



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 24, 2025 – 11:28 pm BST

PDB ID : 8S0G / pdb_00008s0g
Title : Crystal structure of Renilla reniformis luciferase-GFP BRET complex
Authors : Marek, M.
Deposited on : 2024-02-14
Resolution : 2.38 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 Phenix2.0rc1
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.45.1

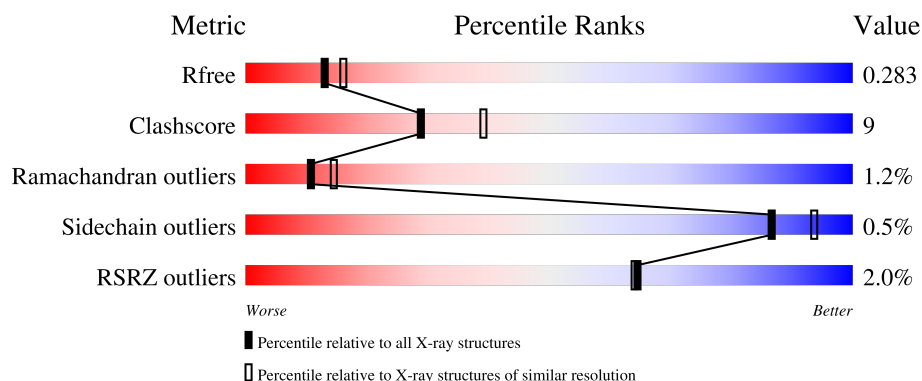
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.38 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}	164625	6699 (2.40-2.36)
Clashscore	180529	7414 (2.40-2.36)
Ramachandran outliers	177936	7337 (2.40-2.36)
Sidechain outliers	177891	7338 (2.40-2.36)
RSRZ outliers	164620	6699 (2.40-2.36)

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	231	<div> <div>0%</div> <div> <div></div> <div>76%</div> <div>23%</div> <div>.</div> </div> </div>
1	B	231	<div> <div>0%</div> <div> <div></div> <div>80%</div> <div>19%</div> <div>.</div> </div> </div>
1	E	231	<div> <div>4%</div> <div> <div></div> <div>76%</div> <div>20%</div> <div>..</div> </div> </div>
1	F	231	<div> <div>4%</div> <div> <div></div> <div>71%</div> <div>26%</div> <div>..</div> </div> </div>
1	I	231	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>21%</div> <div>..</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	J	231	 79%19%
2	C	319	 78%17%
2	D	319	 81%16%
2	G	319	 80%17%
2	H	319	 75%21%
2	K	319	 72%24%
2	L	319	 76%21%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 26441 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 Green fluorescent protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	230	Total	C	N	O	S	0	0	0
			1820	1166	300	345	9			
1	B	231	Total	C	N	O	S	0	0	0
			1828	1171	301	346	10			
1	E	229	Total	C	N	O	S	0	1	0
			1821	1167	300	345	9			
1	F	228	Total	C	N	O	S	0	0	0
			1804	1156	298	341	9			
1	I	226	Total	C	N	O	S	0	0	0
			1790	1147	295	339	9			
1	J	230	Total	C	N	O	S	0	0	0
			1820	1166	300	345	9			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	67	CRO	SER	chromophore	UNP Q963I9
A	67	CRO	TYR	chromophore	UNP Q963I9
A	67	CRO	GLY	chromophore	UNP Q963I9
B	67	CRO	SER	chromophore	UNP Q963I9
B	67	CRO	TYR	chromophore	UNP Q963I9
B	67	CRO	GLY	chromophore	UNP Q963I9
E	67	CRO	SER	chromophore	UNP Q963I9
E	67	CRO	TYR	chromophore	UNP Q963I9
E	67	CRO	GLY	chromophore	UNP Q963I9
F	67	CRO	SER	chromophore	UNP Q963I9
F	67	CRO	TYR	chromophore	UNP Q963I9
F	67	CRO	GLY	chromophore	UNP Q963I9
I	67	CRO	SER	chromophore	UNP Q963I9
I	67	CRO	TYR	chromophore	UNP Q963I9
I	67	CRO	GLY	chromophore	UNP Q963I9
J	67	CRO	SER	chromophore	UNP Q963I9
J	67	CRO	TYR	chromophore	UNP Q963I9

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Chain	Residue	Modelled	Actual	Comment	Reference
J	67	CRO	GLY	chromophore	UNP Q963I9

- Molecule 2 is a protein called Coelenterazine h 2-monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	306	Total	C	N	O	S	0	0	0
			2510	1628	421	452	9			
2	D	309	Total	C	N	O	S	0	0	0
			2532	1640	424	459	9			
2	G	312	Total	C	N	O	S	0	0	0
			2551	1650	428	464	9			
2	H	308	Total	C	N	O	S	0	0	0
			2523	1635	423	456	9			
2	K	307	Total	C	N	O	S	0	0	0
			2515	1631	421	454	9			
2	L	308	Total	C	N	O	S	0	0	0
			2523	1635	423	456	9			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	55	THR	ALA	engineered mutation	UNP P27652
C	124	ALA	CYS	engineered mutation	UNP P27652
C	130	ALA	SER	engineered mutation	UNP P27652
C	136	ARG	LYS	engineered mutation	UNP P27652
C	143	MET	ALA	engineered mutation	UNP P27652
C	185	VAL	MET	engineered mutation	UNP P27652
C	253	LEU	MET	engineered mutation	UNP P27652
C	287	LEU	SER	engineered mutation	UNP P27652
C	312	SER	-	expression tag	UNP P27652
C	313	GLY	-	expression tag	UNP P27652
C	314	LEU	-	expression tag	UNP P27652
C	315	GLU	-	expression tag	UNP P27652
C	316	VAL	-	expression tag	UNP P27652
C	317	LEU	-	expression tag	UNP P27652
C	318	PHE	-	expression tag	UNP P27652
C	319	GLN	-	expression tag	UNP P27652
D	55	THR	ALA	engineered mutation	UNP P27652
D	124	ALA	CYS	engineered mutation	UNP P27652
D	130	ALA	SER	engineered mutation	UNP P27652
D	136	ARG	LYS	engineered mutation	UNP P27652
D	143	MET	ALA	engineered mutation	UNP P27652

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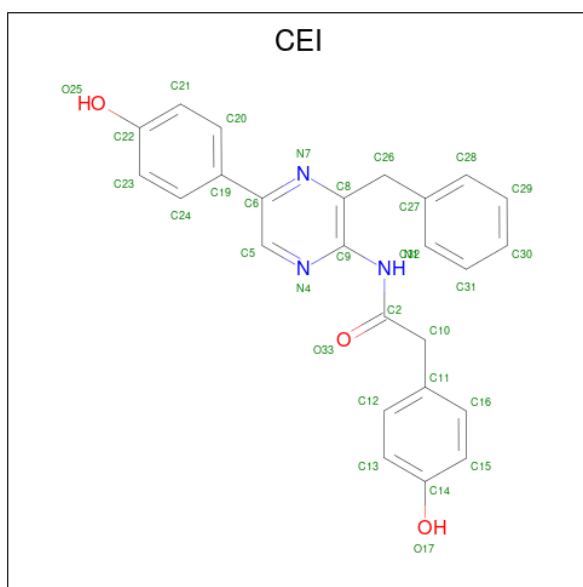
Chain	Residue	Modelled	Actual	Comment	Reference
D	185	VAL	MET	engineered mutation	UNP P27652
D	253	LEU	MET	engineered mutation	UNP P27652
D	287	LEU	SER	engineered mutation	UNP P27652
D	312	SER	-	expression tag	UNP P27652
D	313	GLY	-	expression tag	UNP P27652
D	314	LEU	-	expression tag	UNP P27652
D	315	GLU	-	expression tag	UNP P27652
D	316	VAL	-	expression tag	UNP P27652
D	317	LEU	-	expression tag	UNP P27652
D	318	PHE	-	expression tag	UNP P27652
D	319	GLN	-	expression tag	UNP P27652
G	55	THR	ALA	engineered mutation	UNP P27652
G	124	ALA	CYS	engineered mutation	UNP P27652
G	130	ALA	SER	engineered mutation	UNP P27652
G	136	ARG	LYS	engineered mutation	UNP P27652
G	143	MET	ALA	engineered mutation	UNP P27652
G	185	VAL	MET	engineered mutation	UNP P27652
G	253	LEU	MET	engineered mutation	UNP P27652
G	287	LEU	SER	engineered mutation	UNP P27652
G	312	SER	-	expression tag	UNP P27652
G	313	GLY	-	expression tag	UNP P27652
G	314	LEU	-	expression tag	UNP P27652
G	315	GLU	-	expression tag	UNP P27652
G	316	VAL	-	expression tag	UNP P27652
G	317	LEU	-	expression tag	UNP P27652
G	318	PHE	-	expression tag	UNP P27652
G	319	GLN	-	expression tag	UNP P27652
H	55	THR	ALA	engineered mutation	UNP P27652
H	124	ALA	CYS	engineered mutation	UNP P27652
H	130	ALA	SER	engineered mutation	UNP P27652
H	136	ARG	LYS	engineered mutation	UNP P27652
H	143	MET	ALA	engineered mutation	UNP P27652
H	185	VAL	MET	engineered mutation	UNP P27652
H	253	LEU	MET	engineered mutation	UNP P27652
H	287	LEU	SER	engineered mutation	UNP P27652
H	312	SER	-	expression tag	UNP P27652
H	313	GLY	-	expression tag	UNP P27652
H	314	LEU	-	expression tag	UNP P27652
H	315	GLU	-	expression tag	UNP P27652
H	316	VAL	-	expression tag	UNP P27652
H	317	LEU	-	expression tag	UNP P27652
H	318	PHE	-	expression tag	UNP P27652

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Chain	Residue	Modelled	Actual	Comment	Reference
H	319	GLN	-	expression tag	UNP P27652
K	55	THR	ALA	engineered mutation	UNP P27652
K	124	ALA	CYS	engineered mutation	UNP P27652
K	130	ALA	SER	engineered mutation	UNP P27652
K	136	ARG	LYS	engineered mutation	UNP P27652
K	143	MET	ALA	engineered mutation	UNP P27652
K	185	VAL	MET	engineered mutation	UNP P27652
K	253	LEU	MET	engineered mutation	UNP P27652
K	287	LEU	SER	engineered mutation	UNP P27652
K	312	SER	-	expression tag	UNP P27652
K	313	GLY	-	expression tag	UNP P27652
K	314	LEU	-	expression tag	UNP P27652
K	315	GLU	-	expression tag	UNP P27652
K	316	VAL	-	expression tag	UNP P27652
K	317	LEU	-	expression tag	UNP P27652
K	318	PHE	-	expression tag	UNP P27652
K	319	GLN	-	expression tag	UNP P27652
L	55	THR	ALA	engineered mutation	UNP P27652
L	124	ALA	CYS	engineered mutation	UNP P27652
L	130	ALA	SER	engineered mutation	UNP P27652
L	136	ARG	LYS	engineered mutation	UNP P27652
L	143	MET	ALA	engineered mutation	UNP P27652
L	185	VAL	MET	engineered mutation	UNP P27652
L	253	LEU	MET	engineered mutation	UNP P27652
L	287	LEU	SER	engineered mutation	UNP P27652
L	312	SER	-	expression tag	UNP P27652
L	313	GLY	-	expression tag	UNP P27652
L	314	LEU	-	expression tag	UNP P27652
L	315	GLU	-	expression tag	UNP P27652
L	316	VAL	-	expression tag	UNP P27652
L	317	LEU	-	expression tag	UNP P27652
L	318	PHE	-	expression tag	UNP P27652
L	319	GLN	-	expression tag	UNP P27652

- Molecule 3 is N-[3-BENZYL-5-(4-HYDROXYPHENYL)PYRAZIN-2-YL]-2-(4-HYDROXYPHENYL)ACETAMIDE (CCD ID: CEI) (formula: C₂₅H₂₁N₃O₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	L	1	Total	C	N	O	0	0
			31	25	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	34	Total	O	0	1
			35	35		
4	B	40	Total	O	0	0
			40	40		
4	C	47	Total	O	0	0
			47	47		
4	D	37	Total	O	0	0
			37	37		
4	E	30	Total	O	0	0
			30	30		
4	F	20	Total	O	0	0
			20	20		
4	G	56	Total	O	0	1
			57	57		
4	H	22	Total	O	0	0
			22	22		
4	I	18	Total	O	0	0
			18	18		
4	J	22	Total	O	0	0
			22	22		
4	K	18	Total	O	0	0
			18	18		

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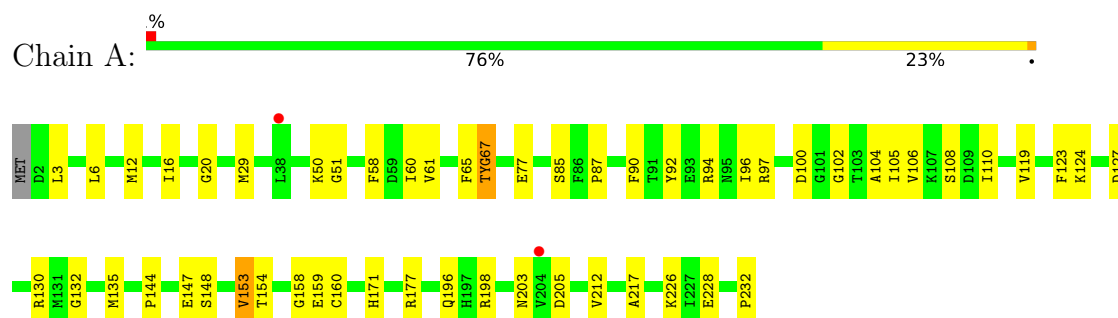
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	27	Total	O	0	0
			27	27		

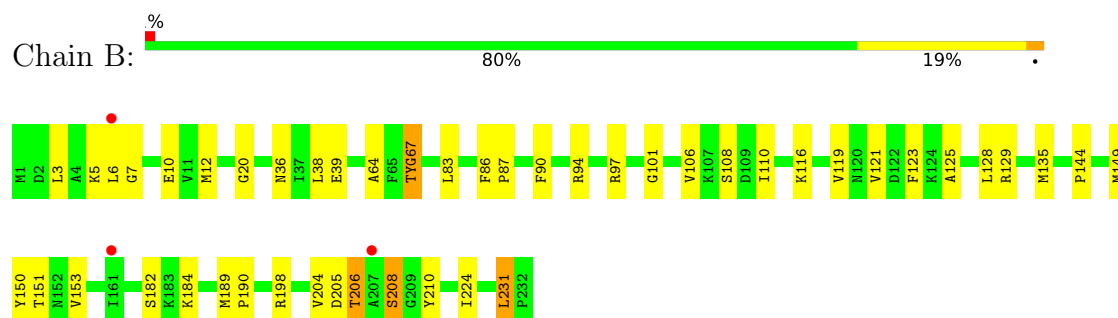
3 Residue-property plots

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.

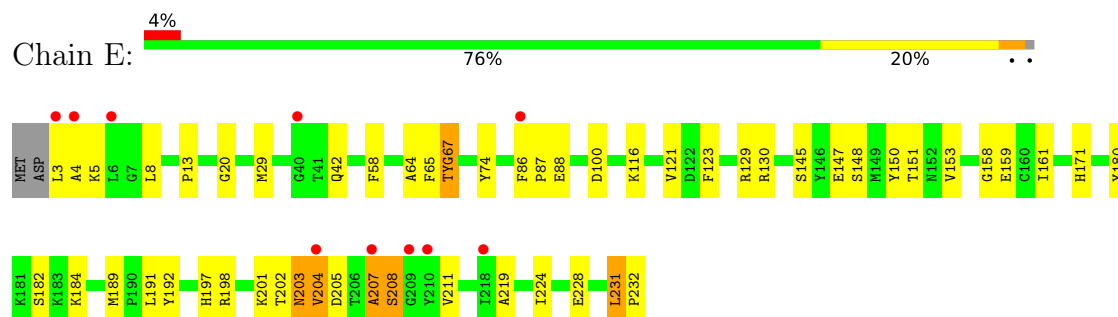
• Molecule 1: Green fluorescent protein



• Molecule 1: Green fluorescent protein

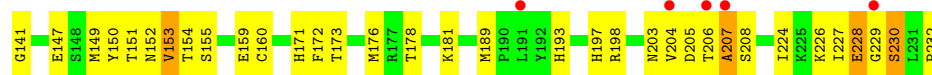


• Molecule 1: Green fluorescent protein

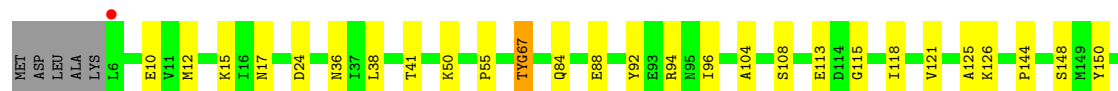
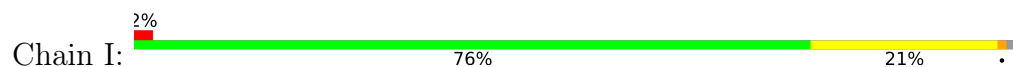


• Molecule 1: Green fluorescent protein

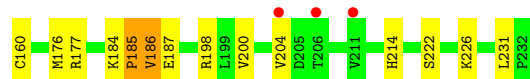




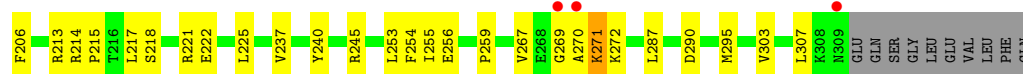
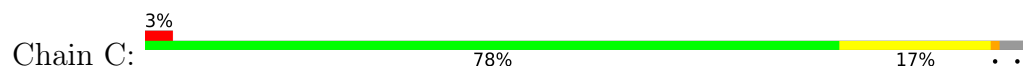
• Molecule 1: Green fluorescent protein



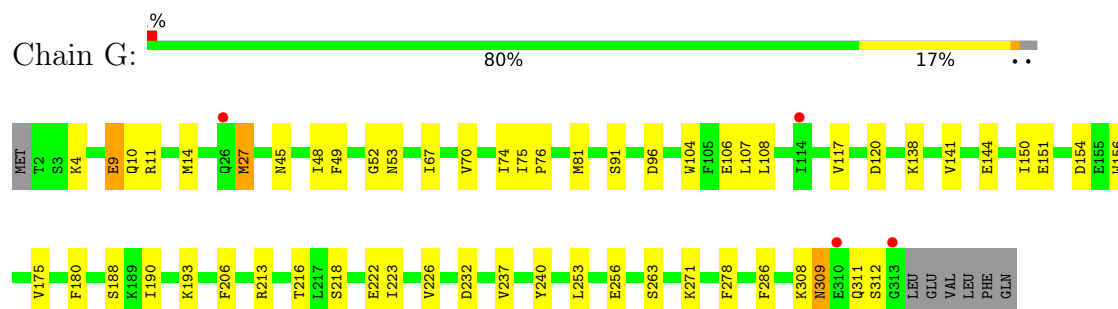
• Molecule 1: Green fluorescent protein



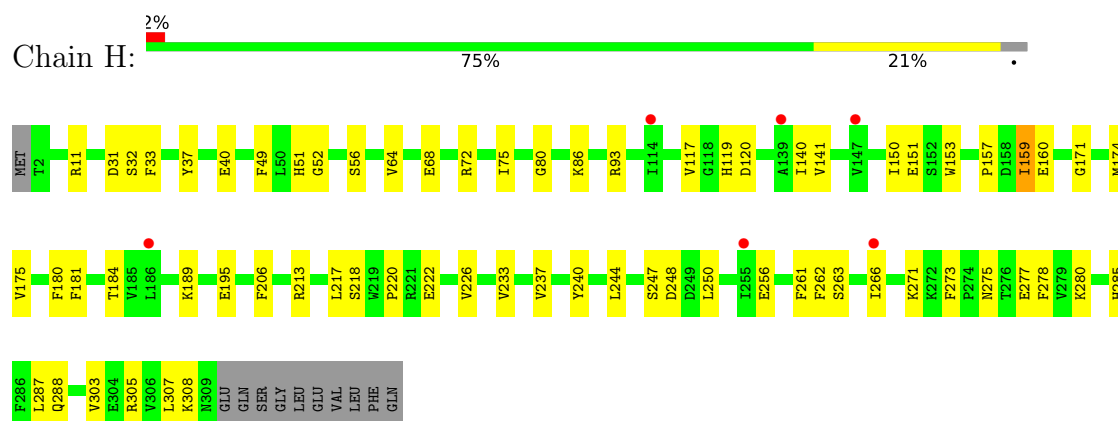
• Molecule 2: Coelenterazine h 2-monooxygenase



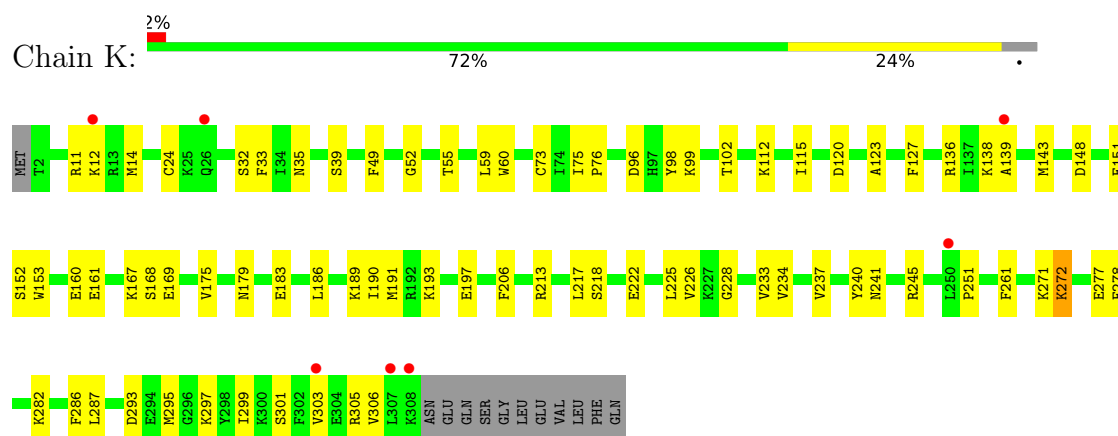
• Molecule 2: Coelenterazine h 2-monooxygenase



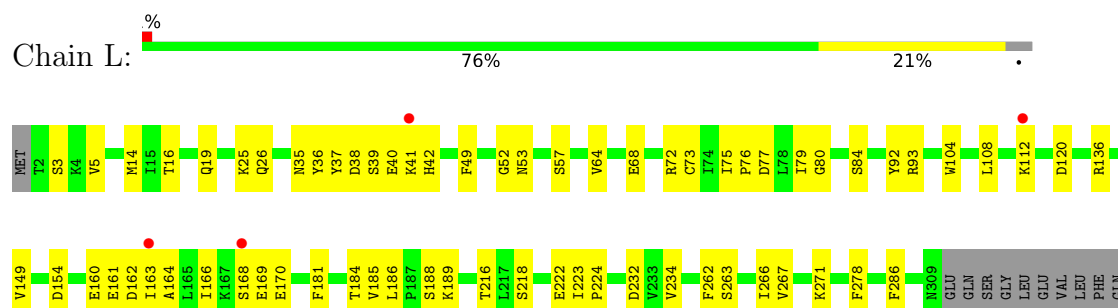
• Molecule 2: Coelenterazine h 2-monooxygenase



• Molecule 2: Coelenterazine h 2-monooxygenase



• Molecule 2: Coelenterazine h 2-monooxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	96.88Å 96.88Å 361.46Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.44 – 2.38 48.44 – 2.38	Depositor EDS
% Data completeness (in resolution range)	99.5 (48.44-2.38) 99.5 (48.44-2.38)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.06 (at 2.37Å)	Xtriage
Refinement program	PHENIX (1.14_3260)	Depositor
R, R_{free}	0.220 , 0.281 0.224 , 0.283	Depositor DCC
R_{free} test set	7233 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å ²)	46.8	Xtriage
Anisotropy	0.571	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.016 for -h,-k,l 0.092 for h,-h-k,-l 0.042 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	26441	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.16% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.21	0/1837	0.42	0/2479
1	B	0.23	0/1845	0.46	0/2489
1	E	0.20	0/1838	0.40	0/2480
1	F	0.23	0/1821	0.51	1/2457 (0.0%)
1	I	0.16	0/1807	0.39	0/2439
1	J	0.22	0/1837	0.45	0/2479
2	C	0.27	0/2582	0.48	2/3496 (0.1%)
2	D	0.13	0/2604	0.34	0/3526
2	G	0.25	0/2623	0.46	3/3551 (0.1%)
2	H	0.14	0/2595	0.37	0/3514
2	K	0.16	0/2587	0.37	0/3503
2	L	0.23	1/2595 (0.0%)	0.40	1/3514 (0.0%)
All	All	0.21	1/26571 (0.0%)	0.42	7/35927 (0.0%)

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	E	0	1
2	G	0	1
2	L	0	1
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	41	LYS	CB-CG	-5.45	1.36	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	138	ASP	CB-CA-C	-8.34	96.63	110.72
2	C	13	ARG	N-CA-C	-7.63	101.13	113.19
2	G	309	ASN	N-CA-CB	7.46	123.09	110.49
2	C	271	LYS	CD-CE-NZ	-7.39	88.24	111.90
2	G	308	LYS	CA-C-N	5.92	132.86	121.54
2	G	308	LYS	C-N-CA	5.92	132.86	121.54
2	L	41	LYS	CA-CB-CG	-5.81	102.49	114.10

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	207	ALA	Peptide
2	G	312	SER	Peptide
2	L	40	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	1820	0	1803	34	0
1	B	1828	0	1815	31	0
1	E	1821	0	1804	44	0
1	F	1804	0	1788	43	0
1	I	1790	0	1770	33	0
1	J	1820	0	1800	31	0
2	C	2510	0	2476	46	0
2	D	2532	0	2494	36	0
2	G	2551	0	2510	34	0
2	H	2523	0	2488	42	0
2	K	2515	0	2482	54	0
2	L	2523	0	2488	43	0
3	L	31	0	21	8	0
4	A	35	0	0	1	0
4	B	40	0	0	0	0
4	C	47	0	0	1	0
4	D	37	0	0	2	0
4	E	30	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	F	20	0	0	0	0
4	G	57	0	0	0	0
4	H	22	0	0	0	0
4	I	18	0	0	1	0
4	J	22	0	0	1	0
4	K	18	0	0	1	0
4	L	27	0	0	0	0
All	All	26441	0	25739	443	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 9.

All (443) 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:E:42:GLN:HE22	1:E:67:CRO:CB1	1.80	0.94
1:E:42:GLN:NE2	1:E:67:CRO:CB1	2.32	0.92
2:C:112:LYS:H	2:C:112:LYS:HD3	1.32	0.92
2:K:148:ASP:OD1	2:K:241:ASN:CG	2.15	0.89
1:A:159:GLU:OE1	4:A:301:HOH:O	1.97	0.83
2:D:256:GLU:HB3	2:D:280:LYS:HG2	1.62	0.81
2:H:52:GLY:HA3	2:H:120:ASP:HB3	1.64	0.80
2:L:168:SER:O	2:L:170:GLU:N	2.14	0.79
2:C:206:PHE:HB3	2:C:213:ARG:HG2	1.66	0.77
2:C:267:VAL:O	2:C:271:LYS:HB2	1.82	0.77
2:C:177:GLU:O	2:C:177:GLU:OE1	2.02	0.77
1:F:8:LEU:HB3	1:F:12:MET:HE1	1.68	0.76
1:F:13:PRO:HG2	1:F:116:LYS:HG3	1.70	0.74
2:H:206:PHE:HB3	2:H:213:ARG:HG2	1.69	0.73
2:K:39:SER:OG	2:K:73:CYS:N	2.20	0.73
1:J:144:PRO:HG3	1:J:198:ARG:HG3	1.69	0.73
1:A:144:PRO:HG3	1:A:198:ARG:HG3	1.71	0.72
2:D:253:LEU:HD11	2:D:279:VAL:HG22	1.71	0.72
1:J:153:VAL:HG23	1:J:154:THR:H	1.55	0.71
1:F:29:MET:HE1	1:F:61:VAL:HG21	1.72	0.71
1:I:144:PRO:HG3	1:I:198:ARG:HG3	1.72	0.71
2:G:141:VAL:HG22	2:G:253:LEU:HB3	1.72	0.71
2:C:112:LYS:HD3	2:C:112:LYS:N	2.06	0.70
2:C:173:LYS:HE2	2:C:177:GLU:HG3	1.71	0.70
1:E:42:GLN:NE2	1:E:67:CRO:OG1	2.24	0.70
1:B:3:LEU:O	1:B:5:LYS:HG3	1.92	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:3:LEU:HA	1:E:5:LYS:HE2	1.75	0.69
2:K:179:ASN:HB3	2:K:183:GLU:HG3	1.74	0.68
1:E:198:ARG:NH1	1:F:224:ILE:O	2.25	0.68
1:E:100:ASP:O	1:E:129:ARG:NH1	2.27	0.68
2:K:143:MET:HG3	2:K:287:LEU:HD21	1.76	0.67
2:G:206:PHE:HB3	2:G:213:ARG:HG2	1.77	0.67
1:E:151:THR:HG21	1:E:189:MET:HB3	1.78	0.65
2:L:49:PHE:HB2	2:L:75:ILE:HG23	1.78	0.65
1:A:148:SER:O	1:A:158:GLY:HA2	1.96	0.65
1:F:153:VAL:HG22	1:F:154:THR:HG23	1.78	0.65
2:G:4:LYS:HZ3	2:G:188:SER:HA	1.61	0.65
2:K:148:ASP:OD1	2:K:241:ASN:ND2	2.29	0.65
2:C:5:VAL:HG23	2:C:6:TYR:H	1.62	0.65
1:E:13:PRO:HG2	1:E:116:LYS:HG2	1.79	0.64
2:C:112:LYS:H	2:C:112:LYS:CD	1.97	0.64
2:L:93:ARG:NH2	2:L:232:ASP:OD2	2.30	0.64
1:F:227:ILE:HB	1:F:230:SER:HB3	1.81	0.63
1:I:88:GLU:HG3	1:I:184:LYS:HB2	1.81	0.63
1:J:48:VAL:HG21	2:L:5:VAL:HG11	1.80	0.63
2:D:52:GLY:HA3	2:D:120:ASP:HB3	1.81	0.62
2:H:49:PHE:HB2	2:H:75:ILE:HG12	1.81	0.62
2:C:255:ILE:HD13	2:C:295:MET:HG3	1.82	0.62
2:D:93:ARG:NH2	2:D:232:ASP:OD2	2.33	0.62
1:I:206:THR:O	2:K:193:LYS:NZ	2.33	0.61
1:I:198:ARG:HH21	1:J:222:SER:HB3	1.65	0.61
1:F:204:VAL:HG22	1:F:205:ASP:H	1.65	0.61
2:C:147:VAL:HG21	2:C:270:ALA:HB2	1.81	0.61
2:G:9:GLU:O	2:G:11:ARG:N	2.31	0.60
1:F:108:SER:HB2	1:F:121:VAL:HG23	1.83	0.60
1:I:94:ARG:HB2	1:I:178:THR:HG23	1.83	0.60
2:K:96:ASP:HA	2:K:99:LYS:HD3	1.84	0.60
1:E:224:ILE:O	1:F:198:ARG:NH1	2.34	0.59
1:B:36:ASN:HD22	1:B:39:GLU:H	1.48	0.59
1:E:231:LEU:HD21	1:F:204:VAL:HG11	1.83	0.59
1:F:206:THR:O	1:F:208:SER:N	2.35	0.59
2:C:141:VAL:HG22	2:C:253:LEU:HB3	1.83	0.59
1:E:231:LEU:CD2	1:F:204:VAL:HG11	2.33	0.58
1:J:185:PRO:O	1:J:186:VAL:HB	2.01	0.58
2:K:52:GLY:HA3	2:K:120:ASP:HB3	1.85	0.58
2:K:167:LYS:NZ	4:K:401:HOH:O	2.35	0.58
2:K:277:GLU:OE1	2:K:305:ARG:NH2	2.35	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:203:ASN:HB2	1:F:229:GLY:O	2.04	0.58
1:F:151:THR:HG21	1:F:189:MET:HB3	1.86	0.58
1:I:84:GLN:HB3	1:I:186:VAL:HG13	1.85	0.57
2:C:147:VAL:HG11	2:C:270:ALA:N	2.19	0.57
1:E:64:ALA:HB1	1:E:121:VAL:HG11	1.85	0.57
2:G:49:PHE:HB2	2:G:75:ILE:HG23	1.85	0.57
2:G:52:GLY:HA3	2:G:120:ASP:HB3	1.86	0.57
1:F:206:THR:OG1	1:F:207:ALA:N	2.35	0.57
1:A:135:MET:HB3	2:C:259:PRO:HG3	1.87	0.57
1:F:69:ARG:NH1	1:F:197:HIS:HE1	2.03	0.57
2:L:39:SER:HB3	2:L:73:CYS:H	1.69	0.57
1:A:58:PHE:HB3	2:C:5:VAL:HG21	1.86	0.56
2:D:189:LYS:NZ	4:D:402:HOH:O	2.37	0.56
2:L:262:PHE:CZ	3:L:401:CEI:H261	2.39	0.56
1:B:151:THR:OG1	1:B:190:PRO:O	2.21	0.56
2:D:301:SER:O	2:D:305:ARG:HG3	2.06	0.55
1:E:201:LYS:NZ	4:E:303:HOH:O	2.38	0.55
2:H:93:ARG:NH1	2:H:222:GLU:OE2	2.38	0.55
2:K:186:LEU:HA	2:K:286:PHE:HE2	1.72	0.55
2:D:49:PHE:HB2	2:D:75:ILE:HG23	1.89	0.55
1:E:130:ARG:NH1	2:G:256:GLU:OE2	2.34	0.55
1:I:231:LEU:HG	1:J:204:VAL:HG21	1.89	0.55
2:D:281:VAL:HG21	2:D:295:MET:HE2	1.87	0.55
1:F:153:VAL:HG13	1:F:154:THR:H	1.70	0.55
1:E:182:SER:OG	1:E:184:LYS:O	2.18	0.55
2:H:150:ILE:HG13	2:H:150:ILE:O	2.07	0.54
2:C:153:TRP:NE1	2:C:160:GLU:HG3	2.22	0.54
2:D:149:VAL:HG12	2:D:224:PRO:HB2	1.88	0.54
1:F:69:ARG:O	1:F:72:THR:OG1	2.22	0.54
2:G:117:VAL:HG22	2:G:141:VAL:HB	1.88	0.54
2:K:168:SER:OG	2:K:169:GLU:N	2.40	0.54
1:E:88[B]:GLU:OE1	1:E:88[B]:GLU:N	2.37	0.54
2:K:123:ALA:O	2:K:127:PHE:N	2.39	0.54
2:L:52:GLY:HA3	2:L:120:ASP:HB3	1.90	0.54
2:C:168:SER:OG	2:C:169:GLU:N	2.40	0.54
1:E:145:SER:HB3	1:E:197:HIS:HB2	1.90	0.54
2:K:206:PHE:HB3	2:K:213:ARG:HG2	1.89	0.54
1:B:10:GLU:HG3	1:B:38:LEU:HD12	1.90	0.53
2:C:171:GLY:O	2:C:175:VAL:HG12	2.08	0.53
2:K:301:SER:OG	2:K:305:ARG:NH1	2.42	0.53
2:G:218:SER:O	2:G:222:GLU:HG2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:6:LEU:HD22	1:B:6:LEU:H	1.73	0.53
2:C:267:VAL:O	2:C:271:LYS:CB	2.55	0.53
1:E:161:ILE:HG21	1:F:159:GLU:HB3	1.90	0.53
1:J:43:GLU:HG3	1:J:214:HIS:HB2	1.90	0.53
1:A:102:GLY:HA2	1:A:127:ASP:HB2	1.91	0.53
1:B:94:ARG:HB3	1:B:106:VAL:HB	1.89	0.53
2:D:117:VAL:HG22	2:D:141:VAL:HB	1.91	0.53
1:I:41:THR:HG22	1:I:216:THR:HG23	1.91	0.53
1:B:182:SER:OG	1:B:184:LYS:O	2.25	0.52
2:D:287:LEU:HD23	2:D:287:LEU:H	1.73	0.52
1:F:14:THR:OG1	1:F:71:TYR:OH	2.20	0.52
1:A:159:GLU:HG3	1:A:177:ARG:HG3	1.90	0.52
2:C:175:VAL:HA	2:C:180:PHE:H	1.74	0.52
1:J:102:GLY:HA2	1:J:127:ASP:HB2	1.92	0.52
2:D:250:LEU:O	2:D:252:LYS:NZ	2.43	0.52
1:A:3:LEU:HD23	1:A:6:LEU:HD12	1.91	0.52
1:E:5:LYS:NZ	2:K:228:GLY:O	2.43	0.52
1:B:20:GLY:HA2	1:B:123:PHE:O	2.10	0.52
2:L:161:GLU:HA	2:L:164:ALA:HB3	1.92	0.52
2:G:232:ASP:OD1	2:G:232:ASP:N	2.41	0.51
2:H:248:ASP:O	2:H:275:ASN:ND2	2.31	0.51
2:H:157:PRO:O	2:H:159:ILE:HG23	2.11	0.51
2:L:14:MET:HE3	2:L:16:THR:HG22	1.92	0.51
2:L:263:SER:HA	2:L:266:ILE:HG12	1.92	0.51
2:L:263:SER:O	2:L:267:VAL:HG23	2.10	0.51
1:B:144:PRO:HG3	1:B:198:ARG:HG3	1.91	0.51
2:D:122:GLY:HA2	2:D:125:LEU:HD12	1.93	0.51
2:C:49:PHE:O	2:C:76:PRO:HD2	2.11	0.51
1:B:125:ALA:HB1	1:B:128:LEU:HD11	1.91	0.51
2:C:189:LYS:NZ	4:C:406:HOH:O	2.43	0.51
2:C:52:GLY:HA3	2:C:120:ASP:HB3	1.93	0.51
1:I:148:SER:O	1:I:158:GLY:HA2	2.10	0.51
2:C:49:PHE:HB2	2:C:75:ILE:HG12	1.93	0.51
1:B:108:SER:HB2	1:B:121:VAL:HG13	1.91	0.50
1:E:150:TYR:HB3	1:E:192:TYR:HD1	1.75	0.50
2:C:49:PHE:HB2	2:C:75:ILE:HG23	1.93	0.50
2:G:175:VAL:HA	2:G:180:PHE:H	1.76	0.50
2:K:303:VAL:HA	2:K:306:VAL:HG12	1.92	0.50
1:B:135:MET:HE3	2:D:259:PRO:HG3	1.94	0.50
1:F:226:LYS:CE	1:F:232:PRO:O	2.60	0.50
2:K:12:LYS:C	2:K:14:MET:H	2.18	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:80:GLY:HA2	2:H:86:LYS:HD3	1.94	0.50
1:E:159:GLU:OE2	1:F:173:THR:OG1	2.26	0.49
2:D:113:LYS:HB3	2:D:138:LYS:HB2	1.93	0.49
2:C:153:TRP:CZ2	2:C:225:LEU:HD21	2.46	0.49
2:D:49:PHE:HE1	2:D:73:CYS:HB3	1.77	0.49
2:C:269:GLY:C	2:C:271:LYS:H	2.19	0.49
2:H:271:LYS:C	2:H:273:PHE:H	2.20	0.49
1:I:10:GLU:HG3	1:I:38:LEU:HD22	1.94	0.49
1:I:183:LYS:O	1:I:184:LYS:HB3	2.13	0.49
2:L:25:LYS:O	2:L:26:GLN:HG3	2.13	0.49
2:L:184:THR:O	2:L:188:SER:OG	2.22	0.49
1:A:130:ARG:NH1	2:C:256:GLU:OE2	2.42	0.49
2:K:287:LEU:HD23	2:K:287:LEU:H	1.78	0.49
2:L:80:GLY:HA3	2:L:92:TYR:CE1	2.47	0.49
2:H:308:LYS:HD3	2:H:308:LYS:C	2.38	0.49
2:H:119:HIS:CG	2:H:288:GLN:HE22	2.30	0.48
1:I:113:GLU:HG3	1:I:118:ILE:HD13	1.95	0.48
1:I:201:LYS:NZ	4:I:304:HOH:O	2.46	0.48
2:D:247:SER:HB2	2:D:252:LYS:HZ1	1.76	0.48
1:E:20:GLY:HA2	1:E:123:PHE:O	2.14	0.48
1:F:105:ILE:HD13	1:F:107:LYS:HE2	1.95	0.48
2:H:11:ARG:HG3	2:H:195:GLU:CD	2.38	0.48
1:F:4:ALA:O	1:F:6:LEU:HD12	2.13	0.48
2:K:218:SER:O	2:K:222:GLU:HG2	2.12	0.48
2:G:4:LYS:HE2	2:G:190:ILE:O	2.14	0.48
2:G:27:MET:HE3	2:G:27:MET:HB2	1.74	0.48
1:J:159:GLU:HG3	1:J:177:ARG:HG2	1.96	0.48
1:J:147:GLU:HA	1:J:160:CYS:HB3	1.95	0.48
2:C:112:LYS:N	2:C:112:LYS:CD	2.72	0.48
2:C:217:LEU:HD11	2:C:221:ARG:CZ	2.44	0.48
2:C:303:VAL:O	2:C:307:LEU:HG	2.14	0.48
2:H:262:PHE:O	2:H:266:ILE:HG12	2.14	0.48
1:F:36:ASN:HB3	1:F:39:GLU:HB2	1.95	0.47
2:H:233:VAL:O	2:H:237:VAL:HG22	2.14	0.47
2:L:149:VAL:HG11	2:L:234:VAL:HG13	1.95	0.47
2:G:151:GLU:HG2	2:G:226:VAL:HG21	1.96	0.47
2:H:256:GLU:HB3	2:H:280:LYS:HG2	1.96	0.47
2:K:277:GLU:CD	2:K:305:ARG:HH22	2.23	0.47
2:L:49:PHE:O	2:L:76:PRO:HD2	2.14	0.47
1:A:147:GLU:HG3	1:A:160:CYS:HB2	1.96	0.47
2:D:32:SER:OG	2:D:33:PHE:N	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:295:MET:HE3	2:D:295:MET:HB2	1.81	0.47
1:F:110:ILE:HG13	1:F:119:VAL:HG13	1.95	0.47
2:H:175:VAL:HG11	2:H:217:LEU:HB2	1.96	0.47
1:J:97:ARG:HG3	1:J:103:THR:HG22	1.95	0.47
2:K:189:LYS:HE3	2:K:261:PHE:CE1	2.49	0.47
1:F:22:VAL:HG23	1:F:27:PHE:HE1	1.79	0.47
2:K:115:ILE:HG12	2:K:138:LYS:HB3	1.97	0.47
2:K:245:ARG:HE	2:K:272:LYS:HB3	1.79	0.47
3:L:401:CEI:C28	3:L:401:CEI:HN1	2.27	0.47
2:L:104:TRP:CZ2	2:L:108:LEU:HD11	2.50	0.47
1:B:6:LEU:HG	1:B:83:LEU:O	2.15	0.47
1:I:125:ALA:C	1:I:126:LYS:HG3	2.40	0.47
1:B:64:ALA:HB1	1:B:121:VAL:HG11	1.97	0.46
1:B:67:CRO:N2	1:B:67:CRO:HD2	2.31	0.46
1:B:189:MET:HE3	1:B:189:MET:HB3	1.85	0.46
2:D:86:LYS:NZ	4:D:403:HOH:O	2.48	0.46
2:D:149:VAL:HG11	2:D:234:VAL:HG13	1.97	0.46
2:K:32:SER:OG	2:K:33:PHE:N	2.46	0.46
2:K:98:TYR:O	2:K:102:THR:HG23	2.15	0.46
2:L:36:TYR:CE2	2:L:38:ASP:HB2	2.49	0.46
2:H:37:TYR:OH	2:H:68:GLU:OE1	2.27	0.46
2:H:261:PHE:HB2	2:H:285:HIS:HB2	1.97	0.46
1:E:3:LEU:HD23	1:E:5:LYS:H	1.79	0.46
1:E:208:SER:HB2	1:J:5:LYS:O	2.16	0.46
1:F:226:LYS:HE3	1:F:232:PRO:O	2.14	0.46
1:B:6:LEU:HB3	1:B:86:PHE:HB2	1.98	0.46
1:E:204:VAL:HG12	1:E:207:ALA:HB3	1.97	0.46
1:F:155:SER:HB3	1:F:181:LYS:HG2	1.96	0.46
2:G:45:ASN:OD1	2:G:138:LYS:NZ	2.49	0.46
2:H:277:GLU:OE1	2:H:305:ARG:NH1	2.48	0.46
1:I:67:CRO:N2	1:I:67:CRO:HD2	2.31	0.46
1:I:198:ARG:NH1	1:J:226:LYS:HB2	2.31	0.46
2:C:218:SER:O	2:C:222:GLU:HG2	2.15	0.46
1:I:96:ILE:HB	1:I:104:ALA:HB3	1.97	0.46
2:K:175:VAL:HG11	2:K:217:LEU:HB2	1.97	0.46
2:K:299:ILE:O	2:K:303:VAL:HG22	2.16	0.46
1:A:29:MET:HB2	1:A:65:PHE:HZ	1.81	0.46
2:H:256:GLU:HG3	2:H:263:SER:HB3	1.98	0.46
1:B:6:LEU:HD12	1:B:86:PHE:O	2.16	0.46
1:F:203:ASN:O	1:F:207:ALA:HB2	2.16	0.46
2:K:60:TRP:HB2	2:K:75:ILE:HD13	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:VAL:HG12	1:A:154:THR:HG23	1.97	0.46
1:B:110:ILE:HG13	1:B:119:VAL:HG13	1.98	0.46
2:K:115:ILE:HG23	2:K:139:ALA:HB3	1.97	0.46
2:K:251:PRO:HG2	2:K:306:VAL:HG23	1.98	0.46
2:L:42:HIS:H	2:L:72:ARG:HB2	1.80	0.46
1:E:171:HIS:HB2	1:F:150:TYR:CD2	2.52	0.45
2:G:190:ILE:HG12	2:G:286:PHE:HB2	1.98	0.45
2:D:2:THR:HG21	2:D:4:LYS:HZ1	1.81	0.45
1:E:150:TYR:CD2	1:F:171:HIS:HB2	2.50	0.45
1:F:22:VAL:HG23	1:F:27:PHE:CE1	2.52	0.45
2:C:191:MET:HB2	2:C:290:ASP:CG	2.41	0.45
1:E:58:PHE:CE2	1:E:211:VAL:HG11	2.51	0.45
1:F:94:ARG:HB2	1:F:178:THR:HG23	1.97	0.45
1:F:99:GLN:HG2	1:F:172:PHE:CE2	2.51	0.45
2:H:51:HIS:CD2	2:H:56:SER:HA	2.52	0.45
1:J:31:GLY:HA2	1:J:45:LYS:O	2.16	0.45
1:J:148:SER:O	1:J:158:GLY:HA2	2.17	0.45
2:C:172:GLU:O	2:C:176:LEU:HB2	2.16	0.45
1:J:129:ARG:HH12	1:J:130:ARG:HB3	1.82	0.45
1:B:86:PHE:HB3	1:B:87:PRO:HA	1.98	0.45
2:L:53:ASN:HB3	3:L:401:CEI:C29	2.46	0.45
1:A:226:LYS:NZ	1:A:232:PRO:O	2.50	0.45
2:D:4:LYS:HZ3	2:D:188:SER:HA	1.81	0.45
1:E:205:ASP:C	1:E:207:ALA:H	2.25	0.45
2:G:4:LYS:NZ	2:G:188:SER:HA	2.30	0.45
2:L:36:TYR:HE2	2:L:38:ASP:HB2	1.82	0.45
2:G:237:VAL:HA	2:G:240:TYR:CE2	2.52	0.45
2:L:149:VAL:HG12	2:L:224:PRO:HB2	1.98	0.45
1:A:50:LYS:HD2	1:A:51:GLY:H	1.82	0.45
1:B:208:SER:HB3	1:B:210:TYR:CD2	2.52	0.45
2:H:151:GLU:HG2	2:H:226:VAL:HG21	1.99	0.45
1:A:92:TYR:CZ	1:A:108:SER:HB3	2.52	0.45
2:D:186:LEU:HA	2:D:286:PHE:HE2	1.82	0.45
1:F:129:ARG:HA	1:F:129:ARG:HD3	1.83	0.45
2:G:49:PHE:O	2:G:76:PRO:HD2	2.17	0.45
2:G:144:GLU:OE2	2:G:263:SER:OG	2.31	0.45
2:L:64:VAL:HG22	2:L:68:GLU:HG3	1.99	0.45
1:I:36:ASN:OD1	1:I:38:LEU:N	2.48	0.44
1:I:159:GLU:HG3	1:I:177:ARG:HG3	1.99	0.44
1:I:230:SER:HB2	1:J:200:VAL:HB	1.98	0.44
2:K:160:GLU:O	2:K:161:GLU:HB3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:26:GLN:HG3	2:L:35:ASN:OD1	2.17	0.44
1:B:204:VAL:O	1:B:206:THR:N	2.50	0.44
2:D:150:ILE:HB	2:D:225:LEU:HD23	1.98	0.44
1:E:202:THR:O	1:E:204:VAL:N	2.51	0.44
2:H:40:GLU:HA	2:H:72:ARG:HD2	1.98	0.44
1:A:110:ILE:HG13	1:A:119:VAL:HG13	1.99	0.44
1:E:86:PHE:HB3	1:E:87:PRO:HA	2.00	0.44
1:F:147:GLU:HA	1:F:160:CYS:HB3	2.00	0.44
2:H:308:LYS:C	2:H:308:LYS:CD	2.90	0.44
1:I:161:ILE:HG21	1:J:159:GLU:HB3	2.00	0.44
1:I:227:ILE:O	1:I:230:SER:OG	2.33	0.44
1:J:81:TYR:O	1:J:85:SER:OG	2.32	0.44
2:K:237:VAL:HA	2:K:240:TYR:CE2	2.52	0.44
2:L:77:ASP:HB2	2:L:84:SER:OG	2.17	0.44
1:A:77:GLU:OE2	1:A:77:GLU:N	2.47	0.44
1:A:67:CRO:HD2	1:A:67:CRO:N2	2.33	0.44
2:C:269:GLY:C	2:C:271:LYS:N	2.75	0.44
1:J:88:GLU:OE2	1:J:184:LYS:HB2	2.18	0.44
1:J:129:ARG:HH11	1:J:130:ARG:N	2.16	0.44
2:K:12:LYS:C	2:K:14:MET:N	2.76	0.44
2:K:49:PHE:HB2	2:K:75:ILE:HG12	2.00	0.44
2:L:218:SER:O	2:L:222:GLU:HG2	2.17	0.44
1:A:20:GLY:HA2	1:A:123:PHE:O	2.17	0.44
1:A:60:ILE:HG13	1:A:61:VAL:HG13	2.00	0.44
1:A:203:ASN:HB2	1:A:212:VAL:HG11	2.00	0.44
2:C:287:LEU:H	2:C:287:LEU:HD23	1.83	0.44
1:E:3:LEU:CD2	1:E:5:LYS:HG2	2.48	0.44
1:E:202:THR:C	1:E:204:VAL:H	2.26	0.44
1:E:208:SER:O	2:G:193:LYS:NZ	2.49	0.44
2:K:148:ASP:OD1	2:K:241:ASN:CB	2.65	0.44
2:K:234:VAL:HA	2:K:237:VAL:HG12	2.00	0.44
2:K:271:LYS:HG2	2:K:278:PHE:CZ	2.52	0.44
1:A:97:ARG:HD3	1:B:97:ARG:HH12	1.82	0.44
2:D:104:TRP:CE2	2:D:108:LEU:HD11	2.53	0.44
2:G:271:LYS:HA	2:G:278:PHE:HZ	1.81	0.44
2:H:140:ILE:HG13	2:H:250:LEU:HD21	2.00	0.44
1:I:94:ARG:HD3	1:I:176:MET:SD	2.58	0.44
2:L:262:PHE:HZ	3:L:401:CEI:H261	1.81	0.44
1:A:87:PRO:HG3	1:I:228:GLU:HG2	2.00	0.44
2:K:55:THR:HG21	2:K:59:LEU:HD22	1.99	0.44
2:K:186:LEU:HA	2:K:286:PHE:CE2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:149:MET:HB2	1:F:193:HIS:CE1	2.53	0.43
1:I:168:THR:HG23	1:I:170:LYS:H	1.82	0.43
1:J:77:GLU:CD	1:J:77:GLU:H	2.26	0.43
1:J:153:VAL:HG23	1:J:154:THR:N	2.27	0.43
2:K:151:GLU:HG3	2:K:226:VAL:HG21	1.99	0.43
1:B:101:GLY:HA3	1:B:129:ARG:HH21	1.83	0.43
2:C:237:VAL:HA	2:C:240:TYR:CE2	2.53	0.43
2:H:218:SER:O	2:H:222:GLU:HG2	2.17	0.43
2:L:181:PHE:HA	2:L:185:VAL:HB	1.98	0.43
1:A:196:GLN:HG2	1:B:224:ILE:HD12	1.99	0.43
2:C:37:TYR:OH	2:C:68:GLU:OE1	2.30	0.43
2:K:11:ARG:NH2	2:K:197:GLU:OE2	2.50	0.43
2:L:186:LEU:HA	2:L:286:PHE:CE2	2.53	0.43
1:J:67:CRO:N2	1:J:67:CRO:HD1	2.33	0.43
1:A:3:LEU:HD23	1:A:3:LEU:HA	1.77	0.43
1:B:5:LYS:C	1:B:7:GLY:H	2.26	0.43
2:G:138:LYS:HB2	2:G:138:LYS:HE3	1.82	0.43
2:H:303:VAL:O	2:H:307:LEU:HG	2.19	0.43
2:K:295:MET:HE3	2:K:295:MET:HB2	1.74	0.43
2:L:223:ILE:HD12	3:L:401:CEI:H16	2.01	0.43
1:A:16:ILE:HG12	1:A:119:VAL:HB	2.01	0.43
1:B:90:PHE:CE2	1:B:110:ILE:HD13	2.54	0.43
2:G:91:SER:OG	2:G:96:ASP:OD2	2.27	0.43
2:H:271:LYS:HA	2:H:278:PHE:HZ	1.84	0.43
1:J:63:VAL:HG13	1:J:142:MET:HE1	2.01	0.43
1:A:96:ILE:HB	1:A:104:ALA:HB3	2.01	0.43
1:E:3:LEU:HD23	1:E:5:LYS:HG2	2.00	0.43
2:L:163:ILE:HA	2:L:166:ILE:HB	2.00	0.43
2:C:117:VAL:HG22	2:C:141:VAL:HB	1.99	0.43
2:H:32:SER:OG	2:H:33:PHE:N	2.51	0.43
1:I:50:LYS:HA	1:I:50:LYS:HD2	1.75	0.43
2:D:291:ALA:O	2:D:295:MET:HE3	2.19	0.42
1:F:67:CRO:O3	1:F:92:TYR:HE2	2.02	0.42
1:F:92:TYR:CE2	1:F:108:SER:HB3	2.54	0.42
2:H:117:VAL:HG22	2:H:141:VAL:HB	2.01	0.42
2:K:293:ASP:O	2:K:297:LYS:HG3	2.19	0.42
1:E:148:SER:O	1:E:158:GLY:HA2	2.18	0.42
2:L:16:THR:OG1	2:L:19:GLN:HG3	2.19	0.42
1:B:12:MET:HB3	1:B:12:MET:HE3	1.75	0.42
2:H:31:ASP:OD1	2:H:31:ASP:N	2.50	0.42
2:H:287:LEU:HD23	2:H:287:LEU:H	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:155:SER:OG	4:J:301:HOH:O	2.22	0.42
2:D:49:PHE:O	2:D:76:PRO:HD2	2.20	0.42
2:G:48:ILE:HA	2:G:74:ILE:O	2.19	0.42
1:I:55:PRO:HA	2:K:191:MET:HE3	2.01	0.42
1:I:150:TYR:HB3	1:I:192:TYR:HD2	1.85	0.42
2:L:160:GLU:C	2:L:162:ASP:H	2.27	0.42
2:H:240:TYR:O	2:H:244:LEU:HD12	2.19	0.42
2:K:49:PHE:O	2:K:76:PRO:HD2	2.19	0.42
2:G:14:MET:HE3	2:G:14:MET:HB3	1.85	0.42
2:G:104:TRP:CE2	2:G:108:LEU:HD11	2.55	0.42
2:H:64:VAL:HG21	2:H:75:ILE:HD11	2.02	0.42
2:C:214:ARG:N	2:C:215:PRO:HD2	2.34	0.42
2:D:263:SER:HA	2:D:266:ILE:HG12	2.02	0.42
1:E:232:PRO:HD3	1:F:141:GLY:HA3	2.00	0.42
2:G:67:ILE:O	2:G:70:VAL:HG12	2.20	0.42
2:H:247:SER:O	2:H:247:SER:OG	2.30	0.42
1:J:94:ARG:HB3	1:J:106:VAL:HB	2.02	0.42
2:K:191:MET:HE1	2:K:282:LYS:HD2	2.02	0.42
1:E:29:MET:HB2	1:E:65:PHE:HZ	1.84	0.42
2:K:24:CYS:HB3	2:K:35:ASN:HD21	1.84	0.42
1:A:105:ILE:HG13	1:A:124:LYS:HB2	2.02	0.42
1:E:147:GLU:OE2	1:E:180:TYR:OH	2.33	0.42
2:H:189:LYS:HA	2:H:189:LYS:HD3	1.68	0.42
2:L:37:TYR:OH	2:L:68:GLU:OE1	2.27	0.42
1:A:12:MET:HE3	1:A:12:MET:HB3	1.87	0.42
2:D:186:LEU:HA	2:D:286:PHE:CE2	2.55	0.42
1:I:108:SER:HB2	1:I:121:VAL:HG13	2.02	0.42
2:L:57:SER:HB2	2:L:75:ILE:HG22	2.02	0.42
2:L:189:LYS:HE2	3:L:401:CEI:H23	2.01	0.42
2:C:177:GLU:C	2:C:177:GLU:CD	2.87	0.41
1:E:67:CRO:HD2	1:E:67:CRO:N2	2.35	0.41
2:G:27:MET:SD	2:G:107:LEU:HB2	2.60	0.41
2:G:150:ILE:HD12	2:G:223:ILE:HD11	2.03	0.41
2:L:53:ASN:OD1	2:L:216:THR:HA	2.19	0.41
2:D:4:LYS:NZ	2:D:188:SER:HA	2.35	0.41
1:I:92:TYR:CZ	1:I:108:SER:HB3	2.54	0.41
2:K:153:TRP:CZ2	2:K:225:LEU:HD21	2.55	0.41
2:H:171:GLY:HA2	2:H:174:MET:HE3	2.02	0.41
2:L:53:ASN:HB3	3:L:401:CEI:C30	2.51	0.41
2:H:180:PHE:O	2:H:184:THR:HB	2.21	0.41
1:I:15:LYS:HE2	1:I:17:ASN:OD1	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:190:ILE:HG12	2:K:286:PHE:HB2	2.02	0.41
2:L:271:LYS:HG3	2:L:278:PHE:CZ	2.56	0.41
1:A:94:ARG:HB3	1:A:106:VAL:HB	2.03	0.41
1:A:171:HIS:HB2	1:B:150:TYR:CD1	2.56	0.41
1:B:231:LEU:HD13	1:B:231:LEU:HA	1.84	0.41
2:G:104:TRP:CZ2	2:G:108:LEU:HD11	2.56	0.41
1:I:12:MET:HE3	1:I:12:MET:HB3	1.92	0.41
1:J:6:LEU:HD13	1:J:83:LEU:O	2.21	0.41
1:J:90:PHE:CE2	1:J:110:ILE:HD13	2.56	0.41
2:C:177:GLU:OE1	2:C:177:GLU:C	2.62	0.41
1:J:160:CYS:SG	1:J:176:MET:HE2	2.61	0.41
2:K:112:LYS:O	2:K:136:ARG:HD2	2.20	0.41
1:A:85:SER:OG	1:A:90:PHE:HB3	2.21	0.41
2:C:295:MET:HE2	2:C:295:MET:HB2	1.90	0.41
1:F:152:ASN:O	1:F:154:THR:N	2.54	0.41
2:G:53:ASN:HA	2:G:81:MET:HE1	2.01	0.41
2:G:213:ARG:O	2:G:216:THR:OG1	2.36	0.41
1:A:100:ASP:OD2	1:A:132:GLY:HA3	2.21	0.41
2:D:6:TYR:CZ	2:D:193:LYS:HD3	2.56	0.41
2:H:181:PHE:CE2	2:H:220:PRO:HG3	2.55	0.41
1:I:84:GLN:NE2	1:I:187:GLU:HB2	2.36	0.41
2:K:233:VAL:O	2:K:237:VAL:HG12	2.21	0.41
2:D:138:LYS:O	2:D:251:PRO:HD2	2.22	0.40
1:E:4:ALA:HB1	1:E:8:LEU:O	2.21	0.40
1:E:74:TYR:HA	1:E:219:ALA:HB3	2.02	0.40
2:H:271:LYS:C	2:H:273:PHE:N	2.80	0.40
2:C:93:ARG:HD2	2:C:222:GLU:OE2	2.21	0.40
1:F:160:CYS:SG	1:F:176:MET:HE2	2.61	0.40
2:K:115:ILE:CG1	2:K:138:LYS:HB3	2.51	0.40
2:K:152:SER:O	2:K:153:TRP:HB2	2.20	0.40
3:L:401:CEI:HN1	3:L:401:CEI:H28	1.84	0.40
1:A:196:GLN:O	1:A:217:ALA:HA	2.22	0.40
2:C:148:ASP:OD1	2:C:149:VAL:N	2.51	0.40
2:D:253:LEU:HD13	2:D:277:GLU:HB3	2.03	0.40
1:J:129:ARG:HH11	1:J:130:ARG:H	1.69	0.40
2:L:19:GLN:HE21	2:L:19:GLN:HB3	1.61	0.40
2:L:49:PHE:HE1	2:L:73:CYS:HB3	1.86	0.40
2:L:112:LYS:O	2:L:136:ARG:HA	2.21	0.40
1:E:42:GLN:HE22	1:E:67:CRO:CA1	2.32	0.40
2:G:154:ASP:C	2:G:156:TRP:H	2.30	0.40
2:H:237:VAL:HA	2:H:240:TYR:CE2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:149:MET:HE3	1:B:149:MET:HB3	1.98	0.40
2:C:245:ARG:CZ	2:C:272:LYS:HD3	2.52	0.40
2:C:254:PHE:HE2	2:C:256:GLU:HG3	1.87	0.40
2:D:237:VAL:HA	2:D:240:TYR:CE2	2.56	0.40
2:D:301:SER:O	2:D:305:ARG:CG	2.69	0.40
2:H:153:TRP:CZ3	2:H:160:GLU:HG3	2.56	0.40
2:L:79:ILE:HD12	2:L:79:ILE:HA	1.93	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	225/231 (97%)	216 (96%)	6 (3%)	3 (1%)	10	13
1	B	226/231 (98%)	217 (96%)	5 (2%)	4 (2%)	7	8
1	E	225/231 (97%)	215 (96%)	4 (2%)	6 (3%)	4	3
1	F	223/231 (96%)	210 (94%)	7 (3%)	6 (3%)	4	3
1	I	221/231 (96%)	210 (95%)	6 (3%)	5 (2%)	5	5
1	J	225/231 (97%)	213 (95%)	8 (4%)	4 (2%)	7	8
2	C	304/319 (95%)	279 (92%)	24 (8%)	1 (0%)	37	49
2	D	307/319 (96%)	288 (94%)	19 (6%)	0	100	100
2	G	310/319 (97%)	286 (92%)	19 (6%)	5 (2%)	8	10
2	H	306/319 (96%)	280 (92%)	25 (8%)	1 (0%)	37	49
2	K	305/319 (96%)	275 (90%)	30 (10%)	0	100	100
2	L	306/319 (96%)	280 (92%)	23 (8%)	3 (1%)	13	18
All	All	3183/3300 (96%)	2969 (93%)	176 (6%)	38 (1%)	11	14

All (38) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	228	GLU
1	F	5	LYS
1	F	153	VAL
1	F	228	GLU
2	G	309	ASN
2	G	311	GLN
1	J	185	PRO
1	A	205	ASP
1	B	205	ASP
1	B	206	THR
2	C	5	VAL
1	E	153	VAL
1	F	207	ALA
1	I	153	VAL
1	I	191	LEU
1	J	186	VAL
1	J	187	GLU
2	L	169	GLU
1	B	153	VAL
1	E	203	ASN
1	F	113	GLU
2	G	10	GLN
2	G	27	MET
1	I	115	GLY
1	J	153	VAL
2	L	3	SER
2	L	154	ASP
1	A	153	VAL
1	A	228	GLU
1	E	204	VAL
1	F	114	ASP
2	H	159	ILE
1	I	184	LYS
1	I	187	GLU
1	E	208	SER
1	B	208	SER
1	E	191	LEU
2	G	9	GLU

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	197/198 (100%)	197 (100%)	0	100	100
1	B	198/198 (100%)	196 (99%)	2 (1%)	73	85
1	E	197/198 (100%)	196 (100%)	1 (0%)	86	93
1	F	195/198 (98%)	193 (99%)	2 (1%)	73	85
1	I	194/198 (98%)	192 (99%)	2 (1%)	73	85
1	J	197/198 (100%)	195 (99%)	2 (1%)	73	85
2	C	270/282 (96%)	268 (99%)	2 (1%)	81	90
2	D	273/282 (97%)	273 (100%)	0	100	100
2	G	275/282 (98%)	274 (100%)	1 (0%)	89	95
2	H	272/282 (96%)	272 (100%)	0	100	100
2	K	271/282 (96%)	270 (100%)	1 (0%)	89	95
2	L	272/282 (96%)	272 (100%)	0	100	100
All	All	2811/2880 (98%)	2798 (100%)	13 (0%)	86	93

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	116	LYS
1	B	231	LEU
2	C	13	ARG
2	C	112	LYS
1	E	231	LEU
1	F	228	GLU
1	F	230	SER
2	G	106	GLU
1	I	24	ASP
1	I	232	PRO
1	J	107	LYS
1	J	231	LEU
2	K	272	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	GLN
1	A	120	ASN
1	A	136	GLN
1	B	167	GLN
1	B	214	HIS
2	C	119	HIS
2	C	235	GLN
2	D	45	ASN
2	D	97	HIS
1	E	42	GLN
1	E	84	GLN
1	F	42	GLN
1	F	84	GLN
1	F	136	GLN
2	G	133	HIS
2	G	235	GLN
2	H	19	GLN
2	H	235	GLN
2	H	239	ASN
1	I	84	GLN
2	K	35	ASN
2	K	178	ASN
2	K	264	ASN
2	L	19	GLN
2	L	178	ASN
2	L	235	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	CRO	B	67	1	22,22,24	2.56	6 (27%)	27,30,34	3.00	8 (29%)
1	CRO	I	67	1	22,22,24	2.59	6 (27%)	27,30,34	2.93	7 (25%)
1	CRO	F	67	1	22,22,24	2.57	6 (27%)	27,30,34	2.83	7 (25%)
1	CRO	J	67	1	22,22,24	2.62	5 (22%)	27,30,34	3.02	7 (25%)
1	CRO	E	67	1	22,22,24	2.60	5 (22%)	27,30,34	3.19	7 (25%)
1	CRO	A	67	1	22,22,24	2.59	6 (27%)	27,30,34	2.98	7 (25%)

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
1	CRO	B	67	1	-	3/9/29/32	0/2/2/2
1	CRO	I	67	1	-	3/9/29/32	0/2/2/2
1	CRO	F	67	1	-	4/9/29/32	0/2/2/2
1	CRO	J	67	1	-	5/9/29/32	0/2/2/2
1	CRO	E	67	1	-	4/9/29/32	0/2/2/2
1	CRO	A	67	1	-	4/9/29/32	0/2/2/2

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	67	CRO	CA2-C2	-9.61	1.39	1.48
1	E	67	CRO	CA2-C2	-9.56	1.39	1.48
1	A	67	CRO	CA2-C2	-9.45	1.39	1.48
1	I	67	CRO	CA2-C2	-9.42	1.39	1.48
1	F	67	CRO	CA2-C2	-9.35	1.39	1.48
1	B	67	CRO	CA2-C2	-9.34	1.39	1.48
1	F	67	CRO	C1-N2	4.07	1.38	1.32
1	J	67	CRO	C1-N2	3.95	1.38	1.32
1	I	67	CRO	C1-N2	3.91	1.38	1.32
1	A	67	CRO	C1-N2	3.89	1.37	1.32
1	E	67	CRO	C1-N2	3.81	1.37	1.32
1	B	67	CRO	C1-N2	3.76	1.37	1.32
1	F	67	CRO	CE2-CZ	3.15	1.44	1.38
1	E	67	CRO	CE2-CZ	3.14	1.44	1.38
1	A	67	CRO	CE2-CZ	3.12	1.44	1.38
1	I	67	CRO	CE2-CZ	3.09	1.44	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	67	CRO	CE2-CZ	3.08	1.44	1.38
1	B	67	CRO	CE2-CZ	3.08	1.44	1.38
1	B	67	CRO	CG2-CB2	2.40	1.51	1.46
1	I	67	CRO	CG2-CB2	2.39	1.51	1.46
1	E	67	CRO	CG2-CB2	2.33	1.51	1.46
1	F	67	CRO	CG2-CB2	2.32	1.51	1.46
1	B	67	CRO	CD2-CG2	2.32	1.43	1.39
1	A	67	CRO	CG2-CB2	2.31	1.51	1.46
1	A	67	CRO	CD2-CG2	2.30	1.43	1.39
1	J	67	CRO	CG2-CB2	2.27	1.51	1.46
1	E	67	CRO	CD2-CG2	2.24	1.43	1.39
1	J	67	CRO	CD2-CG2	2.22	1.43	1.39
1	I	67	CRO	CD2-CG2	2.21	1.43	1.39
1	F	67	CRO	CD2-CG2	2.17	1.43	1.39
1	F	67	CRO	C1-N3	-2.15	1.33	1.37
1	I	67	CRO	C1-N3	-2.13	1.33	1.37
1	A	67	CRO	C1-N3	-2.10	1.33	1.37
1	B	67	CRO	C1-N3	-2.03	1.33	1.37

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	67	CRO	O2-C2-CA2	-10.97	124.80	130.96
1	J	67	CRO	O2-C2-CA2	-9.65	125.54	130.96
1	A	67	CRO	O2-C2-CA2	-9.57	125.58	130.96
1	J	67	CRO	CA2-C2-N3	9.47	107.85	103.37
1	B	67	CRO	CA2-C2-N3	9.45	107.84	103.37
1	E	67	CRO	CA2-C2-N3	9.45	107.84	103.37
1	I	67	CRO	CA2-C2-N3	9.44	107.84	103.37
1	A	67	CRO	CA2-C2-N3	9.41	107.82	103.37
1	F	67	CRO	CA2-C2-N3	9.17	107.71	103.37
1	B	67	CRO	O2-C2-CA2	-8.98	125.92	130.96
1	I	67	CRO	O2-C2-CA2	-8.94	125.94	130.96
1	F	67	CRO	O2-C2-CA2	-8.82	126.01	130.96
1	J	67	CRO	N3-C1-N2	-4.16	108.57	111.45
1	B	67	CRO	CG2-CB2-CA2	-3.94	125.12	129.94
1	I	67	CRO	N3-C1-N2	-3.92	108.74	111.45
1	B	67	CRO	N3-C1-N2	-3.89	108.76	111.45
1	B	67	CRO	CA2-N2-C1	3.78	108.56	105.77
1	A	67	CRO	N3-C1-N2	-3.67	108.91	111.45
1	E	67	CRO	CG2-CB2-CA2	-3.66	125.46	129.94
1	J	67	CRO	CA2-N2-C1	3.60	108.42	105.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	67	CRO	N3-C1-N2	-3.57	108.98	111.45
1	I	67	CRO	CA2-N2-C1	3.50	108.36	105.77
1	F	67	CRO	N3-C1-N2	-3.50	109.03	111.45
1	E	67	CRO	CA2-N2-C1	3.45	108.32	105.77
1	A	67	CRO	CA2-N2-C1	3.36	108.25	105.77
1	A	67	CRO	CG2-CB2-CA2	-3.23	125.99	129.94
1	I	67	CRO	CG2-CB2-CA2	-3.15	126.08	129.94
1	J	67	CRO	CG2-CB2-CA2	-3.10	126.14	129.94
1	F	67	CRO	CG2-CB2-CA2	-3.00	126.27	129.94
1	B	67	CRO	C2-CA2-N2	-2.97	106.86	108.93
1	F	67	CRO	CA2-N2-C1	2.92	107.92	105.77
1	I	67	CRO	C2-CA2-N2	-2.88	106.91	108.93
1	J	67	CRO	C2-CA2-N2	-2.72	107.03	108.93
1	A	67	CRO	C2-CA2-N2	-2.69	107.05	108.93
1	E	67	CRO	C2-CA2-N2	-2.56	107.14	108.93
1	B	67	CRO	CB2-CA2-C2	2.51	125.28	122.28
1	F	67	CRO	C2-CA2-N2	-2.43	107.23	108.93
1	I	67	CRO	O3-C3-CA3	-2.36	119.26	126.39
1	F	67	CRO	O3-C3-CA3	-2.31	119.42	126.39
1	J	67	CRO	O3-C3-CA3	-2.24	119.61	126.39
1	A	67	CRO	O3-C3-CA3	-2.17	119.84	126.39
1	E	67	CRO	O3-C3-CA3	-2.09	120.07	126.39
1	B	67	CRO	O3-C3-CA3	-2.01	120.31	126.39

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	67	CRO	C3-CA3-N3-C2
1	B	67	CRO	C3-CA3-N3-C2
1	B	67	CRO	C2-CA2-CB2-CG2
1	E	67	CRO	C3-CA3-N3-C1
1	E	67	CRO	C3-CA3-N3-C2
1	F	67	CRO	C3-CA3-N3-C1
1	F	67	CRO	C3-CA3-N3-C2
1	F	67	CRO	C2-CA2-CB2-CG2
1	I	67	CRO	C3-CA3-N3-C2
1	J	67	CRO	N1-CA1-CB1-OG1
1	J	67	CRO	C1-CA1-CB1-OG1
1	J	67	CRO	C3-CA3-N3-C2
1	A	67	CRO	N2-CA2-CB2-CG2
1	B	67	CRO	N2-CA2-CB2-CG2

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Mol	Chain	Res	Type	Atoms
1	E	67	CRO	N2-CA2-CB2-CG2
1	F	67	CRO	N2-CA2-CB2-CG2
1	I	67	CRO	C1-CA1-CB1-OG1
1	A	67	CRO	C2-CA2-CB2-CG2
1	E	67	CRO	C2-CA2-CB2-CG2
1	I	67	CRO	N2-CA2-CB2-CG2
1	J	67	CRO	N2-CA2-CB2-CG2
1	A	67	CRO	C3-CA3-N3-C1
1	J	67	CRO	C3-CA3-N3-C1

There are no ring outliers.

6 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	67	CRO	1	0
1	I	67	CRO	1	0
1	F	67	CRO	1	0
1	J	67	CRO	1	0
1	E	67	CRO	5	0
1	A	67	CRO	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	CEI	L	401	-	31,34,34	1.54	4 (12%)	38,46,46	1.07	2 (5%)

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
3	CEI	L	401	-	-	0/12/16/16	0/4/4/4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	401	CEI	C2-N1	4.73	1.46	1.35
3	L	401	CEI	C9-N1	3.69	1.46	1.39
3	L	401	CEI	C6-N7	-3.50	1.31	1.37
3	L	401	CEI	O33-C2	-2.18	1.18	1.23

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	401	CEI	N1-C9-N4	2.68	124.33	117.45
3	L	401	CEI	C8-C9-N4	-2.48	119.94	122.25

There are no chirality outliers.

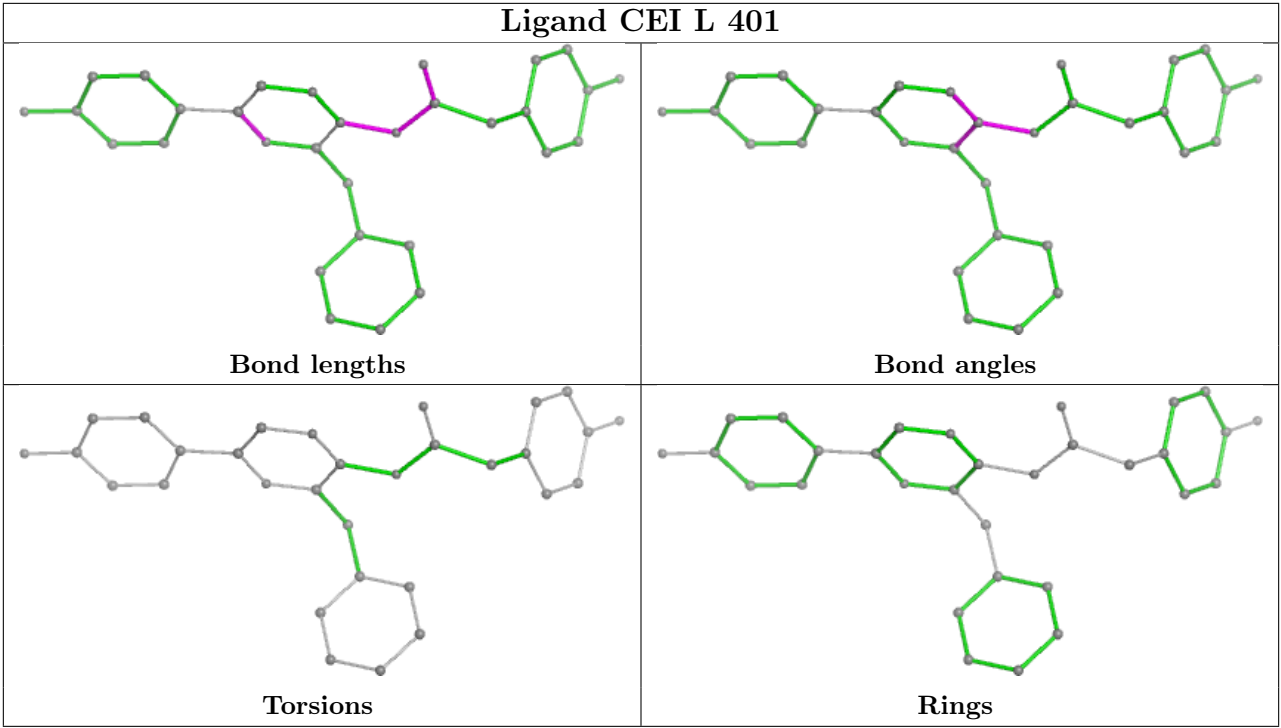
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	L	401	CEI	8	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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	F	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	F	67:CRO	C3	68:ASN	N	1.61

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	229/231 (99%)	0.07	2 (0%) 81 80	32, 51, 79, 138	0
1	B	230/231 (99%)	0.09	3 (1%) 74 74	34, 50, 77, 111	0
1	E	228/231 (98%)	0.22	10 (4%) 39 40	33, 54, 90, 125	1 (0%)
1	F	227/231 (98%)	0.27	9 (3%) 43 43	36, 56, 94, 160	0
1	I	225/231 (97%)	0.22	5 (2%) 62 62	35, 54, 79, 124	0
1	J	229/231 (99%)	0.10	3 (1%) 74 74	34, 50, 83, 150	0
2	C	306/319 (95%)	0.12	9 (2%) 54 53	36, 51, 78, 98	0
2	D	309/319 (96%)	0.11	1 (0%) 90 89	36, 53, 75, 118	0
2	G	312/319 (97%)	0.09	4 (1%) 74 74	33, 52, 75, 112	0
2	H	308/319 (96%)	0.42	6 (1%) 66 65	41, 64, 87, 108	0
2	K	307/319 (96%)	0.35	7 (2%) 61 60	37, 64, 88, 121	0
2	L	308/319 (96%)	0.15	4 (1%) 74 74	37, 56, 80, 107	0
All	All	3218/3300 (97%)	0.19	63 (1%) 64 64	32, 55, 83, 160	1 (0%)

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	270	ALA	8.6
1	A	204	VAL	6.9
2	G	313	GLY	3.8
1	E	3	LEU	3.8
1	I	6	LEU	3.7
1	E	207	ALA	3.6
1	E	209	GLY	3.5
2	H	147	VAL	3.4
1	F	4	ALA	3.4
1	I	207	ALA	3.3
1	E	204	VAL	3.3

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Mol	Chain	Res	Type	RSRZ
2	K	12	LYS	3.3
1	E	4	ALA	3.3
1	J	204	VAL	3.0
1	F	204	VAL	2.9
1	B	207	ALA	2.9
2	C	8	PRO	2.9
1	F	206	THR	2.9
1	E	40	GLY	2.9
2	C	269	GLY	2.9
2	K	26	GLN	2.7
1	I	183	LYS	2.7
1	I	210	TYR	2.7
2	K	303	VAL	2.7
2	K	307	LEU	2.7
2	K	250	LEU	2.6
1	B	161	ILE	2.6
2	C	5	VAL	2.6
2	K	139	ALA	2.6
2	L	168	SER	2.6
1	A	38	LEU	2.6
1	F	6	LEU	2.5
1	E	86	PHE	2.4
1	J	206	THR	2.4
1	E	218	ILE	2.4
1	F	229	GLY	2.4
1	E	6	LEU	2.4
1	B	6	LEU	2.3
2	C	177	GLU	2.3
2	H	139	ALA	2.3
1	F	87	PRO	2.3
2	H	255	ILE	2.3
2	C	14	MET	2.2
2	H	186	LEU	2.2
2	G	310	GLU	2.2
2	D	46	ALA	2.2
1	I	206	THR	2.2
1	J	211	VAL	2.2
2	G	26	GLN	2.2
1	F	191	LEU	2.2
1	E	210	TYR	2.2
2	C	10	GLN	2.2
2	K	308	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	F	7	GLY	2.1
2	L	41	LYS	2.1
1	F	207	ALA	2.1
2	G	114	ILE	2.1
2	H	114	ILE	2.1
2	L	163	ILE	2.1
2	H	266	ILE	2.1
2	C	4	LYS	2.0
2	L	112	LYS	2.0
2	C	309	ASN	2.0

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

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
1	CRO	F	67	21/23	0.87	0.12	41,44,48,52	0
1	CRO	J	67	21/23	0.91	0.10	26,48,64,76	0
1	CRO	I	67	21/23	0.92	0.09	37,48,61,63	0
1	CRO	E	67	21/23	0.92	0.10	44,53,64,71	0
1	CRO	A	67	21/23	0.93	0.09	41,49,54,57	0
1	CRO	B	67	21/23	0.93	0.09	34,48,59,62	0

6.3 Carbohydrates ⓘ

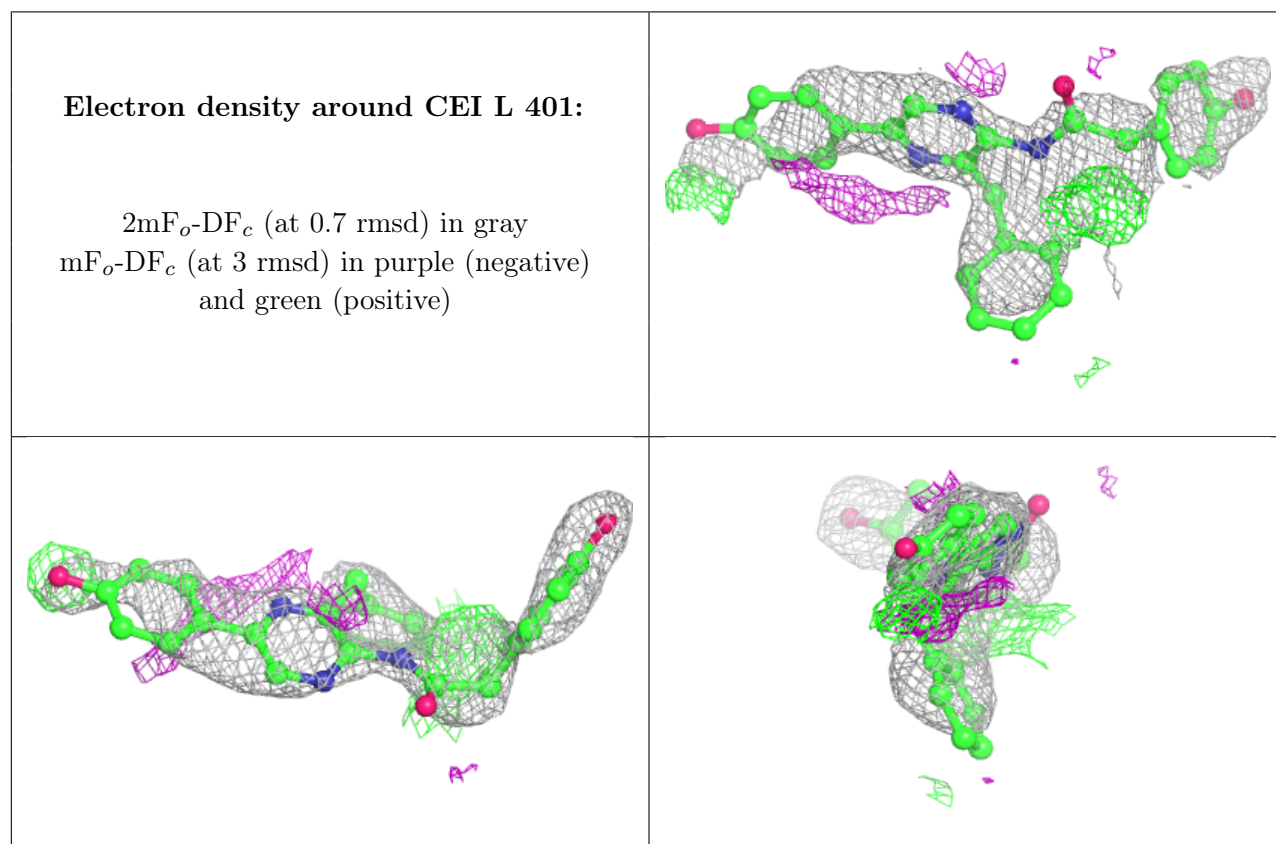
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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	CEI	L	401	31/31	0.79	0.23	64,96,108,111	0

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



6.5 Other polymers ⓘ

There are no such residues in this entry.