



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:44 PM UTC

PDB ID : 2I4N / pdb_00002i4n
Title : Rhodopseudomonas palustris prolyl-tRNA synthetase in complex with CysAMS
Authors : Crepin, T.; Yaremchuk, A.; Tukalo, M.; Cusack, S.
Deposited on : 2006-08-22
Resolution : 2.85 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

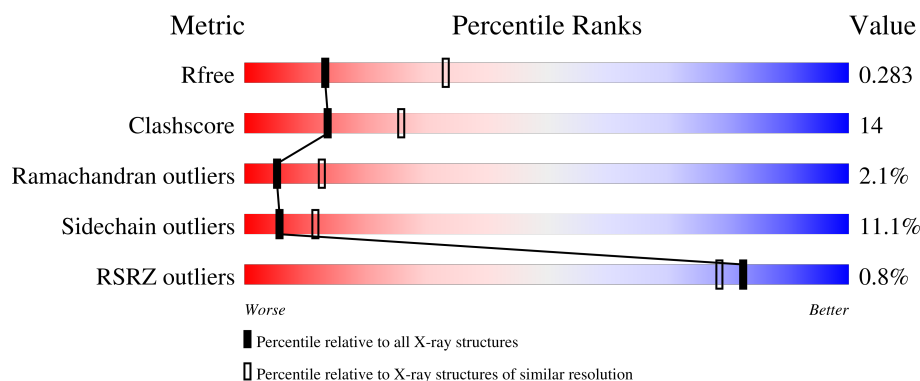
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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	458	% 66% 24% 6% .
1	B	458	66% 26% . .
1	C	458	% 56% 34% 7% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10672 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 Proline-tRNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	441	Total	C	N	O	S	0	0	0
			3495	2211	620	650	14			
1	B	442	Total	C	N	O	S	0	0	0
			3506	2217	624	651	14			
1	C	442	Total	C	N	O	S	0	0	0
			3506	2217	624	651	14			

There are 60 discrepancies between the modelled and reference sequences:

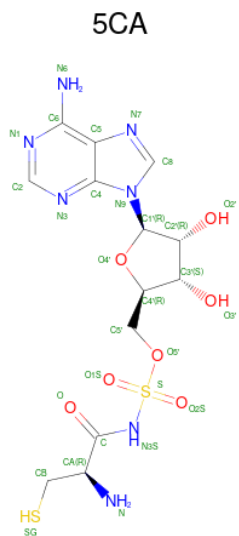
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q6N5P6
A	-18	GLY	-	expression tag	UNP Q6N5P6
A	-17	SER	-	expression tag	UNP Q6N5P6
A	-16	SER	-	expression tag	UNP Q6N5P6
A	-15	HIS	-	expression tag	UNP Q6N5P6
A	-14	HIS	-	expression tag	UNP Q6N5P6
A	-13	HIS	-	expression tag	UNP Q6N5P6
A	-12	HIS	-	expression tag	UNP Q6N5P6
A	-11	HIS	-	expression tag	UNP Q6N5P6
A	-10	HIS	-	expression tag	UNP Q6N5P6
A	-9	SER	-	expression tag	UNP Q6N5P6
A	-8	SER	-	expression tag	UNP Q6N5P6
A	-7	GLY	-	expression tag	UNP Q6N5P6
A	-6	LEU	-	expression tag	UNP Q6N5P6
A	-5	VAL	-	expression tag	UNP Q6N5P6
A	-4	PRO	-	expression tag	UNP Q6N5P6
A	-3	ARG	-	expression tag	UNP Q6N5P6
A	-2	GLY	-	expression tag	UNP Q6N5P6
A	-1	SER	-	expression tag	UNP Q6N5P6
A	0	HIS	-	expression tag	UNP Q6N5P6
B	-19	MET	-	expression tag	UNP Q6N5P6
B	-18	GLY	-	expression tag	UNP Q6N5P6
B	-17	SER	-	expression tag	UNP Q6N5P6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-16	SER	-	expression tag	UNP Q6N5P6
B	-15	HIS	-	expression tag	UNP Q6N5P6
B	-14	HIS	-	expression tag	UNP Q6N5P6
B	-13	HIS	-	expression tag	UNP Q6N5P6
B	-12	HIS	-	expression tag	UNP Q6N5P6
B	-11	HIS	-	expression tag	UNP Q6N5P6
B	-10	HIS	-	expression tag	UNP Q6N5P6
B	-9	SER	-	expression tag	UNP Q6N5P6
B	-8	SER	-	expression tag	UNP Q6N5P6
B	-7	GLY	-	expression tag	UNP Q6N5P6
B	-6	LEU	-	expression tag	UNP Q6N5P6
B	-5	VAL	-	expression tag	UNP Q6N5P6
B	-4	PRO	-	expression tag	UNP Q6N5P6
B	-3	ARG	-	expression tag	UNP Q6N5P6
B	-2	GLY	-	expression tag	UNP Q6N5P6
B	-1	SER	-	expression tag	UNP Q6N5P6
B	0	HIS	-	expression tag	UNP Q6N5P6
C	-19	MET	-	expression tag	UNP Q6N5P6
C	-18	GLY	-	expression tag	UNP Q6N5P6
C	-17	SER	-	expression tag	UNP Q6N5P6
C	-16	SER	-	expression tag	UNP Q6N5P6
C	-15	HIS	-	expression tag	UNP Q6N5P6
C	-14	HIS	-	expression tag	UNP Q6N5P6
C	-13	HIS	-	expression tag	UNP Q6N5P6
C	-12	HIS	-	expression tag	UNP Q6N5P6
C	-11	HIS	-	expression tag	UNP Q6N5P6
C	-10	HIS	-	expression tag	UNP Q6N5P6
C	-9	SER	-	expression tag	UNP Q6N5P6
C	-8	SER	-	expression tag	UNP Q6N5P6
C	-7	GLY	-	expression tag	UNP Q6N5P6
C	-6	LEU	-	expression tag	UNP Q6N5P6
C	-5	VAL	-	expression tag	UNP Q6N5P6
C	-4	PRO	-	expression tag	UNP Q6N5P6
C	-3	ARG	-	expression tag	UNP Q6N5P6
C	-2	GLY	-	expression tag	UNP Q6N5P6
C	-1	SER	-	expression tag	UNP Q6N5P6
C	0	HIS	-	expression tag	UNP Q6N5P6

- Molecule 2 is 5'-O-(N-(L-CYSTEINYL)-SULFAMOYL)ADENOSINE (CCD ID: 5CA) (formula: C₁₃H₁₉N₇O₇S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 29	C 13	N 7	O 7	S 2	0	0
2	B	1	Total 29	C 13	N 7	O 7	S 2	0	0
2	C	1	Total 29	C 13	N 7	O 7	S 2	0	0

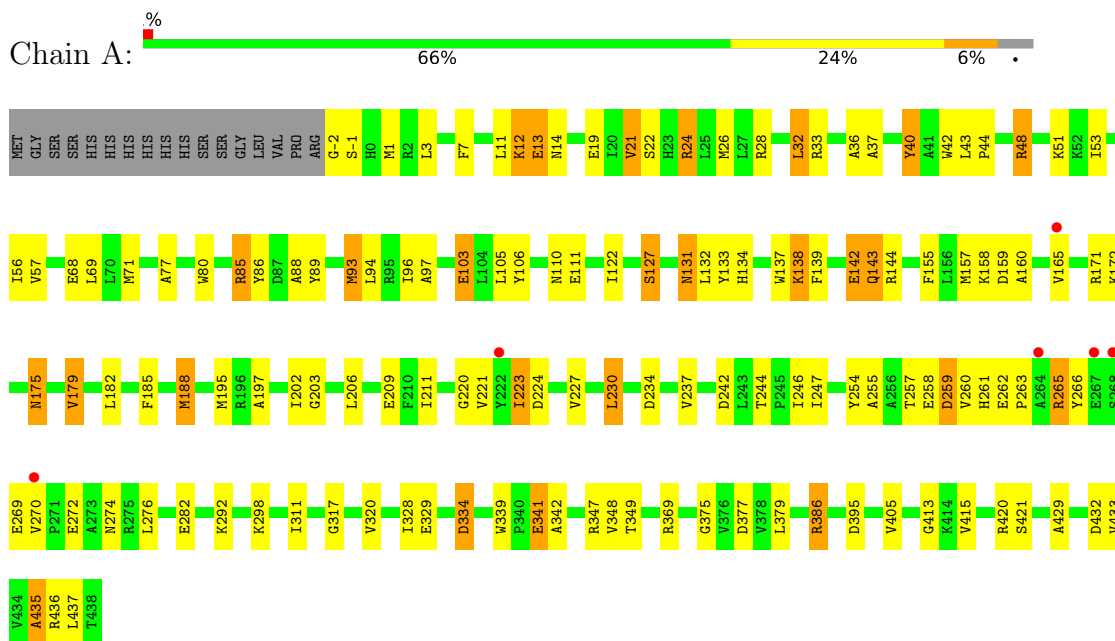
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	28	Total O 28 28	0	0
3	B	30	Total O 30 30	0	0
3	C	20	Total O 20 20	0	0

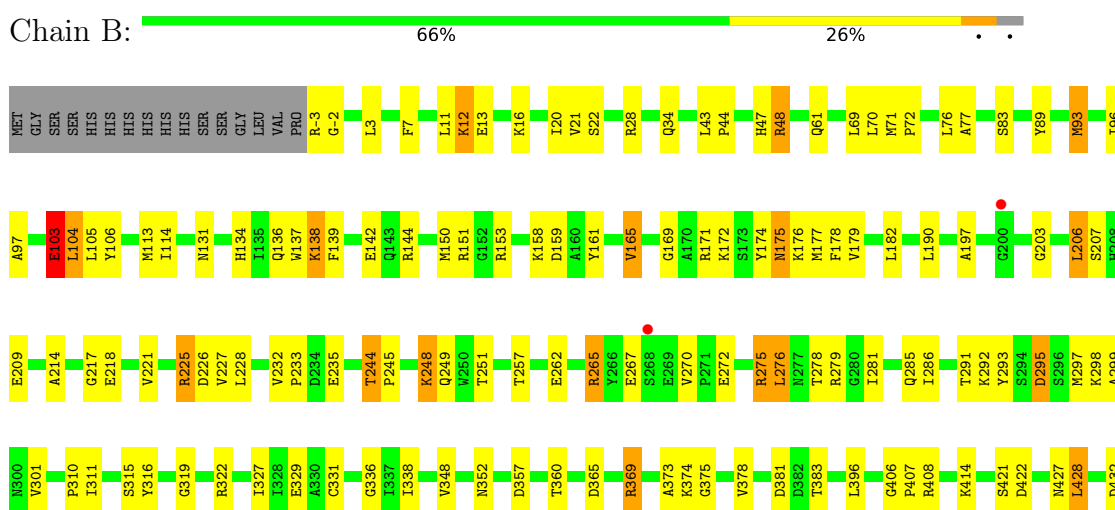
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Proline-tRNA ligase

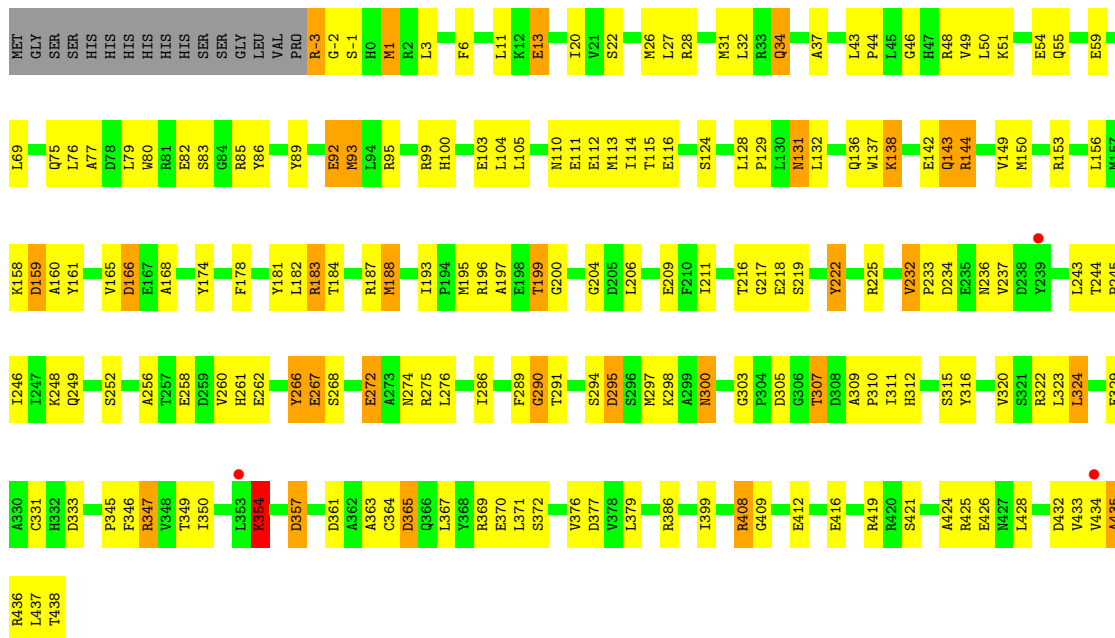


• Molecule 1: Proline-tRNA ligase



V433
V434
A435
R436
L437
T438

• Molecule 1: Proline-tRNA ligase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	217.90Å 107.47Å 110.66Å 90.00° 120.43° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.85 95.42 – 2.85	Depositor EDS
% Data completeness (in resolution range)	94.6 ((Not available)-2.85) 94.6 (95.42-2.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.01	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.20 (at 2.68Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.211 , 0.277 (Not available) , 0.283	Depositor DCC
R_{free} test set	2903 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	63.7	Xtriage
Anisotropy	0.144	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 43.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.047 for -h+k-l,-l,-k 0.034 for -h-k-l,l,k 0.064 for -h-2*l,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10672	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.69% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 5CA

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	0.79	0/3570	1.07	12/4828 (0.2%)
1	B	0.78	0/3581	1.02	6/4842 (0.1%)
1	C	0.77	0/3581	1.06	13/4842 (0.3%)
All	All	0.78	0/10732	1.05	31/14512 (0.2%)

There are no bond length outliers.

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	168	ALA	N-CA-C	-7.74	102.47	112.23
1	A	230	LEU	CA-C-N	7.23	127.23	119.78
1	A	230	LEU	C-N-CA	7.23	127.23	119.78
1	C	437	LEU	N-CA-C	-6.76	105.66	114.04
1	B	378	VAL	N-CA-C	6.64	117.41	108.11
1	A	85	ARG	N-CA-C	-6.64	104.32	112.88
1	A	40	TYR	N-CA-C	6.51	119.79	109.50
1	A	272	GLU	N-CA-C	6.45	118.89	111.02
1	B	406	GLY	CA-C-N	5.93	127.26	119.84
1	B	406	GLY	C-N-CA	5.93	127.26	119.84
1	A	19	GLU	N-CA-C	5.91	117.59	111.03
1	C	1	MET	N-CA-C	5.87	117.95	108.32
1	B	275	ARG	N-CA-C	5.58	119.83	113.19
1	C	200	GLY	CA-C-N	5.57	126.80	119.84
1	C	200	GLY	C-N-CA	5.57	126.80	119.84
1	A	246	ILE	N-CA-C	-5.47	107.11	111.81
1	A	88	ALA	N-CA-C	5.46	117.98	111.71
1	B	83	SER	N-CA-C	-5.37	106.57	113.23
1	C	232	VAL	N-CA-C	5.36	112.85	107.55
1	C	217	GLY	N-CA-C	5.33	118.06	111.93
1	A	276	LEU	N-CA-C	5.32	116.25	107.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	43	LEU	CA-C-N	5.29	126.46	119.84
1	C	43	LEU	C-N-CA	5.29	126.46	119.84
1	C	232	VAL	CA-C-N	5.27	126.43	119.84
1	C	232	VAL	C-N-CA	5.27	126.43	119.84
1	B	165	VAL	CB-CA-C	-5.18	105.07	112.22
1	C	331	CYS	N-CA-C	5.14	117.01	107.60
1	A	342	ALA	N-CA-C	5.09	117.56	111.71
1	A	179	VAL	CB-CA-C	-5.09	105.27	112.14
1	C	256	ALA	N-CA-C	5.03	116.58	108.32
1	A	110	ASN	N-CA-C	5.03	118.49	111.56

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3495	0	3466	100	0
1	B	3506	0	3479	103	0
1	C	3506	0	3479	106	0
2	A	29	0	19	1	0
2	B	29	0	19	1	0
2	C	29	0	19	1	0
3	A	28	0	0	3	0
3	B	30	0	0	0	0
3	C	20	0	0	1	0
All	All	10672	0	10481	305	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 14.

All (305) 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:C:-3:ARG:HG3	1:C:-3:ARG:HH11	1.10	1.13

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:436:ARG:HG2	1:B:436:ARG:HH11	0.95	1.09
1:B:48:ARG:HG2	1:B:48:ARG:HH11	1.18	1.00
1:C:188:MET:HE1	1:C:323:LEU:HD13	1.43	0.98
1:B:72:PRO:HD2	1:B:113:MET:HE1	1.42	0.97
1:A:-2:GLY:HA3	1:A:375:GLY:HA2	1.48	0.96
1:B:175:ASN:N	1:B:175:ASN:HD22	1.63	0.96
1:A:12:LYS:HD3	1:A:12:LYS:H	1.29	0.95
1:A:12:LYS:HD3	1:A:12:LYS:N	1.84	0.92
1:C:188:MET:CE	1:C:323:LEU:HD13	2.00	0.92
1:B:436:ARG:HG2	1:B:436:ARG:NH1	1.74	0.92
1:B:436:ARG:HH11	1:B:436:ARG:CG	1.84	0.91
1:C:-3:ARG:HG3	1:C:-3:ARG:NH1	1.86	0.88
1:C:199:THR:HG21	1:C:204:GLY:O	1.72	0.88
1:B:175:ASN:HD22	1:B:175:ASN:H	1.18	0.86
1:A:259:ASP:O	1:A:261:HIS:N	2.08	0.86
1:C:272:GLU:HA	1:C:275:ARG:HD2	1.59	0.85
1:A:386:ARG:H	1:A:386:ARG:HH11	1.22	0.85
1:C:-3:ARG:HH11	1:C:-3:ARG:CG	1.89	0.83
1:B:48:ARG:HH11	1:B:48:ARG:CG	1.91	0.83
1:C:158:LYS:HD2	1:C:158:LYS:C	2.06	0.81
1:B:171:ARG:O	1:B:175:ASN:ND2	2.13	0.80
1:B:22:SER:HB3	1:B:329:GLU:HG2	1.63	0.79
1:A:48:ARG:CG	1:A:48:ARG:HH11	1.96	0.78
1:B:175:ASN:H	1:B:175:ASN:ND2	1.80	0.78
1:C:76:LEU:HD23	1:C:104:LEU:HD23	1.63	0.78
1:B:293:TYR:O	1:B:297:MET:HB2	1.85	0.77
1:B:267:GLU:HA	1:B:275:ARG:HH12	1.49	0.76
1:A:435:ALA:O	1:A:437:LEU:N	2.19	0.76
1:A:259:ASP:C	1:A:261:HIS:H	1.93	0.76
1:A:32:LEU:HD11	1:A:40:TYR:HB3	1.69	0.73
1:A:48:ARG:HH11	1:A:48:ARG:HG2	1.53	0.73
1:A:14:ASN:HD22	1:A:24:ARG:HG3	1.53	0.73
1:C:138:LYS:NZ	1:C:159:ASP:OD1	2.21	0.73
1:C:305:ASP:OD1	1:C:307:THR:HG23	1.88	0.72
1:A:265:ARG:NH2	1:A:269:GLU:OE1	2.21	0.71
1:B:225:ARG:HG3	1:B:225:ARG:HH11	1.53	0.71
1:A:22:SER:HB3	1:A:329:GLU:HG2	1.72	0.70
1:A:-2:GLY:CA	1:A:375:GLY:HA2	2.22	0.70
1:C:131:ASN:C	1:C:131:ASN:HD22	2.01	0.68
1:B:71:MET:HB3	1:B:113:MET:CE	2.25	0.66
1:B:175:ASN:N	1:B:175:ASN:ND2	2.35	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:GLY:O	1:A:258:GLU:HB2	1.96	0.66
1:A:386:ARG:H	1:A:386:ARG:NH1	1.93	0.66
1:B:72:PRO:HD2	1:B:113:MET:CE	2.23	0.65
1:A:11:LEU:HB2	1:A:28:ARG:HA	1.79	0.64
1:A:234:ASP:O	1:A:237:VAL:HG23	1.96	0.64
1:A:26:MET:HE2	1:A:32:LEU:HD23	1.80	0.64
1:B:93:MET:HE1	1:B:105:LEU:HD22	1.78	0.64
1:A:175:ASN:O	1:A:179:VAL:HG23	1.98	0.64
1:A:68:GLU:HB2	1:A:133:TYR:CZ	2.34	0.63
1:C:51:LYS:HD2	1:C:54:GLU:OE2	1.99	0.63
1:B:76:LEU:HD23	1:B:104:LEU:HD22	1.81	0.62
1:B:71:MET:HB3	1:B:113:MET:HE3	1.81	0.62
1:A:138:LYS:HD3	1:A:157:MET:HB3	1.81	0.62
1:B:77:ALA:HB3	1:B:103:GLU:HB2	1.81	0.62
1:C:408:ARG:HG3	1:C:409:GLY:H	1.64	0.62
1:A:224:ASP:HB2	1:A:274:ASN:HB3	1.81	0.61
1:C:432:ASP:O	1:C:435:ALA:HB2	1.99	0.61
1:B:272:GLU:HA	1:B:275:ARG:HE	1.64	0.61
1:C:294:SER:HB3	1:C:311:ILE:HG13	1.82	0.61
1:B:267:GLU:HA	1:B:275:ARG:NH1	2.15	0.61
1:B:207:SER:HA	1:B:286:ILE:O	2.00	0.61
1:A:89:TYR:HB2	1:A:93:MET:HG3	1.83	0.61
1:A:-1:SER:O	1:A:377:ASP:HA	2.00	0.61
1:B:138:LYS:HE3	1:B:159:ASP:OD1	2.01	0.61
1:C:188:MET:HE1	1:C:323:LEU:CD1	2.24	0.60
1:C:433:VAL:C	1:C:435:ALA:H	2.08	0.60
1:C:158:LYS:HD2	1:C:158:LYS:O	2.00	0.60
1:C:354:LYS:HE2	1:C:386:ARG:HH12	1.66	0.60
1:C:188:MET:HE2	1:C:323:LEU:HD22	1.84	0.60
1:A:257:THR:O	1:A:261:HIS:HB2	2.01	0.59
1:A:21:VAL:HA	1:A:24:ARG:HD2	1.83	0.59
1:B:226:ASP:C	1:B:228:LEU:H	2.11	0.59
1:A:223:ILE:HG22	1:A:255:ALA:HA	1.84	0.59
1:C:11:LEU:HB2	1:C:28:ARG:HA	1.85	0.59
1:B:48:ARG:HG2	1:B:48:ARG:NH1	1.97	0.59
1:B:365:ASP:O	1:B:369:ARG:HD3	2.03	0.59
1:C:77:ALA:HB3	1:C:103:GLU:HB3	1.84	0.59
1:A:106:TYR:CD2	1:A:139:PHE:HB2	2.38	0.58
1:A:263:PRO:C	1:A:265:ARG:H	2.11	0.58
1:B:267:GLU:CA	1:B:275:ARG:HH12	2.15	0.58
1:A:311:ILE:HD12	1:A:311:ILE:C	2.28	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:363:ALA:O	1:C:367:LEU:HB2	2.02	0.58
1:A:131:ASN:C	1:A:131:ASN:HD22	2.12	0.57
1:A:224:ASP:HB3	1:A:270:VAL:HG11	1.85	0.57
1:A:158:LYS:O	1:A:317:GLY:HA2	2.04	0.57
1:C:99:ARG:HG3	1:C:100:HIS:ND1	2.20	0.57
1:C:76:LEU:HD23	1:C:104:LEU:CD2	2.34	0.57
1:C:44:PRO:O	1:C:48:ARG:HG3	2.05	0.57
1:C:93:MET:HE1	1:C:105:LEU:HB2	1.86	0.57
1:C:196:ARG:HB3	3:C:451:HOH:O	2.05	0.56
1:C:361:ASP:O	1:C:365:ASP:HB2	2.06	0.56
1:B:72:PRO:CD	1:B:113:MET:HE1	2.27	0.56
1:A:395:ASP:HA	1:A:420:ARG:HE	1.71	0.56
1:C:349:THR:HA	1:C:379:LEU:O	2.05	0.56
1:A:197:ALA:HB2	1:A:209:GLU:HB2	1.87	0.56
1:B:61:GLN:HE22	1:B:177:MET:HE1	1.70	0.56
1:A:265:ARG:O	1:A:265:ARG:HD3	2.06	0.55
1:C:346:PHE:O	1:C:399:ILE:HG23	2.06	0.55
1:B:48:ARG:CG	1:B:48:ARG:NH1	2.58	0.55
1:A:1:MET:O	1:A:379:LEU:HD12	2.06	0.55
1:A:421:SER:HB3	3:A:453:HOH:O	2.07	0.55
1:C:83:SER:HB3	1:C:297:MET:HE3	1.88	0.55
1:C:225:ARG:HG3	1:C:274:ASN:O	2.07	0.55
1:C:-1:SER:O	1:C:377:ASP:HA	2.06	0.54
1:B:217:GLY:O	1:B:279:ARG:NH1	2.35	0.54
1:A:137:TRP:CZ2	1:B:70:LEU:HD21	2.41	0.54
1:B:244:THR:HB	1:B:245:PRO:HD3	1.88	0.54
1:B:144:ARG:NH1	1:B:144:ARG:HB2	2.22	0.54
1:B:262:GLU:HB3	1:B:265:ARG:HB2	1.89	0.54
1:C:49:VAL:HG22	1:C:345:PRO:HD3	1.90	0.54
1:A:185:PHE:HA	1:A:188:MET:HG3	1.90	0.54
1:A:195:MET:HE3	1:A:211:ILE:HB	1.89	0.53
1:A:13:GLU:H	1:A:13:GLU:CD	2.16	0.53
1:C:294:SER:CB	1:C:311:ILE:HG13	2.39	0.53
1:C:234:ASP:O	1:C:237:VAL:HG23	2.09	0.53
1:B:373:ALA:C	1:B:375:GLY:H	2.16	0.53
1:C:1:MET:HG2	1:C:6:PHE:CD1	2.43	0.53
1:C:93:MET:SD	1:C:105:LEU:HD22	2.48	0.53
1:A:7:PHE:HB3	1:A:44:PRO:HG2	1.91	0.53
1:B:61:GLN:NE2	1:B:177:MET:HE1	2.23	0.52
1:A:12:LYS:N	1:A:12:LYS:CD	2.66	0.52
1:B:139:PHE:HA	1:B:153:ARG:O	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:ILE:HG22	1:A:320:VAL:HG13	1.91	0.52
1:A:429:ALA:O	1:A:432:ASP:HB2	2.08	0.52
1:B:11:LEU:HB2	1:B:28:ARG:HA	1.91	0.52
1:B:214:ALA:HB2	1:B:281:ILE:HD12	1.91	0.52
1:C:75:GLN:HG2	1:C:113:MET:CE	2.40	0.52
1:C:75:GLN:HG2	1:C:113:MET:HE3	1.91	0.52
1:C:13:GLU:CD	1:C:13:GLU:H	2.18	0.51
1:C:89:TYR:O	1:C:92:GLU:HB2	2.09	0.51
1:C:291:THR:O	1:C:295:ASP:HB2	2.10	0.51
1:B:179:VAL:HG13	1:B:233:PRO:HG2	1.93	0.51
1:A:421:SER:CB	3:A:453:HOH:O	2.59	0.51
1:C:246:ILE:HA	1:C:249:GLN:HB3	1.92	0.50
1:C:289:PHE:O	1:C:290:GLY:C	2.54	0.50
1:B:432:ASP:O	1:B:433:VAL:C	2.54	0.50
1:A:203:GLY:HA3	1:A:292:LYS:HE3	1.92	0.50
1:B:311:ILE:C	1:B:311:ILE:HD12	2.36	0.50
1:A:334:ASP:OD2	1:A:334:ASP:N	2.43	0.50
1:C:244:THR:O	1:C:248:LYS:HG3	2.12	0.50
1:A:32:LEU:HD12	1:A:33:ARG:N	2.26	0.50
2:B:439:5CA:O	2:B:439:5CA:O1S	2.28	0.50
1:C:110:ASN:O	1:C:114:ILE:HG22	2.12	0.50
1:C:22:SER:HB3	1:C:329:GLU:HG2	1.94	0.50
1:C:85:ARG:O	1:C:86:TYR:C	2.53	0.50
1:C:142:GLU:H	1:C:153:ARG:HG2	1.77	0.50
1:A:-1:SER:HB3	1:A:377:ASP:OD1	2.12	0.49
1:A:386:ARG:HH11	1:A:386:ARG:N	1.99	0.49
1:B:144:ARG:HB2	1:B:144:ARG:HH11	1.77	0.49
1:C:34:GLN:OE1	1:C:37:ALA:HA	2.11	0.49
1:C:183:ARG:HD2	1:C:233:PRO:O	2.12	0.49
1:C:233:PRO:HB3	1:C:237:VAL:HG21	1.93	0.49
1:C:323:LEU:O	1:C:324:LEU:C	2.53	0.49
1:B:150:MET:HE2	1:B:329:GLU:HG3	1.94	0.49
1:C:89:TYR:HB2	1:C:93:MET:HG2	1.95	0.49
1:C:183:ARG:O	1:C:184:THR:C	2.56	0.49
1:C:193:ILE:HB	1:C:195:MET:HE3	1.94	0.49
1:B:275:ARG:O	1:B:276:LEU:C	2.54	0.48
1:C:50:LEU:HD13	1:C:324:LEU:CD1	2.43	0.48
1:B:291:THR:O	1:B:295:ASP:HB2	2.14	0.48
1:C:-2:GLY:HA3	1:C:372:SER:O	2.14	0.48
1:B:114:ILE:HG21	1:B:161:TYR:CG	2.48	0.48
1:B:209:GLU:HB3	1:B:285:GLN:HG3	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:299:ALA:O	1:B:310:PRO:HA	2.13	0.48
1:B:315:SER:C	1:B:316:TYR:CG	2.91	0.48
1:A:48:ARG:HG2	1:A:48:ARG:NH1	2.23	0.48
1:B:131:ASN:ND2	1:B:177:MET:HE3	2.28	0.48
1:B:131:ASN:HD21	1:B:177:MET:HE3	1.78	0.48
1:C:128:LEU:HB3	1:C:129:PRO:HA	1.94	0.48
1:B:174:TYR:CZ	1:B:206:LEU:HD22	2.48	0.48
1:B:336:GLY:HA3	1:B:396:LEU:HA	1.95	0.48
1:C:26:MET:HE1	1:C:149:VAL:HG22	1.94	0.48
1:C:131:ASN:HD21	1:C:160:ALA:HB1	1.79	0.48
1:C:416:GLU:OE2	1:C:425:ARG:HD3	2.14	0.48
1:C:174:TYR:CE2	1:C:286:ILE:HD12	2.48	0.47
1:B:244:THR:HG22	1:B:248:LYS:HD3	1.95	0.47
1:A:122:ILE:HD12	1:A:127:SER:OG	2.13	0.47
1:B:76:LEU:HD23	1:B:104:LEU:CD2	2.43	0.47
1:B:217:GLY:HA3	1:B:279:ARG:HB3	1.95	0.47
1:A:242:ASP:OD1	1:A:244:THR:HG22	2.15	0.47
1:A:311:ILE:HD12	1:A:311:ILE:O	2.15	0.47
1:B:174:TYR:OH	1:B:206:LEU:HD22	2.15	0.47
1:B:221:VAL:HG22	1:B:278:THR:O	2.15	0.47
1:A:131:ASN:HD21	1:A:160:ALA:HB1	1.79	0.47
1:C:131:ASN:C	1:C:131:ASN:ND2	2.72	0.47
1:C:159:ASP:OD1	1:C:159:ASP:N	2.47	0.47
1:A:96:ILE:HG22	1:A:97:ALA:N	2.30	0.47
1:A:142:GLU:O	1:A:143:GLN:C	2.57	0.47
1:B:96:ILE:HG22	1:B:97:ALA:N	2.29	0.47
1:C:31:MET:HE3	1:C:46:GLY:HA2	1.96	0.47
1:C:136:GLN:HG3	1:C:137:TRP:N	2.30	0.47
1:C:50:LEU:CD1	1:C:320:VAL:HG12	2.46	0.46
1:A:341:GLU:OE1	1:A:347:ARG:NH1	2.48	0.46
1:C:20:ILE:HD12	1:C:329:GLU:OE2	2.15	0.46
1:B:182:LEU:HB3	1:B:232:VAL:HG13	1.97	0.46
1:A:254:TYR:CE2	1:A:266:TYR:HD2	2.34	0.46
1:C:178:PHE:C	1:C:178:PHE:CD2	2.93	0.46
1:C:261:HIS:C	1:C:261:HIS:CD2	2.94	0.46
1:C:137:TRP:CE2	1:C:156:LEU:HD13	2.51	0.45
1:A:103:GLU:HG3	3:A:450:HOH:O	2.16	0.45
1:A:234:ASP:C	1:A:234:ASP:OD1	2.60	0.45
1:B:175:ASN:O	1:B:176:LYS:C	2.58	0.45
1:C:244:THR:N	1:C:245:PRO:HD2	2.31	0.45
1:A:413:GLY:O	1:A:429:ALA:HA	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:ARG:HH11	1:A:48:ARG:HG3	1.80	0.45
1:A:13:GLU:CD	1:A:13:GLU:N	2.75	0.45
1:C:346:PHE:HA	1:C:377:ASP:HB3	1.99	0.45
1:B:357:ASP:HB3	1:B:360:THR:OG1	2.17	0.45
1:A:22:SER:HA	1:A:328:ILE:HG21	1.98	0.45
1:B:190:LEU:HD21	1:B:327:ILE:HA	1.99	0.45
1:C:77:ALA:O	1:C:80:TRP:N	2.35	0.45
1:C:82:GLU:HB3	1:C:297:MET:HE2	1.98	0.45
1:A:111:GLU:OE2	2:A:439:5CA:N	2.50	0.45
1:B:158:LYS:C	1:B:158:LYS:HD2	2.41	0.45
1:A:137:TRP:HZ2	1:B:70:LEU:HD21	1.82	0.45
1:B:197:ALA:HB2	1:B:209:GLU:HG2	1.98	0.44
1:A:80:TRP:HB3	1:A:86:TYR:HA	1.99	0.44
1:A:94:LEU:HD12	1:A:94:LEU:HA	1.73	0.44
1:A:175:ASN:OD1	1:A:247:ILE:HG13	2.17	0.44
1:B:151:ARG:HD3	1:B:218:GLU:OE2	2.17	0.44
1:B:226:ASP:C	1:B:228:LEU:N	2.75	0.44
1:A:77:ALA:HB2	1:A:103:GLU:HB2	1.99	0.44
1:A:43:LEU:O	1:A:44:PRO:C	2.58	0.44
1:B:267:GLU:HG2	1:B:275:ARG:HH12	1.82	0.44
1:B:136:GLN:HG3	1:B:137:TRP:N	2.33	0.44
1:C:333:ASP:C	1:C:333:ASP:OD1	2.60	0.44
1:A:227:VAL:O	1:A:230:LEU:HD12	2.18	0.44
1:A:259:ASP:C	1:A:261:HIS:N	2.60	0.44
1:B:150:MET:O	1:B:322:ARG:NH1	2.51	0.44
1:B:169:GLY:O	1:B:172:LYS:HB2	2.17	0.44
1:A:68:GLU:HG2	1:B:47:HIS:CD2	2.54	0.43
1:C:110:ASN:HB2	1:C:161:TYR:OH	2.18	0.43
1:C:354:LYS:O	1:C:357:ASP:HB3	2.19	0.43
1:B:297:MET:C	1:B:298:LYS:HG2	2.43	0.43
1:B:7:PHE:HB3	1:B:44:PRO:HG2	2.00	0.43
1:C:114:ILE:HD11	1:C:132:LEU:HD12	2.01	0.43
1:C:303:GLY:C	1:C:305:ASP:H	2.26	0.43
1:A:42:TRP:CD1	1:A:42:TRP:N	2.86	0.43
1:C:357:ASP:OD2	1:C:357:ASP:C	2.62	0.43
1:B:69:LEU:O	1:B:134:HIS:HB3	2.19	0.43
1:B:225:ARG:HG3	1:B:225:ARG:NH1	2.28	0.43
1:A:71:MET:HB2	1:A:134:HIS:CE1	2.53	0.43
1:A:348:VAL:HG22	1:A:349:THR:N	2.34	0.43
1:B:209:GLU:OE2	1:B:257:THR:HG21	2.18	0.42
1:C:55:GLN:NE2	1:C:59:GLU:OE2	2.45	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:419:ARG:HB3	1:C:424:ALA:HB3	2.00	0.42
1:A:262:GLU:HA	1:A:263:PRO:HD3	1.82	0.42
1:B:225:ARG:HH11	1:B:225:ARG:CG	2.28	0.42
1:B:381:ASP:OD1	1:B:383:THR:OG1	2.30	0.42
1:A:254:TYR:HB3	1:A:270:VAL:HG21	2.02	0.42
1:B:105:LEU:C	1:B:105:LEU:HD12	2.44	0.42
1:C:27:LEU:HD23	1:C:32:LEU:HB3	2.01	0.42
1:C:244:THR:N	1:C:245:PRO:CD	2.82	0.42
1:A:405:VAL:HG22	1:A:415:VAL:HG11	2.01	0.42
1:C:112:GLU:O	1:C:115:THR:HB	2.19	0.42
1:A:339:TRP:HA	1:A:339:TRP:CE3	2.55	0.42
1:B:203:GLY:HA3	1:B:292:LYS:HD2	2.01	0.42
1:C:93:MET:HB3	1:C:93:MET:HE2	1.79	0.42
1:B:-3:ARG:N	1:B:-3:ARG:HE	2.18	0.42
1:B:428:LEU:HD12	1:B:428:LEU:C	2.45	0.42
1:A:137:TRP:HZ2	1:B:70:LEU:CD2	2.33	0.41
1:A:137:TRP:HA	1:A:155:PHE:O	2.20	0.41
1:A:171:ARG:O	1:A:175:ASN:ND2	2.53	0.41
1:B:178:PHE:CD2	1:B:178:PHE:C	2.98	0.41
1:C:347:ARG:H	1:C:377:ASP:HB2	1.85	0.41
1:B:331:CYS:HB3	1:B:338:ILE:O	2.20	0.41
1:C:315:SER:C	1:C:316:TYR:CG	2.99	0.41
1:A:57:VAL:HG11	1:A:158:LYS:HG3	2.02	0.41
1:C:222:TYR:CD2	1:C:222:TYR:N	2.88	0.41
1:A:36:ALA:O	1:A:37:ALA:C	2.62	0.41
1:B:436:ARG:NH1	1:B:436:ARG:CG	2.54	0.41
1:C:350:ILE:HD11	1:C:371:LEU:HD12	2.03	0.41
1:B:175:ASN:O	1:B:179:VAL:HG23	2.20	0.41
1:C:197:ALA:HB2	1:C:209:GLU:HB2	2.03	0.41
1:B:434:VAL:O	1:B:438:THR:HG23	2.21	0.41
1:A:22:SER:HA	1:A:328:ILE:CG2	2.51	0.41
1:A:-1:SER:CB	1:A:377:ASP:OD1	2.69	0.41
1:A:85:ARG:NH2	1:A:111:GLU:HB2	2.35	0.41
1:A:131:ASN:HD22	1:A:132:LEU:N	2.19	0.41
1:B:43:LEU:HB3	1:B:44:PRO:CD	2.51	0.41
1:C:131:ASN:HD22	1:C:132:LEU:N	2.19	0.41
1:C:181:TYR:O	1:C:182:LEU:C	2.62	0.41
1:A:105:LEU:C	1:A:105:LEU:HD12	2.46	0.41
1:A:57:VAL:CG1	1:A:158:LYS:HG3	2.50	0.40
1:B:89:TYR:HB2	1:B:93:MET:HG3	2.01	0.40
1:B:106:TYR:CD2	1:B:139:PHE:HB2	2.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:311:ILE:HD12	1:B:311:ILE:O	2.21	0.40
1:C:93:MET:CE	1:C:105:LEU:HD22	2.52	0.40
1:A:32:LEU:HD13	1:A:32:LEU:HA	1.80	0.40
1:A:48:ARG:CG	1:A:48:ARG:NH1	2.66	0.40
1:A:137:TRP:CZ2	1:B:70:LEU:CD2	3.04	0.40
1:A:182:LEU:HA	1:A:182:LEU:HD23	1.91	0.40
1:B:297:MET:HB3	1:B:297:MET:HE2	1.87	0.40
1:C:166:ASP:O	1:C:312:HIS:CG	2.75	0.40
1:C:266:TYR:CE1	1:C:275:ARG:HG2	2.57	0.40
1:C:300:ASN:HB3	1:C:309:ALA:C	2.46	0.40
1:C:364:CYS:O	1:C:367:LEU:N	2.54	0.40
1:B:352:ASN:OD1	1:B:352:ASN:C	2.64	0.40
1:B:414:LYS:HE2	1:B:414:LYS:HB3	1.89	0.40
1:B:428:LEU:C	1:B:428:LEU:CD1	2.95	0.40
1:C:111:GLU:OE2	2:C:439:5CA:N	2.55	0.40
1:C:143:GLN:HB3	1:C:144:ARG:HG2	2.03	0.40
1:C:311:ILE:HD12	1:C:311:ILE:C	2.46	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	439/458 (96%)	393 (90%)	41 (9%)	5 (1%)	11	24
1	B	440/458 (96%)	392 (89%)	36 (8%)	12 (3%)	4	9
1	C	440/458 (96%)	374 (85%)	55 (12%)	11 (2%)	4	10
All	All	1319/1374 (96%)	1159 (88%)	132 (10%)	28 (2%)	5	13

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	260	VAL
1	A	436	ARG
1	B	12	LYS
1	B	103	GLU
1	C	408	ARG
1	A	175	ASN
1	A	435	ALA
1	B	-2	GLY
1	B	227	VAL
1	B	276	LEU
1	B	421	SER
1	C	267	GLU
1	C	354	LYS
1	C	434	VAL
1	C	435	ALA
1	B	142	GLU
1	B	235	GLU
1	C	266	TYR
1	C	252	SER
1	A	143	GLN
1	B	295	ASP
1	B	374	LYS
1	C	166	ASP
1	C	258	GLU
1	C	290	GLY
1	B	407	PRO
1	C	310	PRO
1	B	319	GLY

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	361/376 (96%)	327 (91%)	34 (9%)	8	18
1	B	362/376 (96%)	332 (92%)	30 (8%)	10	22
1	C	362/376 (96%)	306 (84%)	56 (16%)	2	5

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1085/1128 (96%)	965 (89%)	120 (11%)	6 12

All (120) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	12	LYS
1	A	13	GLU
1	A	21	VAL
1	A	24	ARG
1	A	32	LEU
1	A	48	ARG
1	A	51	LYS
1	A	56	ILE
1	A	69	LEU
1	A	93	MET
1	A	103	GLU
1	A	127	SER
1	A	131	ASN
1	A	138	LYS
1	A	142	GLU
1	A	144	ARG
1	A	159	ASP
1	A	165	VAL
1	A	172	LYS
1	A	188	MET
1	A	202	ILE
1	A	206	LEU
1	A	221	VAL
1	A	223	ILE
1	A	259	ASP
1	A	265	ARG
1	A	282	GLU
1	A	298	LYS
1	A	334	ASP
1	A	341	GLU
1	A	369	ARG
1	A	386	ARG
1	A	433	VAL
1	B	3	LEU
1	B	12	LYS
1	B	13	GLU

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Mol	Chain	Res	Type
1	B	16	LYS
1	B	20	ILE
1	B	21	VAL
1	B	34	GLN
1	B	48	ARG
1	B	93	MET
1	B	103	GLU
1	B	104	LEU
1	B	138	LYS
1	B	165	VAL
1	B	175	ASN
1	B	206	LEU
1	B	225	ARG
1	B	244	THR
1	B	248	LYS
1	B	249	GLN
1	B	251	THR
1	B	265	ARG
1	B	270	VAL
1	B	301	VAL
1	B	348	VAL
1	B	369	ARG
1	B	408	ARG
1	B	422	ASP
1	B	427	ASN
1	B	428	LEU
1	B	436	ARG
1	C	-3	ARG
1	C	3	LEU
1	C	13	GLU
1	C	34	GLN
1	C	69	LEU
1	C	79	LEU
1	C	92	GLU
1	C	93	MET
1	C	95	ARG
1	C	116	GLU
1	C	124	SER
1	C	131	ASN
1	C	138	LYS
1	C	143	GLN
1	C	144	ARG

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Mol	Chain	Res	Type
1	C	150	MET
1	C	159	ASP
1	C	165	VAL
1	C	183	ARG
1	C	187	ARG
1	C	188	MET
1	C	199	THR
1	C	206	LEU
1	C	211	ILE
1	C	216	THR
1	C	218	GLU
1	C	219	SER
1	C	222	TYR
1	C	232	VAL
1	C	236	ASN
1	C	243	LEU
1	C	260	VAL
1	C	262	GLU
1	C	267	GLU
1	C	268	SER
1	C	272	GLU
1	C	276	LEU
1	C	295	ASP
1	C	298	LYS
1	C	300	ASN
1	C	307	THR
1	C	322	ARG
1	C	324	LEU
1	C	347	ARG
1	C	354	LYS
1	C	357	ASP
1	C	365	ASP
1	C	369	ARG
1	C	370	GLU
1	C	376	VAL
1	C	412	GLU
1	C	421	SER
1	C	426	GLU
1	C	428	LEU
1	C	436	ARG
1	C	438	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	ASN
1	A	131	ASN
1	A	277	ASN
1	A	285	GLN
1	A	332	HIS
1	A	335	ASN
1	B	47	HIS
1	B	61	GLN
1	B	175	ASN
1	B	236	ASN
1	B	285	GLN
1	B	300	ASN
1	C	0	HIS
1	C	23	HIS
1	C	131	ASN
1	C	143	GLN
1	C	229	ASN
1	C	236	ASN
1	C	249	GLN
1	C	332	HIS
1	C	352	ASN
1	C	355	GLN
1	C	402	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 ligands 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
2	5CA	A	439	-	31,31,31	2.52	8 (25%)	42,46,46	2.20	13 (30%)
2	5CA	B	439	-	31,31,31	2.17	6 (19%)	42,46,46	2.35	13 (30%)
2	5CA	C	439	-	31,31,31	2.29	6 (19%)	42,46,46	2.00	10 (23%)

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	5CA	A	439	-	-	3/20/37/37	0/3/3/3
2	5CA	B	439	-	-	9/20/37/37	0/3/3/3
2	5CA	C	439	-	-	3/20/37/37	0/3/3/3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	439	5CA	O2S-S	8.49	1.49	1.42
2	A	439	5CA	O1S-S	8.21	1.49	1.42
2	C	439	5CA	O2S-S	7.74	1.49	1.42
2	B	439	5CA	O2S-S	7.51	1.49	1.42
2	C	439	5CA	O1S-S	7.16	1.48	1.42
2	B	439	5CA	O1S-S	6.03	1.47	1.42
2	B	439	5CA	S-N3S	-3.92	1.53	1.59
2	C	439	5CA	S-N3S	-3.35	1.54	1.59
2	B	439	5CA	C5-C6	3.29	1.50	1.41
2	A	439	5CA	CB-CA	-3.09	1.49	1.53
2	A	439	5CA	C5-C6	3.00	1.49	1.41
2	C	439	5CA	C8-N7	2.77	1.37	1.31
2	C	439	5CA	C5-N7	-2.59	1.34	1.39
2	A	439	5CA	CB-SG	-2.27	1.76	1.81
2	B	439	5CA	C8-N7	2.23	1.36	1.31
2	A	439	5CA	C8-N7	2.23	1.36	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	439	5CA	S-N3S	-2.19	1.56	1.59
2	A	439	5CA	C5-N7	-2.12	1.35	1.39
2	C	439	5CA	C5-C6	2.10	1.46	1.41
2	B	439	5CA	C4-N9	-2.05	1.33	1.37

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	439	5CA	O2S-S-O1S	-6.67	110.86	120.85
2	B	439	5CA	O2S-S-O1S	-6.42	111.24	120.85
2	A	439	5CA	C5-C4-N3	-5.59	119.01	126.72
2	A	439	5CA	CA-CB-SG	-5.48	102.69	114.40
2	B	439	5CA	O5'-S-N3S	5.46	119.82	105.69
2	C	439	5CA	C5-C4-N3	-5.06	119.75	126.72
2	B	439	5CA	CA-CB-SG	-4.61	104.53	114.40
2	B	439	5CA	N3-C2-N1	-4.53	121.72	128.58
2	B	439	5CA	C5-C4-N3	-4.51	120.51	126.72
2	A	439	5CA	N3-C2-N1	-4.41	121.91	128.58
2	A	439	5CA	C2-N3-C4	4.33	122.40	111.83
2	A	439	5CA	O2S-S-O1S	-4.06	114.78	120.85
2	B	439	5CA	C2-N3-C4	4.04	121.69	111.83
2	B	439	5CA	N9-C8-N7	-3.86	108.46	113.94
2	A	439	5CA	N3-C4-N9	3.74	133.53	127.17
2	A	439	5CA	N9-C8-N7	-3.43	109.07	113.94
2	C	439	5CA	N3-C2-N1	-3.41	123.43	128.58
2	C	439	5CA	C2-N3-C4	3.40	120.13	111.83
2	B	439	5CA	C4-N9-C8	3.22	109.12	105.74
2	A	439	5CA	C2'-C1'-N9	-3.00	105.84	113.30
2	C	439	5CA	N9-C8-N7	-2.98	109.70	113.94
2	C	439	5CA	N3-C4-N9	2.97	132.22	127.17
2	B	439	5CA	C5'-O5'-S	2.96	123.17	116.97
2	B	439	5CA	C4-C5-N7	-2.85	107.33	110.58
2	B	439	5CA	N3-C4-N9	2.78	131.90	127.17
2	C	439	5CA	CA-CB-SG	-2.75	108.52	114.40
2	A	439	5CA	C4-C5-N7	-2.58	107.63	110.58
2	B	439	5CA	C5-N7-C8	2.57	107.50	103.45
2	A	439	5CA	CB-CA-C	-2.56	103.98	109.52
2	A	439	5CA	C5-N7-C8	2.55	107.46	103.45
2	A	439	5CA	C4-N9-C8	2.53	108.39	105.74
2	C	439	5CA	O5'-S-N3S	2.48	112.11	105.69
2	A	439	5CA	O5'-S-N3S	2.30	111.64	105.69
2	B	439	5CA	C6-C5-N7	2.13	136.19	132.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	439	5CA	C5-C4-N9	2.03	108.02	105.81
2	C	439	5CA	C5-N7-C8	2.02	106.62	103.45

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	439	5CA	N-CA-CB-SG
2	A	439	5CA	C-CA-CB-SG
2	B	439	5CA	N-CA-CB-SG
2	B	439	5CA	C-CA-CB-SG
2	B	439	5CA	C5'-O5'-S-N3S
2	B	439	5CA	C5'-O5'-S-O2S
2	C	439	5CA	N-CA-CB-SG
2	C	439	5CA	C-CA-CB-SG
2	B	439	5CA	O-C-CA-CB
2	B	439	5CA	C-N3S-S-O2S
2	B	439	5CA	C5'-O5'-S-O1S
2	B	439	5CA	O-C-CA-N
2	B	439	5CA	N3S-C-CA-CB
2	C	439	5CA	O-C-CA-N
2	A	439	5CA	C5'-O5'-S-N3S

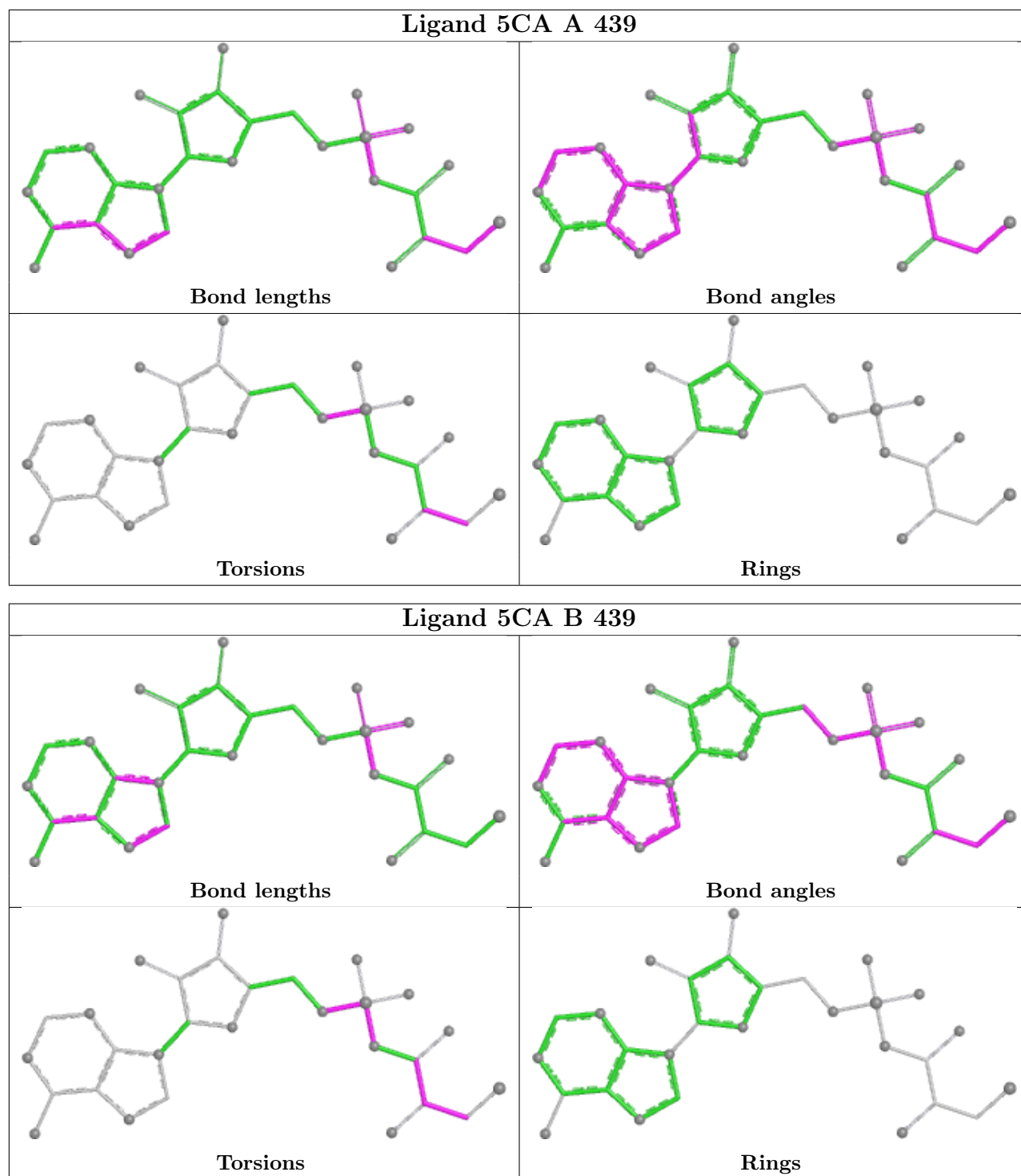
There are no ring outliers.

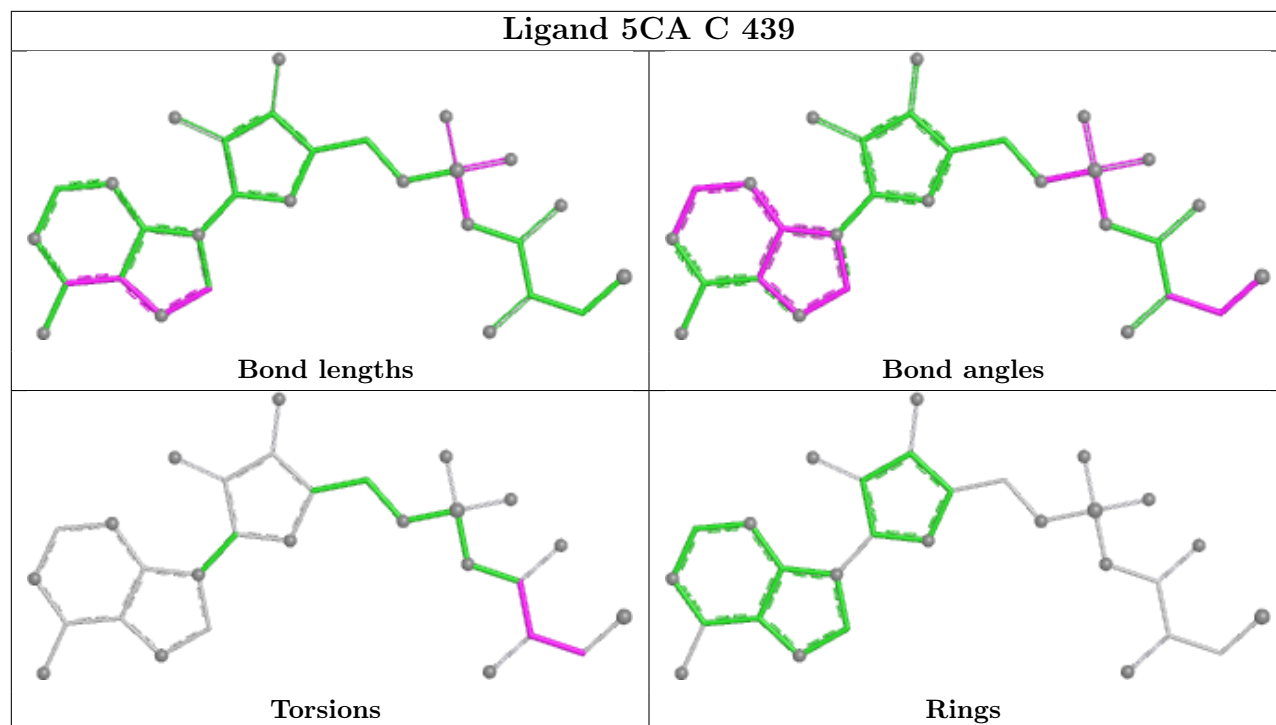
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	439	5CA	1	0
2	B	439	5CA	1	0
2	C	439	5CA	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient

equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	441/458 (96%)	0.23	6 (1%) 73 67	44, 55, 77, 84	0
1	B	442/458 (96%)	0.16	2 (0%) 87 85	45, 58, 69, 81	0
1	C	442/458 (96%)	0.05	3 (0%) 84 80	32, 63, 86, 101	0
All	All	1325/1374 (96%)	0.15	11 (0%) 82 78	32, 58, 81, 101	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	264	ALA	3.0
1	A	270	VAL	3.0
1	A	268	SER	2.5
1	B	268	SER	2.5
1	C	434	VAL	2.3
1	A	222	TYR	2.3
1	A	267	GLU	2.2
1	A	165	VAL	2.2
1	B	200	GLY	2.2
1	C	353	LEU	2.1
1	C	239	TYR	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](#)

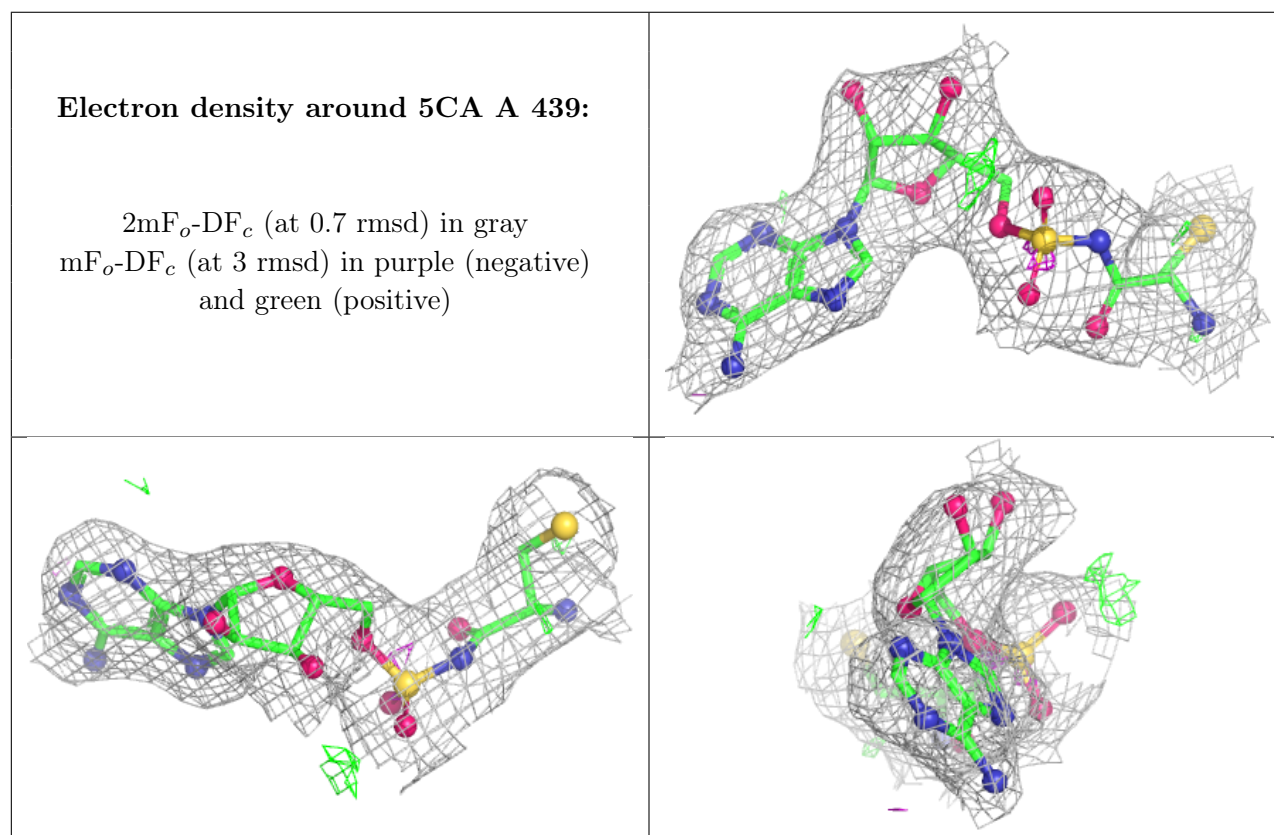
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

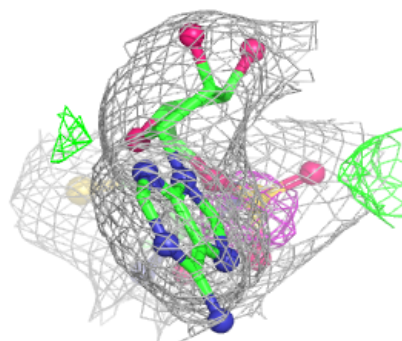
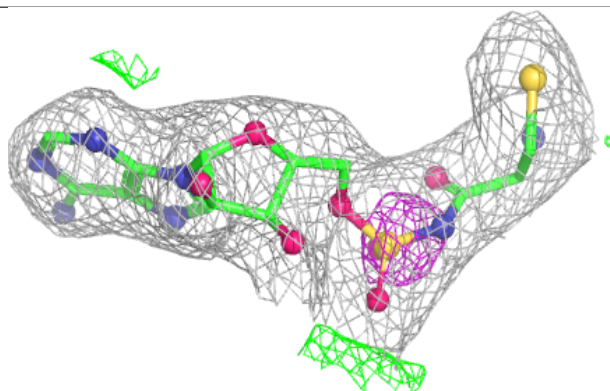
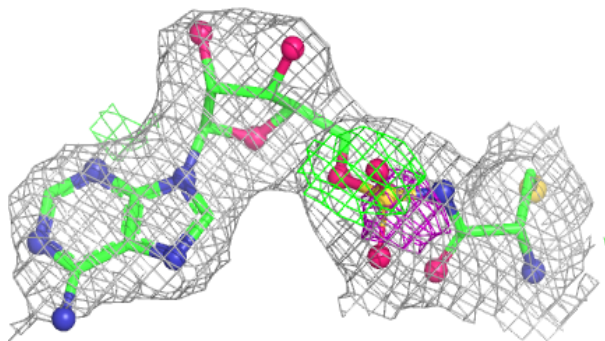
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	5CA	A	439	29/29	0.95	0.08	46,48,50,51	0
2	5CA	B	439	29/29	0.95	0.09	44,50,53,54	0
2	5CA	C	439	29/29	0.97	0.07	37,43,46,48	0

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.

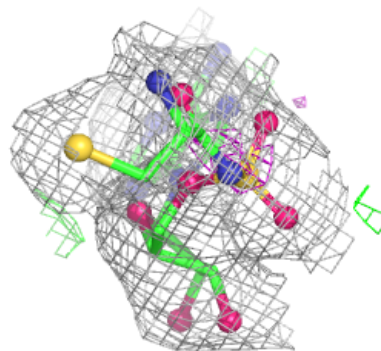
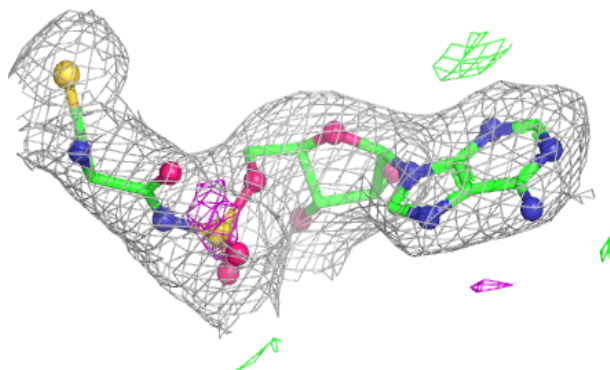
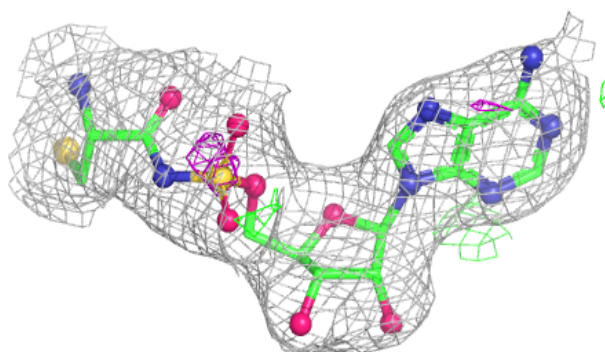


Electron density around 5CA B 439:

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

**Electron density around 5CA C 439:**

$2mF_o-DF_c$ (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.