	1:B:318:ASN:H	2.24	0.46
1:C:305:LEU:N	1:C:305:LEU:HD23	2.31	0.46
1:D:118:ARG:O	1:D:120:THR:HG22	2.16	0.46
1:A:408:MET:HE2	1:A:408:MET:HB3	1.70	0.45
1:B:123:ILE:HD12	1:B:147:THR:HG22	1.98	0.45
1:C:107:LEU:HD22	1:C:107:LEU:C	2.41	0.45
1:A:247:HIS:N	4:A:483:HOH:O	2.49	0.45
1:B:199:LEU:HD13	1:B:345:LYS:HD3	1.98	0.45
1:C:72:VAL:CG1	1:C:80:LYS:HG2	2.44	0.45
1:D:241:LYS:O	1:D:242:VAL:HB	2.16	0.45
1:A:222:PRO:HG3	1:A:227:TYR:CZ	2.51	0.45
1:B:32:ALA:HA	1:B:127:SER:OG	2.16	0.45
1:B:339:GLU:O	1:B:340:GLU:CB	2.63	0.45
1:C:414:SER:HA	1:C:417:TYR:CD1	2.39	0.45
1:A:106:PRO:O	1:A:107:LEU:C	2.59	0.45
1:A:382:LEU:HD23	1:A:384:PRO:HD3	1.98	0.45
1:B:407:LYS:NZ	1:B:407:LYS:HB2	2.32	0.45
1:D:249:GLU:CB	1:D:250:LEU:HD23	2.47	0.45
1:A:67:VAL:HG22	1:A:68:VAL:N	2.32	0.45
1:B:132:ILE:HG13	1:B:133:ASN:N	2.30	0.45
1:D:298:LYS:O	1:D:302:PRO:HD2	2.17	0.45
1:A:376:GLY:O	1:A:377:GLY:C	2.59	0.45
1:B:257:ILE:O	1:B:258:HIS:CG	2.69	0.45
1:D:269:LEU:HD13	1:D:293:PHE:CE2	2.52	0.45
1:B:56:ILE:HD12	1:B:125:ILE:HD11	1.99	0.45
1:C:110:PHE:HB3	1:C:126:TRP:CH2	2.52	0.45
1:A:207:LYS:HA	1:A:208:PRO:HD3	1.71	0.45
1:B:134:LYS:HB3	1:B:134:LYS:HE2	1.57	0.45
1:B:347:MET:O	1:B:351:THR:HG22	2.15	0.45
1:D:132:ILE:CD1	1:D:307:PRO:HG3	2.47	0.45
1:B:272:PRO:CG	2:B:593:GDP:C6	2.99	0.45
1:C:435:HIS:CD2	1:C:435:HIS:C	2.95	0.45
1:D:156:THR:HG23	1:D:157:LEU:H	1.81	0.45
1:D:432:TYR:N	1:D:432:TYR:CD1	2.85	0.45
1:A:119:GLU:O	1:A:121:THR:N	2.50	0.45
1:A:220:SER:C	1:A:222:PRO:HD3	2.42	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:280:PRO:HD3	1:B:223:TYR:CE2	2.52	0.45
1:B:198:ARG:HD2	1:B:256:HIS:CD2	2.52	0.45
1:B:238:LYS:HB2	1:B:238:LYS:HZ3	1.80	0.45
1:C:56:ILE:HD12	1:C:125:ILE:HD11	1.98	0.45
1:C:134:LYS:HB3	1:C:134:LYS:HE2	1.78	0.45
1:C:162:THR:CG2	1:C:163:VAL:N	2.80	0.45
1:C:272:PRO:HB2	1:C:276:VAL:HB	1.98	0.45
1:A:207:LYS:HG2	1:A:259:SER:O	2.17	0.44
1:A:49:ASP:OD2	1:A:49:ASP:C	2.60	0.44
1:B:33:GLY:O	1:B:127:SER:HB3	2.17	0.44
1:C:39:ILE:HD12	1:C:39:ILE:HA	1.62	0.44
1:D:81:SER:HB2	2:D:593:GDP:O2A	2.18	0.44
1:D:232:GLY:O	1:D:236:LEU:HB2	2.18	0.44
1:D:239:ARG:HG2	1:D:239:ARG:NH1	2.31	0.44
1:D:242:VAL:HG11	1:D:247:HIS:N	2.33	0.44
1:A:356:ASN:HD21	1:A:414:SER:HB3	1.82	0.44
1:B:95:GLU:O	1:B:96:SER:O	2.35	0.44
1:C:343:HIS:CG	1:C:344:PRO:HD2	2.52	0.44
1:A:122:GLY:C	1:A:123:ILE:HG22	2.43	0.44
1:B:107:LEU:H	1:B:107:LEU:CD1	2.30	0.44
1:B:155:SER:O	1:B:156:THR:CB	2.66	0.44
1:B:301:ILE:HB	1:B:302:PRO:HD3	2.00	0.44
1:B:399:VAL:HG13	1:B:418:LEU:HD21	2.00	0.44
1:C:212:LEU:HB2	1:C:261:PHE:CE2	2.53	0.44
1:C:129:ILE:HG21	1:C:141:ALA:HB1	1.98	0.44
1:D:178:LEU:HD13	1:D:182:VAL:HA	1.99	0.44
1:A:123:ILE:HD11	1:A:145:MET:HE3	2.00	0.44
1:C:60:GLU:CA	1:C:63:ARG:HH21	2.27	0.44
1:C:162:THR:HG22	1:C:163:VAL:N	2.33	0.44
1:C:382:LEU:O	1:C:383:ALA:HB3	2.17	0.44
1:D:76:PHE:CE2	1:D:77:ARG:HG3	2.53	0.44
1:D:134:LYS:HE3	1:D:310:LEU:O	2.17	0.44
1:A:422:GLU:HA	1:A:425:ILE:HD11	2.00	0.44
1:B:279:ASN:HA	1:B:280:PRO:HD3	1.74	0.44
1:B:63:ARG:HD2	1:B:323:ARG:HH21	1.83	0.44
1:B:267:PHE:HE1	1:B:299:ILE:HD11	1.82	0.44
1:B:296:ASN:N	1:B:296:ASN:ND2	2.61	0.44
1:C:269:LEU:HD13	1:C:293:PHE:CE2	2.52	0.44
1:A:80:LYS:HB2	1:A:80:LYS:NZ	2.33	0.44
1:A:154:GLN:C	1:A:156:THR:H	2.26	0.44
1:A:343:HIS:C	1:A:345:LYS:N	2.74	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:ASP:CG	1:B:111:SER:HA	2.43	0.44
1:C:410:GLY:O	1:C:413:PHE:O	2.35	0.44
1:D:184:GLU:HG3	1:D:251:GLN:HG3	2.00	0.43
1:A:255:LYS:HG3	1:A:256:HIS:CD2	2.53	0.43
1:A:313:LYS:HE3	1:A:315:ILE:HD11	1.99	0.43
1:B:190:LEU:HD13	1:B:257:ILE:HD11	2.00	0.43
1:B:380:PRO:HB2	1:B:381:PHE:H	1.50	0.43
1:C:184:GLU:O	1:C:188:GLN:HB2	2.18	0.43
1:C:363:ALA:HA	1:C:366:THR:HG22	2.00	0.43
1:B:367:TYR:O	1:B:367:TYR:CG	2.71	0.43
1:A:279:ASN:HA	1:A:280:PRO:HD3	1.75	0.43
1:A:348:LEU:CD2	1:A:408:MET:HE1	2.48	0.43
1:A:356:ASN:ND2	1:A:414:SER:HB3	2.34	0.43
1:A:402:PHE:O	1:A:407:LYS:HE3	2.18	0.43
1:C:391:HIS:CE1	1:C:429:TYR:HB2	2.53	0.43
1:D:80:LYS:HB2	1:D:80:LYS:NZ	2.34	0.43
1:D:79:GLY:CA	2:D:593:GDP:O1A	2.67	0.43
1:D:86:PHE:CG	1:D:297:LEU:HD21	2.53	0.43
1:D:258:HIS:HB3	1:D:264:ILE:CG2	2.48	0.43
1:B:176:TYR:CZ	1:B:178:LEU:HD21	2.53	0.43
1:B:302:PRO:O	1:B:305:LEU:O	2.37	0.43
1:C:60:GLU:HA	1:C:63:ARG:NH2	2.28	0.43
1:C:209:PHE:O	1:C:262:THR:CG2	2.59	0.43
1:C:258:HIS:CD2	1:C:264:ILE:CD1	3.02	0.43
1:C:319:LYS:HA	1:C:319:LYS:HD2	1.86	0.43
1:B:134:LYS:HE3	1:B:307:PRO:O	2.19	0.43
1:C:428:LEU:HD23	1:C:428:LEU:HA	1.81	0.43
1:B:402:PHE:CD2	1:B:402:PHE:C	2.97	0.43
1:D:271:HIS:CG	1:D:272:PRO:HD2	2.54	0.43
1:A:35:VAL:O	1:A:36:GLN:C	2.62	0.43
1:A:335:ILE:HG23	1:A:342:PRO:CG	2.49	0.43
1:B:162:THR:HG22	1:B:163:VAL:H	1.80	0.43
1:B:241:LYS:HG2	1:B:242:VAL:N	2.33	0.43
1:C:153:SER:C	1:C:155:SER:H	2.26	0.43
1:D:54:ASN:C	1:D:56:ILE:N	2.77	0.43
1:B:360:VAL:HG23	1:B:402:PHE:CE1	2.54	0.43
1:B:388:GLN:HA	1:B:429:TYR:OH	2.18	0.43
1:C:402:PHE:CE2	1:C:407:LYS:HE2	2.54	0.43
1:C:430:ILE:C	1:C:432:TYR:H	2.27	0.43
1:C:314:GLU:HG3	1:C:319:LYS:HD2	2.01	0.42
1:C:338:GLY:C	1:C:340:GLU:N	2.76	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:370:LYS:HD3	1:C:394:LEU:CD2	2.49	0.42
1:C:434:LYS:O	1:C:435:HIS:C	2.61	0.42
1:D:57:LEU:HD23	1:D:57:LEU:HA	1.81	0.42
1:D:227:TYR:CG	1:D:270:PRO:HG3	2.53	0.42
1:B:316:ASN:HD22	1:B:408:MET:HG3	1.84	0.42
1:C:145:MET:HE1	1:C:170:ILE:HD11	2.02	0.42
1:C:413:PHE:O	1:C:414:SER:OG	2.36	0.42
1:D:207:LYS:HA	1:D:208:PRO:HD3	1.81	0.42
1:C:136:ASP:OD1	1:C:137:GLY:N	2.52	0.42
1:D:248:GLU:O	1:D:249:GLU:O	2.38	0.42
1:A:343:HIS:O	1:A:345:LYS:N	2.53	0.42
1:B:38:LEU:CD1	1:B:333:ILE:HD11	2.49	0.42
1:B:39:ILE:HG13	1:B:40:VAL:N	2.32	0.42
1:B:160:SER:C	1:B:162:THR:N	2.77	0.42
1:C:73:ALA:C	1:C:80:LYS:HD3	2.44	0.42
1:C:196:TYR:CE2	1:C:351:THR:HG21	2.55	0.42
1:C:362:THR:O	1:C:366:THR:HG22	2.20	0.42
1:D:84:MET:HG2	1:D:87:MET:HE2	2.01	0.42
1:D:85:ASP:CG	1:D:111:SER:HA	2.44	0.42
1:D:193:PHE:O	1:D:195:GLU:N	2.52	0.42
1:A:220:SER:O	1:A:221:PHE:CD2	2.72	0.42
1:A:419:GLN:O	1:A:420:GLN:C	2.62	0.42
1:B:348:LEU:HD23	1:B:348:LEU:HA	1.85	0.42
1:D:421:LEU:HD22	1:D:425:ILE:CD1	2.49	0.42
1:A:129:ILE:HG22	1:A:130:PHE:N	2.35	0.42
1:C:59:SER:O	1:C:63:ARG:HB2	2.20	0.42
1:C:271:HIS:ND1	1:C:272:PRO:HD2	2.35	0.42
1:C:359:ALA:O	1:C:362:THR:HG23	2.20	0.42
1:A:374:ILE:HD11	1:A:394:LEU:HD11	2.02	0.42
1:B:120:THR:HG23	1:B:120:THR:O	2.19	0.42
1:B:258:HIS:C	1:B:258:HIS:ND1	2.78	0.42
1:C:74:GLY:O	1:C:80:LYS:HE2	2.19	0.42
1:C:200:ALA:O	1:C:204:THR:HB	2.19	0.42
1:A:35:VAL:HG21	1:A:56:ILE:HD11	2.01	0.42
1:B:212:LEU:HB3	1:B:264:ILE:HB	2.01	0.42
1:A:422:GLU:HA	1:A:425:ILE:CG1	2.49	0.42
1:B:68:VAL:HG11	1:B:310:LEU:HG	2.00	0.42
1:B:228:GLY:HA2	1:B:267:PHE:CE2	2.55	0.42
1:A:269:LEU:HD22	1:A:270:PRO:HD2	2.02	0.42
1:C:132:ILE:HD11	1:C:307:PRO:HA	2.01	0.42
1:C:228:GLY:HA2	1:C:267:PHE:CE2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:252:ASN:N	1:C:252:ASN:ND2	2.65	0.42
1:C:257:ILE:C	1:C:259:SER:H	2.28	0.42
1:D:188:GLN:O	1:D:191:GLN:HB3	2.19	0.42
1:B:94:GLN:NE2	1:B:307:PRO:HD3	2.35	0.41
1:B:157:LEU:C	1:B:159:ASP:N	2.77	0.41
1:B:252:ASN:C	1:B:254:ARG:N	2.76	0.41
1:B:401:LEU:HD22	1:B:401:LEU:HA	1.82	0.41
1:B:110:PHE:HB3	1:B:126:TRP:CH2	2.55	0.41
1:B:181:ASN:OD1	1:B:235:PHE:HZ	2.04	0.41
1:D:113:ARG:HD2	1:D:113:ARG:HA	1.91	0.41
1:D:129:ILE:CG2	1:D:141:ALA:HB1	2.50	0.41
1:B:333:ILE:HG22	1:B:334:LYS:N	2.35	0.41
1:B:134:LYS:HB3	1:B:135:PRO:CD	2.50	0.41
1:B:158:ARG:O	1:B:162:THR:HB	2.21	0.41
1:B:238:LYS:NZ	1:B:238:LYS:CB	2.83	0.41
1:B:239:ARG:HD3	1:B:239:ARG:HA	1.85	0.41
1:B:264:ILE:HD13	1:B:264:ILE:O	2.21	0.41
1:B:286:LEU:HD22	1:B:294:ILE:CD1	2.47	0.41
1:C:326:VAL:CG1	1:C:330:LYS:HE3	2.50	0.41
1:D:57:LEU:HD22	1:D:322:CYS:HB3	2.02	0.41
1:D:271:HIS:HA	1:D:272:PRO:HD3	1.87	0.41
1:D:421:LEU:HD22	1:D:425:ILE:HD13	2.02	0.41
1:A:305:LEU:HA	1:A:305:LEU:HD23	1.73	0.41
1:A:370:LYS:HB3	1:A:394:LEU:HD13	2.03	0.41
1:B:76:PHE:HD1	1:B:148:GLN:OE1	2.03	0.41
1:B:84:MET:SD	1:B:146:ASP:HB2	2.60	0.41
1:B:393:GLN:O	1:B:397:GLU:HG3	2.20	0.41
1:A:123:ILE:O	1:A:123:ILE:HG12	2.20	0.41
1:A:170:ILE:HD13	1:A:329:PHE:CZ	2.55	0.41
1:A:221:PHE:HA	1:A:223:TYR:CE2	2.56	0.41
1:A:228:GLY:HA2	1:A:267:PHE:CE1	2.56	0.41
1:A:242:VAL:HG22	1:C:401:LEU:HD22	2.02	0.41
1:A:422:GLU:HA	1:A:425:ILE:CD1	2.50	0.41
1:B:160:SER:O	1:B:161:ALA:C	2.63	0.41
1:D:312:ILE:HG21	1:D:319:LYS:CG	2.51	0.41
1:A:167:SER:O	1:A:171:SER:HB3	2.21	0.41
1:A:210:GLN:HB3	1:A:211:SER:H	1.53	0.41
1:B:150:THR:C	1:B:152:ASP:N	2.76	0.41
1:B:196:TYR:O	1:B:197:GLY:C	2.62	0.41
1:B:366:THR:C	1:B:368:ASN:N	2.78	0.41
1:C:276:VAL:CG2	1:C:289:ILE:HG12	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:206:LEU:HD12	1:A:316:ASN:HA	2.02	0.41
1:A:313:LYS:HG3	1:A:320:ILE:HD13	2.03	0.41
1:A:402:PHE:CE2	1:A:407:LYS:HE2	2.56	0.41
1:B:374:ILE:HG22	1:B:374:ILE:O	2.21	0.41
1:D:338:GLY:O	1:D:340:GLU:N	2.54	0.41
1:A:58:LEU:O	1:A:59:SER:C	2.64	0.41
1:B:193:PHE:O	1:B:197:GLY:N	2.43	0.41
1:C:258:HIS:HD2	1:C:264:ILE:CD1	2.34	0.41
1:D:56:ILE:CD1	1:D:127:SER:HA	2.46	0.41
1:D:80:LYS:HG3	1:D:177:ASN:ND2	2.31	0.41
1:D:219:TRP:HD1	1:D:227:TYR:CE2	2.39	0.41
1:D:332:TYR:O	1:D:333:ILE:C	2.62	0.41
1:D:333:ILE:CG2	1:D:334:LYS:N	2.81	0.41
1:D:367:TYR:CD2	1:D:367:TYR:C	2.99	0.41
1:A:151:PHE:N	1:A:160:SER:OG	2.44	0.41
1:B:221:PHE:N	1:B:222:PRO:HD3	2.36	0.41
1:D:331:ALA:HB3	1:D:354:ALA:HB2	2.03	0.41
1:A:84:MET:SD	1:A:146:ASP:HB2	2.61	0.40
1:A:371:MET:HE2	1:A:391:HIS:HA	2.04	0.40
1:A:386:ASP:C	1:A:388:GLN:H	2.29	0.40
1:B:196:TYR:CD1	1:B:196:TYR:C	2.98	0.40
1:B:220:SER:C	1:B:222:PRO:HD3	2.46	0.40
1:C:58:LEU:HD12	1:C:58:LEU:HA	1.85	0.40
1:C:80:LYS:HB2	1:C:80:LYS:NZ	2.36	0.40
1:C:165:ALA:HA	1:C:193:PHE:CZ	2.55	0.40
1:D:425:ILE:HG12	1:D:425:ILE:H	1.67	0.40
1:A:126:TRP:HB3	4:A:452:HOH:O	2.19	0.40
1:B:84:MET:HA	1:B:87:MET:HE2	2.04	0.40
1:B:360:VAL:HG23	1:B:402:PHE:HE1	1.86	0.40
1:D:315:ILE:HB	1:D:320:ILE:HD11	2.02	0.40
1:C:79:GLY:CA	2:C:593:GDP:O1A	2.69	0.40
1:C:290:ASP:OD2	1:C:292:GLU:N	2.54	0.40
1:D:274:LEU:HA	1:D:274:LEU:HD23	1.82	0.40
1:B:59:SER:O	1:B:63:ARG:CB	2.68	0.40
1:B:212:LEU:HD23	1:B:212:LEU:HA	1.91	0.40
1:C:184:GLU:HG3	1:C:251:GLN:HG3	2.03	0.40
1:C:335:ILE:HD13	1:C:335:ILE:HG21	1.81	0.40
1:A:364:LYS:O	1:A:364:LYS:CG	2.70	0.40
1:B:103:TYR:O	1:B:105:GLU:N	2.55	0.40
1:D:160:SER:C	1:D:162:THR:N	2.79	0.40
1:D:163:VAL:O	1:D:164:PHE:C	2.65	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	400/459 (87%)	323 (81%)	57 (14%)	20 (5%)	1	5
1	B	393/459 (86%)	333 (85%)	37 (9%)	23 (6%)	1	3
1	C	401/459 (87%)	337 (84%)	43 (11%)	21 (5%)	1	5
1	D	394/459 (86%)	312 (79%)	62 (16%)	20 (5%)	1	5
All	All	1588/1836 (86%)	1305 (82%)	199 (12%)	84 (5%)	1	5

All (84) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	123	ILE
1	A	249	GLU
1	A	258	HIS
1	A	307	PRO
1	A	382	LEU
1	B	96	SER
1	B	97	VAL
1	B	102	ASP
1	B	103	TYR
1	B	104	ASN
1	B	111	SER
1	B	154	GLN
1	B	249	GLU
1	B	306	SER
1	B	380	PRO
1	B	381	PHE
1	B	408	MET
1	C	96	SER
1	C	97	VAL
1	C	249	GLU

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Mol	Chain	Res	Type
1	C	382	LEU
1	D	96	SER
1	D	97	VAL
1	D	120	THR
1	D	249	GLU
1	D	374	ILE
1	D	408	MET
1	A	97	VAL
1	A	120	THR
1	A	153	SER
1	A	284	GLY
1	A	377	GLY
1	A	393	GLN
1	B	156	THR
1	B	241	LYS
1	C	102	ASP
1	C	103	TYR
1	C	159	ASP
1	C	377	GLY
1	D	75	ALA
1	D	103	TYR
1	D	111	SER
1	D	153	SER
1	D	410	GLY
1	A	127	SER
1	A	157	LEU
1	A	159	ASP
1	A	375	CYS
1	A	380	PRO
1	B	32	ALA
1	B	158	ARG
1	C	111	SER
1	C	115	GLY
1	C	119	GLU
1	C	338	GLY
1	C	379	LYS
1	C	434	LYS
1	C	435	HIS
1	D	164	PHE
1	A	36	GLN
1	B	152	ASP
1	B	155	SER

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Mol	Chain	Res	Type
1	B	340	GLU
1	C	158	ARG
1	C	306	SER
1	D	95	GLU
1	D	306	SER
1	D	339	GLU
1	A	32	ALA
1	A	364	LYS
1	B	180	GLN
1	C	43	ASP
1	C	383	ALA
1	D	32	ALA
1	D	113	ARG
1	D	116	SER
1	A	306	SER
1	B	95	GLU
1	B	375	CYS
1	B	430	ILE
1	D	430	ILE
1	C	380	PRO
1	D	383	ALA
1	C	374	ILE

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	320/405 (79%)	262 (82%)	58 (18%)	2	6
1	B	323/405 (80%)	258 (80%)	65 (20%)	1	4
1	C	321/405 (79%)	262 (82%)	59 (18%)	1	6
1	D	306/405 (76%)	256 (84%)	50 (16%)	2	8
All	All	1270/1620 (78%)	1038 (82%)	232 (18%)	2	6

All (232) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	LYS
1	A	100	VAL
1	A	120	THR
1	A	121	THR
1	A	123	ILE
1	A	124	GLN
1	A	125	ILE
1	A	132	ILE
1	A	143	LEU
1	A	145	MET
1	A	147	THR
1	A	148	GLN
1	A	159	ASP
1	A	160	SER
1	A	166	LEU
1	A	170	ILE
1	A	184	GLU
1	A	203	GLU
1	A	210	GLN
1	A	212	LEU
1	A	213	ILE
1	A	236	LEU
1	A	238	LYS
1	A	239	ARG
1	A	252	ASN
1	A	253	VAL
1	A	255	LYS
1	A	257	ILE
1	A	263	ASN
1	A	264	ILE
1	A	269	LEU
1	A	278	THR
1	A	286	LEU
1	A	299	ILE
1	A	300	LEU
1	A	306	SER
1	A	309	SER
1	A	310	LEU
1	A	312	ILE
1	A	319	LYS
1	A	333	ILE
1	A	335	ILE
1	A	337	GLN

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Mol	Chain	Res	Type
1	A	348	LEU
1	A	349	GLN
1	A	351	THR
1	A	362	THR
1	A	366	THR
1	A	382	LEU
1	A	387	LEU
1	A	396	GLU
1	A	398	SER
1	A	399	VAL
1	A	407	LYS
1	A	412	GLU
1	A	418	LEU
1	A	421	LEU
1	A	425	ILE
1	B	37	VAL
1	B	39	ILE
1	B	42	ASP
1	B	51	THR
1	B	67	VAL
1	B	80	LYS
1	B	97	VAL
1	B	98	ASP
1	B	100	VAL
1	B	102	ASP
1	B	107	LEU
1	B	121	THR
1	B	123	ILE
1	B	125	ILE
1	B	132	ILE
1	B	134	LYS
1	B	138	LYS
1	B	140	VAL
1	B	143	LEU
1	B	145	MET
1	B	146	ASP
1	B	148	GLN
1	B	157	LEU
1	B	160	SER
1	B	162	THR
1	B	166	LEU
1	B	170	ILE

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Mol	Chain	Res	Type
1	B	172	SER
1	B	188	GLN
1	B	204	THR
1	B	212	LEU
1	B	236	LEU
1	B	238	LYS
1	B	250	LEU
1	B	253	VAL
1	B	255	LYS
1	B	257	ILE
1	B	258	HIS
1	B	259	SER
1	B	262	THR
1	B	263	ASN
1	B	264	ILE
1	B	278	THR
1	B	287	LYS
1	B	296	ASN
1	B	299	ILE
1	B	300	LEU
1	B	306	SER
1	B	310	LEU
1	B	312	ILE
1	B	333	ILE
1	B	335	ILE
1	B	337	GLN
1	B	345	LYS
1	B	348	LEU
1	B	351	THR
1	B	362	THR
1	B	397	GLU
1	B	401	LEU
1	B	407	LYS
1	B	411	GLU
1	B	412	GLU
1	B	418	LEU
1	B	420	GLN
1	B	421	LEU
1	C	37	VAL
1	C	39	ILE
1	C	58	LEU
1	C	62	VAL

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Mol	Chain	Res	Type
1	C	80	LYS
1	C	97	VAL
1	C	100	VAL
1	C	107	LEU
1	C	120	THR
1	C	123	ILE
1	C	125	ILE
1	C	132	ILE
1	C	138	LYS
1	C	140	VAL
1	C	143	LEU
1	C	145	MET
1	C	162	THR
1	C	166	LEU
1	C	172	SER
1	C	184	GLU
1	C	188	GLN
1	C	201	MET
1	C	203	GLU
1	C	204	THR
1	C	213	ILE
1	C	236	LEU
1	C	242	VAL
1	C	262	THR
1	C	263	ASN
1	C	264	ILE
1	C	269	LEU
1	C	278	THR
1	C	286	LEU
1	C	299	ILE
1	C	300	LEU
1	C	306	SER
1	C	310	LEU
1	C	319	LYS
1	C	320	ILE
1	C	333	ILE
1	C	335	ILE
1	C	337	GLN
1	C	340	GLU
1	C	345	LYS
1	C	348	LEU
1	C	349	GLN

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Mol	Chain	Res	Type
1	C	351	THR
1	C	362	THR
1	C	366	THR
1	C	371	MET
1	C	389	THR
1	C	401	LEU
1	C	405	VAL
1	C	407	LYS
1	C	412	GLU
1	C	418	LEU
1	C	421	LEU
1	C	425	ILE
1	C	430	ILE
1	D	38	LEU
1	D	51	THR
1	D	58	LEU
1	D	77	ARG
1	D	80	LYS
1	D	97	VAL
1	D	100	VAL
1	D	120	THR
1	D	123	ILE
1	D	132	ILE
1	D	140	VAL
1	D	143	LEU
1	D	147	THR
1	D	148	GLN
1	D	162	THR
1	D	166	LEU
1	D	170	ILE
1	D	184	GLU
1	D	185	ASP
1	D	204	THR
1	D	212	LEU
1	D	220	SER
1	D	239	ARG
1	D	242	VAL
1	D	250	LEU
1	D	262	THR
1	D	264	ILE
1	D	269	LEU
1	D	278	THR

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Mol	Chain	Res	Type
1	D	286	LEU
1	D	299	ILE
1	D	300	LEU
1	D	301	ILE
1	D	306	SER
1	D	310	LEU
1	D	333	ILE
1	D	335	ILE
1	D	348	LEU
1	D	366	THR
1	D	382	LEU
1	D	387	LEU
1	D	391	HIS
1	D	398	SER
1	D	401	LEU
1	D	407	LYS
1	D	414	SER
1	D	418	LEU
1	D	421	LEU
1	D	425	ILE
1	D	432	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	44	HIS
1	A	54	ASN
1	A	124	GLN
1	A	188	GLN
1	A	191	GLN
1	A	252	ASN
1	A	256	HIS
1	A	337	GLN
1	A	355	ASN
1	A	356	ASN
1	A	368	ASN
1	A	420	GLN
1	B	36	GLN
1	B	44	HIS
1	B	54	ASN
1	B	94	GLN

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Mol	Chain	Res	Type
1	B	174	GLN
1	B	189	HIS
1	B	252	ASN
1	B	279	ASN
1	B	296	ASN
1	B	355	ASN
1	B	356	ASN
1	B	420	GLN
1	C	36	GLN
1	C	54	ASN
1	C	104	ASN
1	C	124	GLN
1	C	133	ASN
1	C	252	ASN
1	C	318	ASN
1	C	355	ASN
1	C	356	ASN
1	C	420	GLN
1	D	54	ASN
1	D	124	GLN
1	D	148	GLN
1	D	177	ASN
1	D	191	GLN
1	D	318	ASN
1	D	355	ASN
1	D	356	ASN
1	D	431	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

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
2	GDP	A	593	3	29,30,30	1.20	3 (10%)	45,47,47	1.84	7 (15%)
2	GDP	D	593	3	29,30,30	1.21	2 (6%)	45,47,47	1.64	6 (13%)
2	GDP	B	593	3	29,30,30	1.26	4 (13%)	45,47,47	1.78	7 (15%)
2	GDP	C	593	3	29,30,30	1.22	4 (13%)	45,47,47	1.71	6 (13%)

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	GDP	A	593	3	-	4/16/32/32	0/3/3/3
2	GDP	D	593	3	-	8/16/32/32	0/3/3/3
2	GDP	B	593	3	-	10/16/32/32	0/3/3/3
2	GDP	C	593	3	-	3/16/32/32	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	593	GDP	C5-C4	3.62	1.48	1.38
2	D	593	GDP	C5-C4	3.40	1.48	1.38
2	C	593	GDP	C5-C4	3.10	1.47	1.38
2	A	593	GDP	C6-N1	-3.09	1.33	1.38
2	A	593	GDP	C5-C4	3.01	1.47	1.38
2	C	593	GDP	C6-N1	-2.88	1.33	1.38
2	B	593	GDP	C6-N1	-2.55	1.34	1.38
2	D	593	GDP	C6-N1	-2.54	1.34	1.38
2	B	593	GDP	C5-N7	-2.28	1.34	1.39
2	A	593	GDP	C5-N7	-2.25	1.34	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	593	GDP	PA-O3A	2.17	1.61	1.59
2	C	593	GDP	C4-N9	-2.14	1.32	1.38
2	C	593	GDP	C5-N7	-2.07	1.34	1.39

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	593	GDP	C5-C4-N3	-6.13	118.63	128.39
2	B	593	GDP	C5-C4-N3	-6.06	118.75	128.39
2	C	593	GDP	C5-C4-N3	-5.50	119.64	128.39
2	D	593	GDP	C5-C4-N3	-5.36	119.85	128.39
2	A	593	GDP	C2-N3-C4	4.87	120.68	112.30
2	D	593	GDP	C2-N3-C4	4.57	120.17	112.30
2	B	593	GDP	N9-C4-N3	4.53	135.01	125.95
2	B	593	GDP	C2-N3-C4	4.35	119.79	112.30
2	C	593	GDP	C2-N3-C4	4.32	119.73	112.30
2	D	593	GDP	C6-C5-N7	3.94	137.47	130.29
2	A	593	GDP	C6-C5-N7	3.93	137.44	130.29
2	C	593	GDP	N9-C4-N3	3.88	133.71	125.95
2	A	593	GDP	N9-C4-N3	3.74	133.43	125.95
2	D	593	GDP	N9-C4-N3	3.73	133.41	125.95
2	C	593	GDP	C6-C5-N7	3.30	136.30	130.29
2	B	593	GDP	C6-C5-N7	3.28	136.25	130.29
2	A	593	GDP	C4-C5-N7	-3.22	105.56	110.67
2	D	593	GDP	C4-C5-N7	-2.90	106.07	110.67
2	C	593	GDP	O3B-PB-O3A	2.89	114.34	104.64
2	B	593	GDP	C4-C5-N7	-2.61	106.54	110.67
2	C	593	GDP	C4-C5-N7	-2.54	106.64	110.67
2	A	593	GDP	O5'-PA-O1A	2.09	117.22	108.94
2	B	593	GDP	C5-C6-N1	2.09	118.57	113.25
2	D	593	GDP	C8-N7-C5	2.06	107.93	104.26
2	B	593	GDP	C2-N1-C6	-2.06	121.38	125.11
2	A	593	GDP	C5-C6-N1	2.05	118.47	113.25

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	593	GDP	C5'-O5'-PA-O3A
2	A	593	GDP	C5'-O5'-PA-O1A
2	B	593	GDP	C5'-O5'-PA-O3A
2	B	593	GDP	C5'-O5'-PA-O1A

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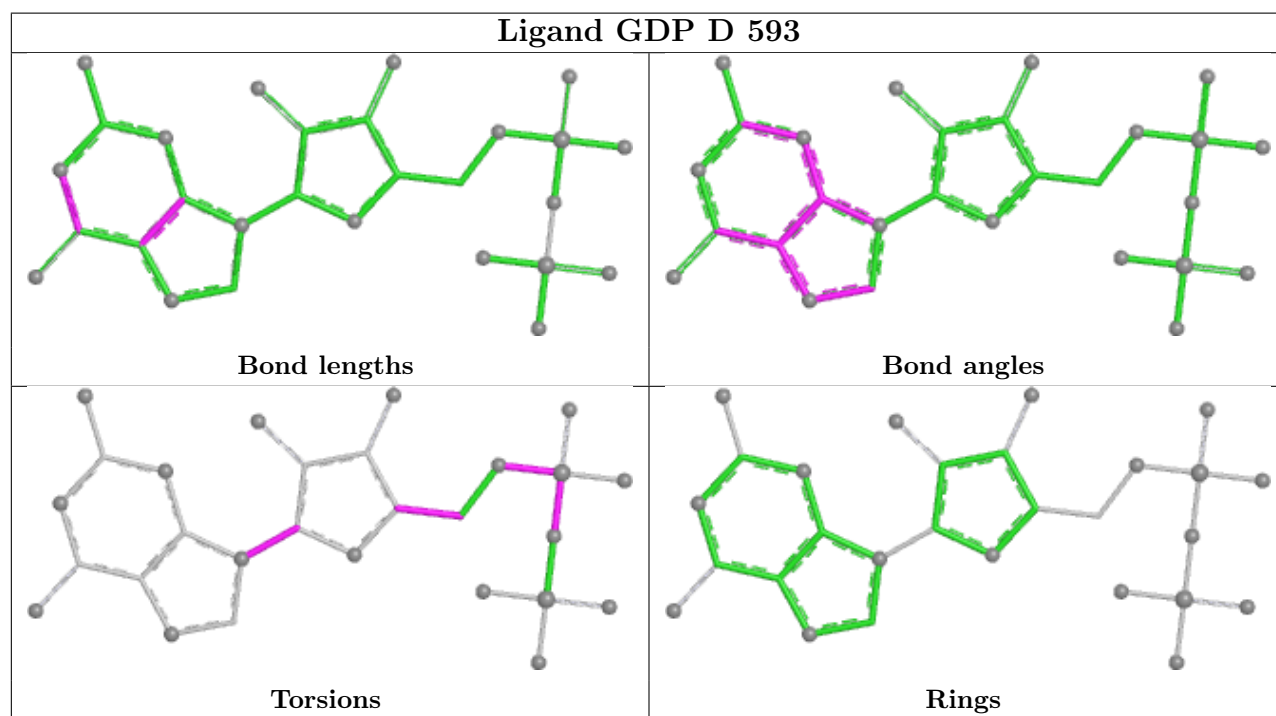
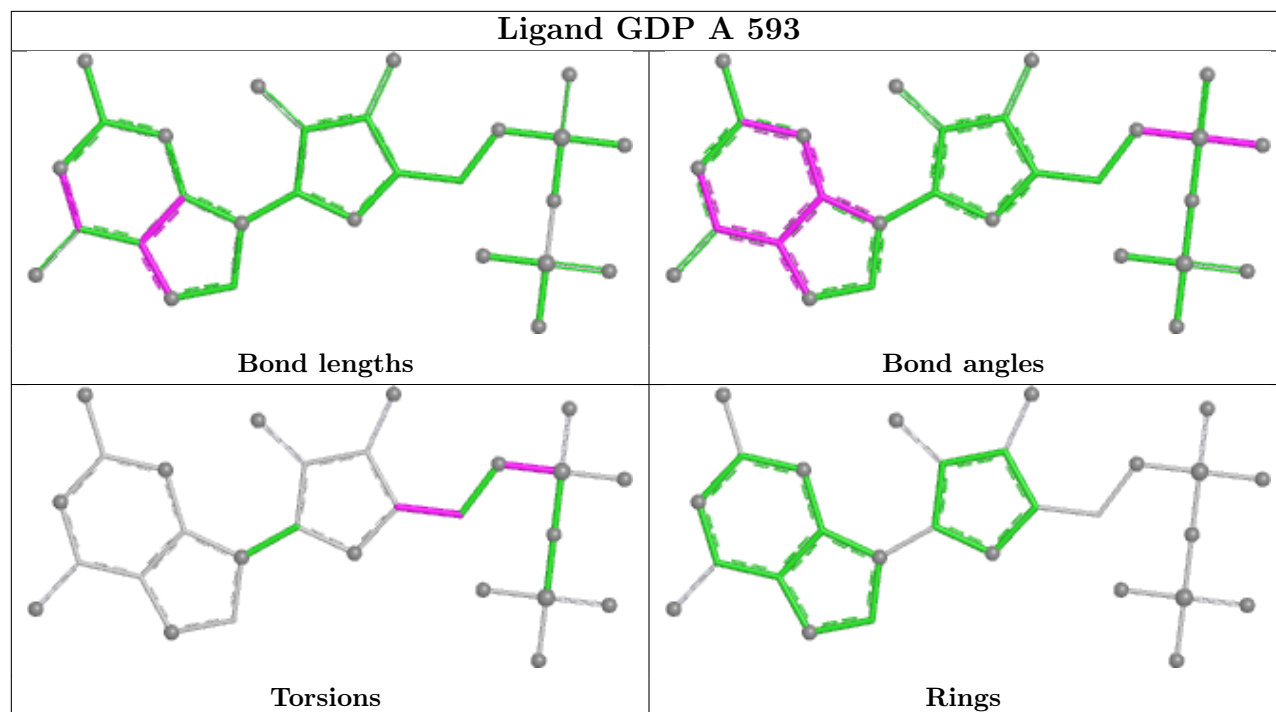
Mol	Chain	Res	Type	Atoms
2	B	593	GDP	C5'-O5'-PA-O2A
2	C	593	GDP	C5'-O5'-PA-O3A
2	C	593	GDP	C5'-O5'-PA-O1A
2	D	593	GDP	C5'-O5'-PA-O3A
2	D	593	GDP	C5'-O5'-PA-O1A
2	A	593	GDP	O4'-C4'-C5'-O5'
2	B	593	GDP	O4'-C4'-C5'-O5'
2	D	593	GDP	O4'-C4'-C5'-O5'
2	B	593	GDP	C3'-C4'-C5'-O5'
2	A	593	GDP	C3'-C4'-C5'-O5'
2	B	593	GDP	PA-O3A-PB-O3B
2	D	593	GDP	PB-O3A-PA-O1A
2	C	593	GDP	C5'-O5'-PA-O2A
2	D	593	GDP	C5'-O5'-PA-O2A
2	D	593	GDP	C2'-C1'-N9-C8
2	D	593	GDP	C2'-C1'-N9-C4
2	B	593	GDP	C2'-C1'-N9-C4
2	B	593	GDP	PB-O3A-PA-O1A
2	B	593	GDP	C2'-C1'-N9-C8
2	B	593	GDP	PB-O3A-PA-O2A
2	D	593	GDP	PB-O3A-PA-O2A

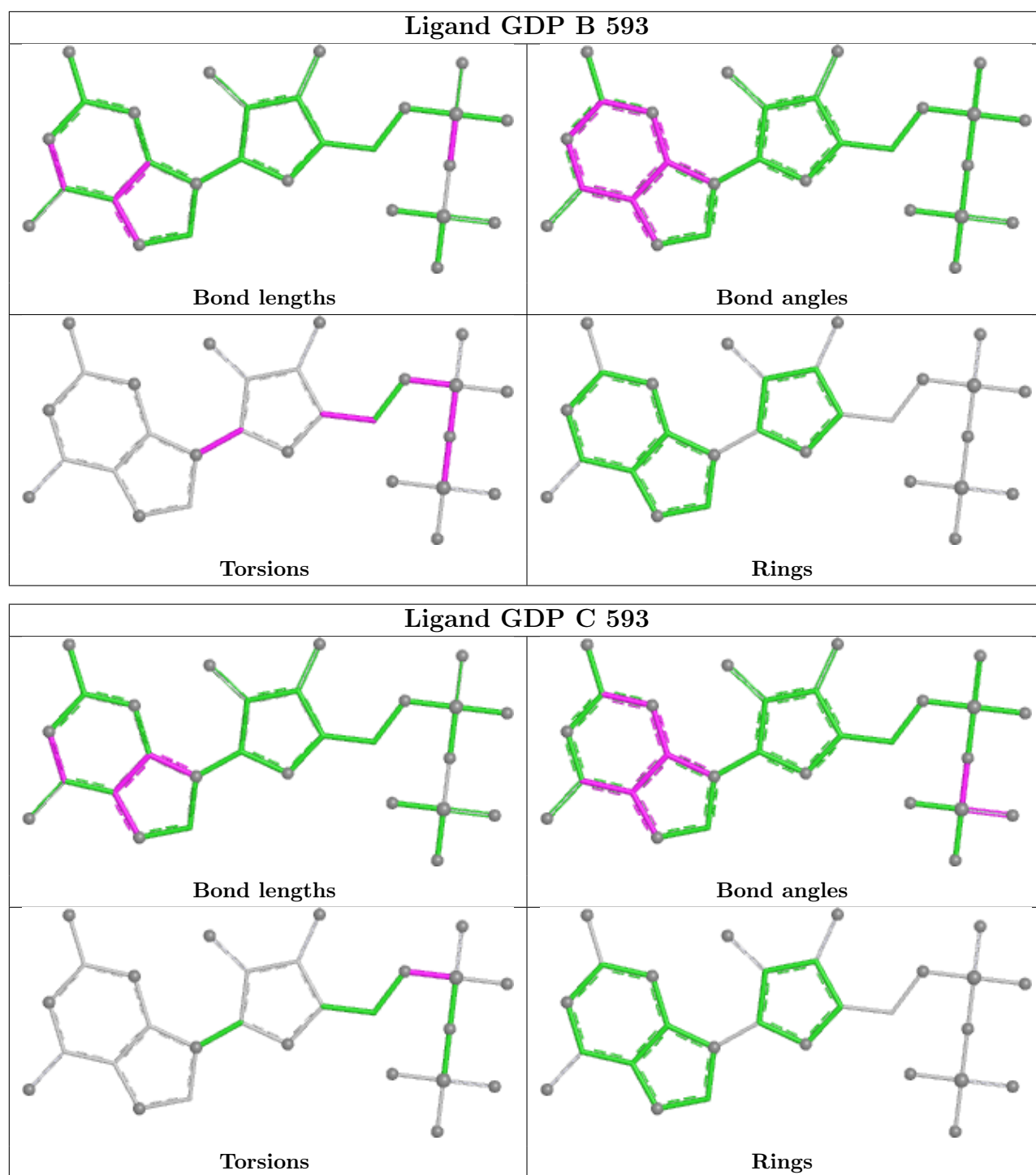
There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	593	GDP	4	0
2	B	593	GDP	2	0
2	C	593	GDP	3	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	404/459 (88%)	0.53	32 (7%) 18 13	27, 71, 119, 147	0
1	B	397/459 (86%)	0.42	24 (6%) 27 21	31, 68, 138, 152	0
1	C	405/459 (88%)	0.22	28 (6%) 23 16	32, 53, 106, 158	0
1	D	398/459 (86%)	0.51	34 (8%) 16 12	34, 69, 140, 159	0
All	All	1604/1836 (87%)	0.42	118 (7%) 20 15	27, 65, 132, 159	0

All (118) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	382	LEU	5.1
1	A	157	LEU	4.8
1	B	156	THR	4.8
1	D	159	ASP	4.7
1	A	152	ASP	4.6
1	C	150	THR	4.5
1	C	157	LEU	4.5
1	D	157	LEU	4.5
1	B	155	SER	4.5
1	D	247	HIS	4.4
1	C	242	VAL	4.3
1	C	153	SER	4.3
1	D	153	SER	4.2
1	D	40	VAL	4.1
1	D	432	TYR	4.0
1	B	242	VAL	4.0
1	A	438	SER	3.9
1	C	156	THR	3.8
1	D	338	GLY	3.8
1	C	155	SER	3.8
1	C	114	GLY	3.8

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Mol	Chain	Res	Type	RSRZ
1	D	155	SER	3.8
1	D	337	GLN	3.7
1	C	152	ASP	3.6
1	D	341	LEU	3.5
1	D	251	GLN	3.5
1	A	40	VAL	3.5
1	A	120	THR	3.5
1	A	252	ASN	3.5
1	C	247	HIS	3.4
1	B	432	TYR	3.4
1	B	104	ASN	3.3
1	C	118	ARG	3.2
1	A	156	THR	3.2
1	D	152	ASP	3.2
1	D	39	ILE	3.1
1	B	40	VAL	3.1
1	B	117	GLU	3.1
1	D	380	PRO	3.0
1	B	107	LEU	3.0
1	C	154	GLN	3.0
1	B	382	LEU	2.9
1	A	39	ILE	2.9
1	C	383	ALA	2.9
1	B	430	ILE	2.8
1	D	151	PHE	2.8
1	C	258	HIS	2.8
1	B	375	CYS	2.7
1	A	148	GLN	2.7
1	B	385	ASN	2.7
1	D	242	VAL	2.7
1	B	46	PHE	2.7
1	B	383	ALA	2.7
1	C	382	LEU	2.7
1	D	281	ASN	2.7
1	B	157	LEU	2.7
1	A	341	LEU	2.6
1	C	385	ASN	2.6
1	C	107	LEU	2.6
1	A	154	GLN	2.6
1	D	46	PHE	2.5
1	C	113	ARG	2.5
1	C	110	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	335	ILE	2.4
1	D	149	GLY	2.4
1	D	339	GLU	2.4
1	D	383	ALA	2.4
1	A	47	GLU	2.4
1	A	38	LEU	2.4
1	B	58	LEU	2.4
1	C	116	SER	2.3
1	D	430	ILE	2.3
1	B	340	GLU	2.3
1	C	117	GLU	2.3
1	D	42	ASP	2.3
1	D	374	ILE	2.3
1	A	363	ALA	2.3
1	D	375	CYS	2.3
1	A	161	ALA	2.3
1	A	283	ASP	2.3
1	A	401	LEU	2.3
1	C	148	GLN	2.3
1	D	283	ASP	2.3
1	D	150	THR	2.3
1	C	338	GLY	2.3
1	D	154	GLN	2.2
1	A	116	SER	2.2
1	D	336	TYR	2.2
1	D	115	GLY	2.2
1	B	31	LYS	2.2
1	A	123	ILE	2.2
1	C	283	ASP	2.2
1	D	386	ASP	2.2
1	A	332	TYR	2.2
1	A	382	LEU	2.2
1	A	118	ARG	2.2
1	A	170	ILE	2.2
1	C	151	PHE	2.2
1	B	194	THR	2.2
1	C	109	GLY	2.2
1	B	76	PHE	2.2
1	A	36	GLN	2.1
1	C	149	GLY	2.1
1	B	250	LEU	2.1
1	A	117	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	113	ARG	2.1
1	A	159	ASP	2.1
1	A	153	SER	2.1
1	A	44	HIS	2.1
1	D	187	LEU	2.1
1	B	120	THR	2.1
1	A	383	ALA	2.1
1	A	339	GLU	2.1
1	A	406	LYS	2.0
1	C	104	ASN	2.0
1	B	372	GLU	2.0
1	A	77	ARG	2.0
1	C	40	VAL	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.

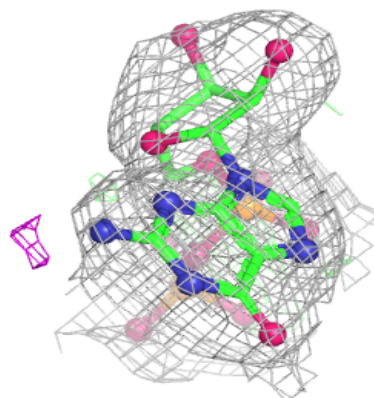
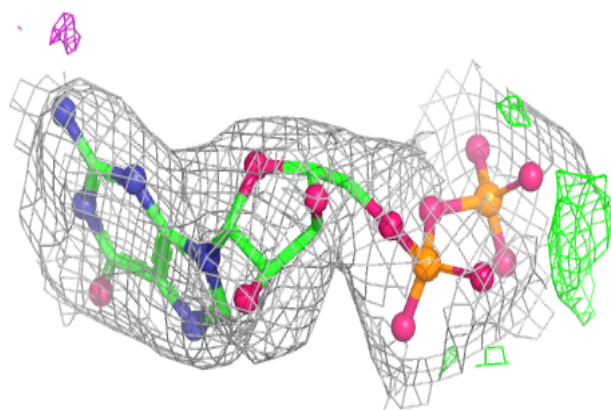
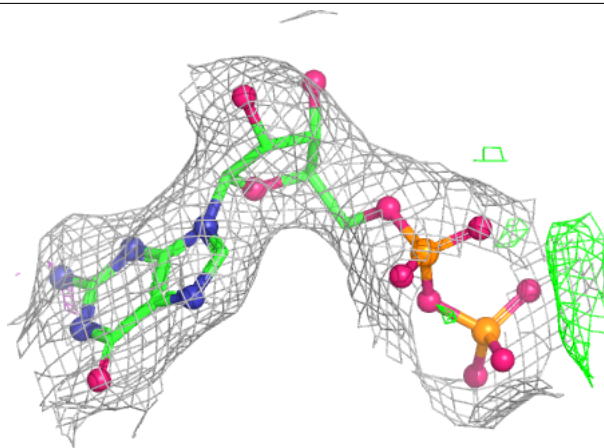
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	MG	C	595	1/1	0.83	0.17	57,57,57,57	0
3	MG	D	595	1/1	0.86	0.15	69,69,69,69	0
3	MG	A	595	1/1	0.90	0.26	68,68,68,68	0
3	MG	B	595	1/1	0.91	0.23	70,70,70,70	0
2	GDP	A	593	28/28	0.94	0.07	35,54,69,75	0
2	GDP	B	593	28/28	0.94	0.08	32,46,70,76	0
2	GDP	C	593	28/28	0.95	0.08	38,47,56,65	0
2	GDP	D	593	28/28	0.96	0.07	35,49,61,69	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.

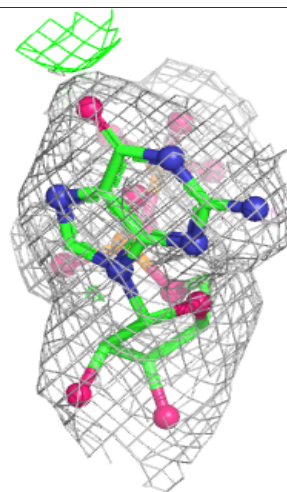
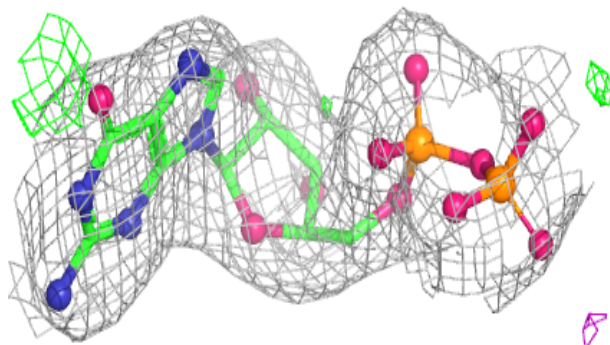
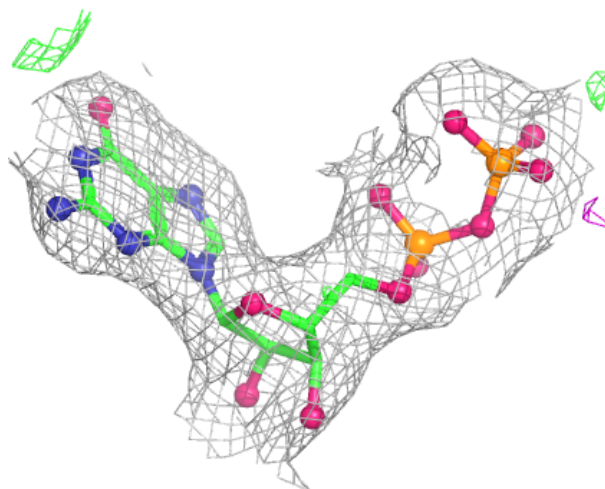
Electron density around GDP A 593:

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



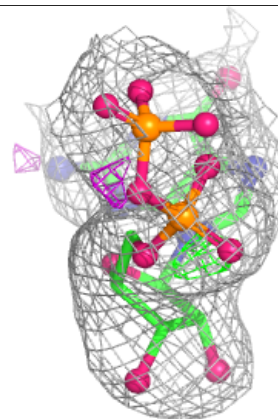
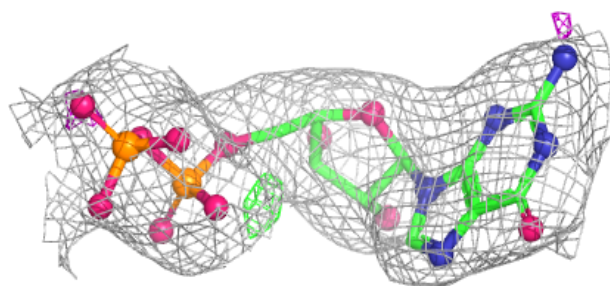
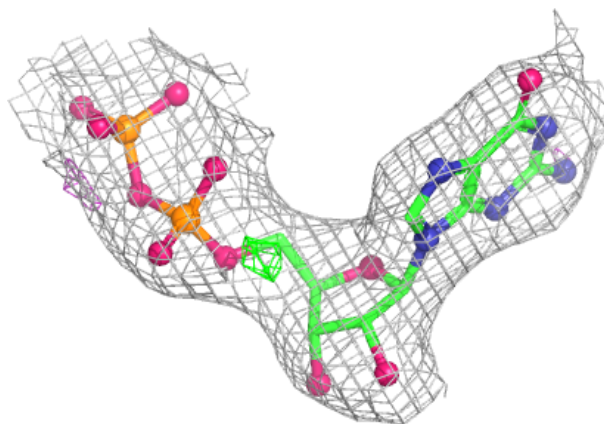
Electron density around GDP B 593:

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

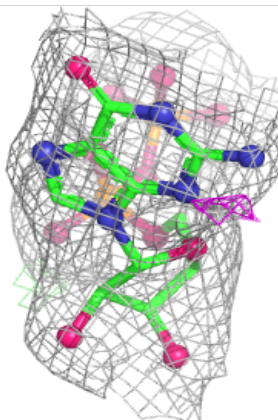
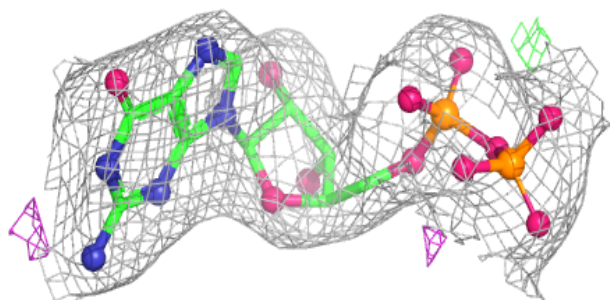
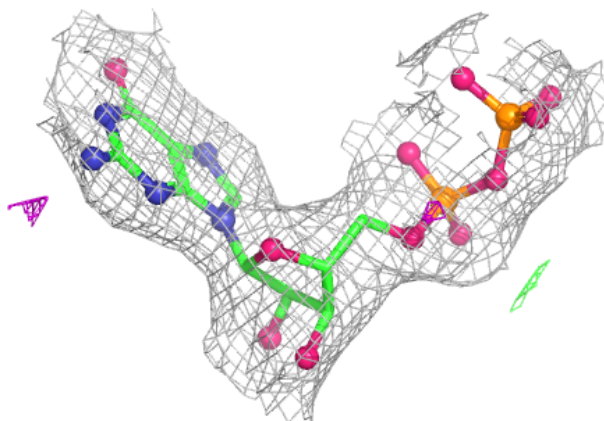


Electron density around GDP C 593:

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

**Electron density around GDP D 593:**

$2mF_o-DF_c$ (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.