



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:19 AM UTC

PDB ID : 5EZV / pdb_00005ezv
Title : X-ray crystal structure of AMP-activated protein kinase alpha-2/alpha-1 RIM chimaera (alpha-2(1-347)/alpha-1(349-401)/alpha-2(397-end) beta-1 gamma-1) co-crystallized with C2 (5-(5-hydroxyl-isoxazol-3-yl)-furan-2-phosphonic acid)
Authors : Langendorf, C.G.; Kemp, B.E.
Deposited on : 2015-11-26
Resolution : 2.99 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](#) ①) 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

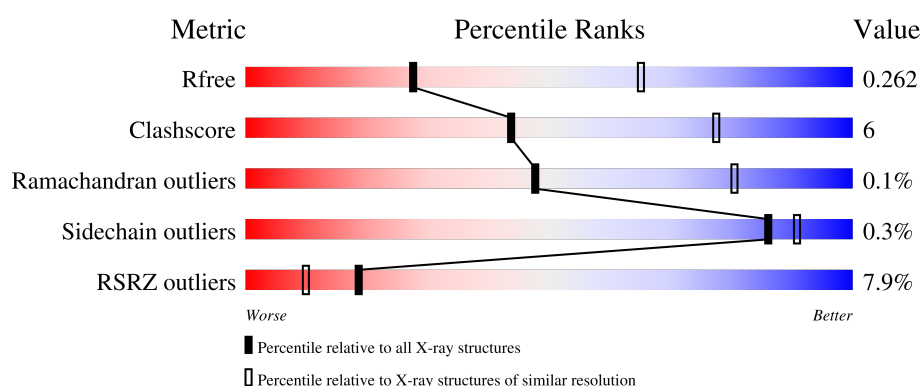
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.99 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	560	6% 71% 8% 21%
1	C	560	8% 70% 8% 21%
2	B	270	4% 55% 13% 31%
2	D	270	5% 59% 10% 31%
3	E	336	6% 82% 8% 10%

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Mol	Chain	Length	Quality of chain
3	F	336	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
1	TPO	A	172	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 14682 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 5'-AMP-activated protein kinase catalytic subunit alpha-2/alpha-1 RIM SWAP chimera.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	442	Total	C	N	O	P	S	0	0	0
			3462	2214	589	634	1	24			
1	C	441	Total	C	N	O	P	S	0	0	0
			3419	2186	588	621	1	23			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	initiating methionine	UNP P54646
A	-12	GLY	-	expression tag	UNP P54646
A	-11	SER	-	expression tag	UNP P54646
A	-10	SER	-	expression tag	UNP P54646
A	-9	HIS	-	expression tag	UNP P54646
A	-8	HIS	-	expression tag	UNP P54646
A	-7	HIS	-	expression tag	UNP P54646
A	-6	HIS	-	expression tag	UNP P54646
A	-5	HIS	-	expression tag	UNP P54646
A	-4	HIS	-	expression tag	UNP P54646
A	-3	HIS	-	expression tag	UNP P54646
A	-2	SER	-	expression tag	UNP P54646
A	-1	GLN	-	expression tag	UNP P54646
A	0	ASP	-	expression tag	UNP P54646
A	1	PRO	-	expression tag	UNP P54646
C	-13	MET	-	initiating methionine	UNP P54646
C	-12	GLY	-	expression tag	UNP P54646
C	-11	SER	-	expression tag	UNP P54646
C	-10	SER	-	expression tag	UNP P54646
C	-9	HIS	-	expression tag	UNP P54646
C	-8	HIS	-	expression tag	UNP P54646
C	-7	HIS	-	expression tag	UNP P54646
C	-6	HIS	-	expression tag	UNP P54646
C	-5	HIS	-	expression tag	UNP P54646

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-4	HIS	-	expression tag	UNP P54646
C	-3	HIS	-	expression tag	UNP P54646
C	-2	SER	-	expression tag	UNP P54646
C	-1	GLN	-	expression tag	UNP P54646
C	0	ASP	-	expression tag	UNP P54646
C	1	PRO	-	expression tag	UNP P54646

- Molecule 2 is a protein called 5'-AMP-activated protein kinase subunit beta-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	185	Total	C	N	O	P	S	0	0	0
			1400	904	233	257	1	5			
2	D	186	Total	C	N	O	P	S	0	0	0
			1422	914	234	267	1	6			

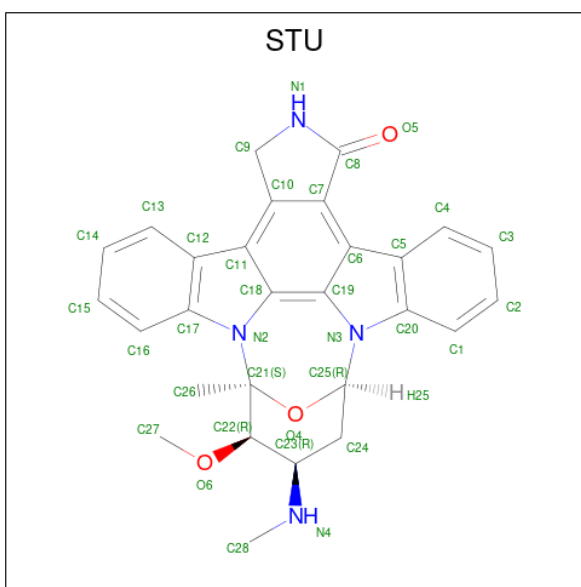
- Molecule 3 is a protein called 5'-AMP-activated protein kinase subunit gamma-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	E	302	Total	C	N	O	S		0	0	0
			2362	1538	391	426	7				
3	F	300	Total	C	N	O	S		0	1	0
			2351	1529	393	422	7				

There are 12 discrepancies between the modelled and reference sequences:

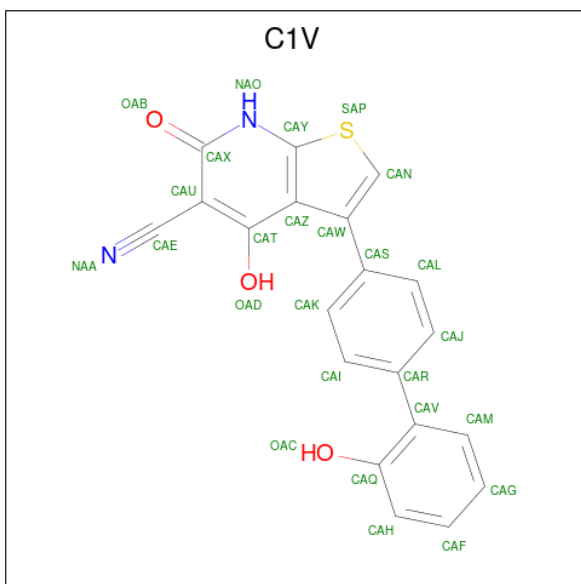
Chain	Residue	Modelled	Actual	Comment	Reference
E	-4	MET	-	initiating methionine	UNP P54619
E	-3	ALA	-	expression tag	UNP P54619
E	-2	ASP	-	expression tag	UNP P54619
E	-1	LEU	-	expression tag	UNP P54619
E	0	ASN	-	expression tag	UNP P54619
E	1	TRP	-	expression tag	UNP P54619
F	-4	MET	-	initiating methionine	UNP P54619
F	-3	ALA	-	expression tag	UNP P54619
F	-2	ASP	-	expression tag	UNP P54619
F	-1	LEU	-	expression tag	UNP P54619
F	0	ASN	-	expression tag	UNP P54619
F	1	TRP	-	expression tag	UNP P54619

- Molecule 4 is STAUROSPORINE (CCD ID: STU) (formula: C₂₈H₂₆N₄O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			35	28	4	3		
4	C	1	Total	C	N	O	0	0
			35	28	4	3		

- Molecule 5 is 3-[4-(2-hydroxyphenyl)phenyl]-4-oxidanyl-6-oxidanylidene-7H-thieno[2,3-b]pyridine-5-carbonitrile (CCD ID: C1V) (formula: $C_{20}H_{12}N_2O_3S$).



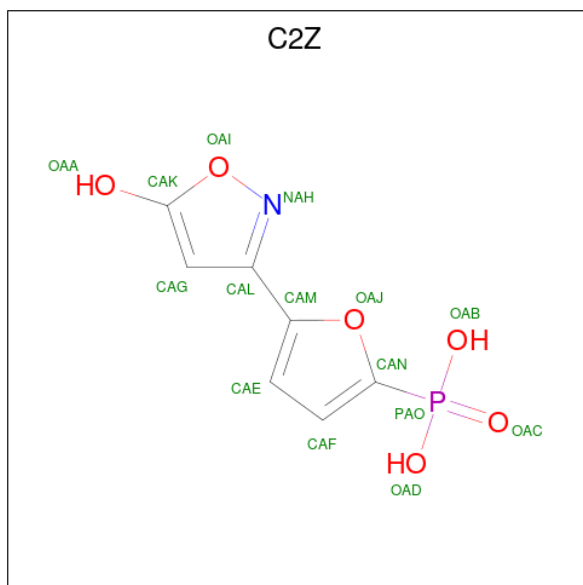
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			26	20	2	3	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	C	1	Total	C	N	O	S	0	0
			26	20	2	3	1		

- Molecule 6 is 5-(5-hydroxyl-isoxazol-3-yl)-furan-2-phosphonic acid (CCD ID: C2Z) (formula: $C_7H_6NO_6P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	E	1	Total	C	N	O	P	0	0
			15	7	1	6	1		
6	E	1	Total	C	N	O	P	0	0
			15	7	1	6	1		
6	F	1	Total	C	N	O	P	0	0
			15	7	1	6	1		
6	F	1	Total	C	N	O	P	0	0
			15	7	1	6	1		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	32	Total	O	0	0
			32	32		
7	B	8	Total	O	0	0
			8	8		
7	C	25	Total	O	0	0
			25	25		
7	D	7	Total	O	0	0
			7	7		

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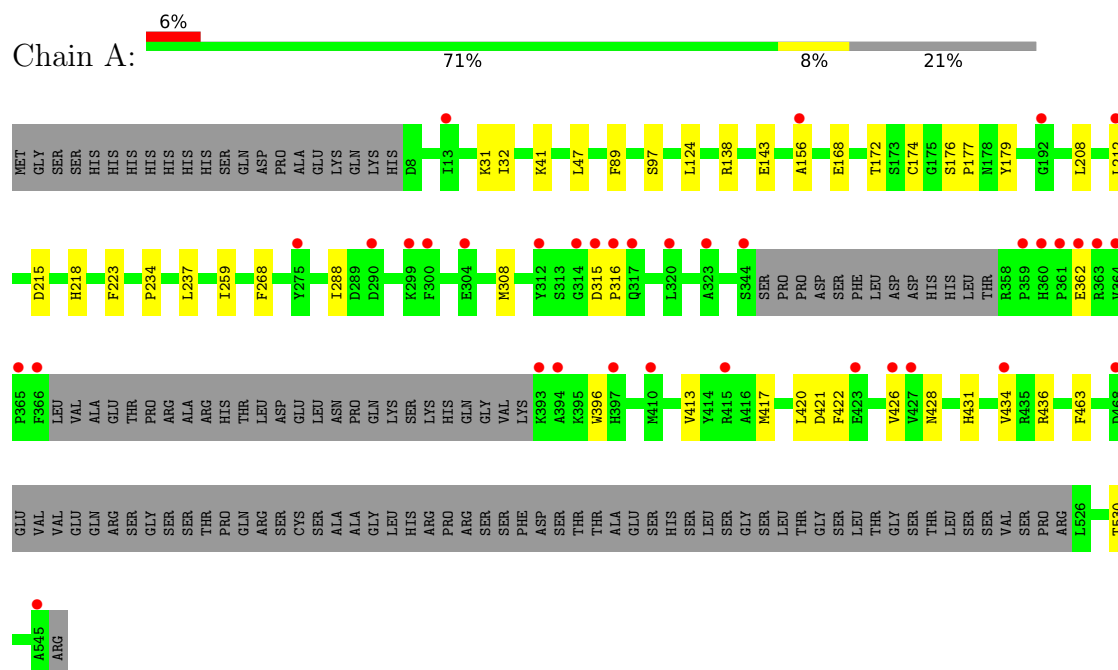
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	E	7	Total	O	0	0
			7	7		
7	F	5	Total	O	0	0
			5	5		

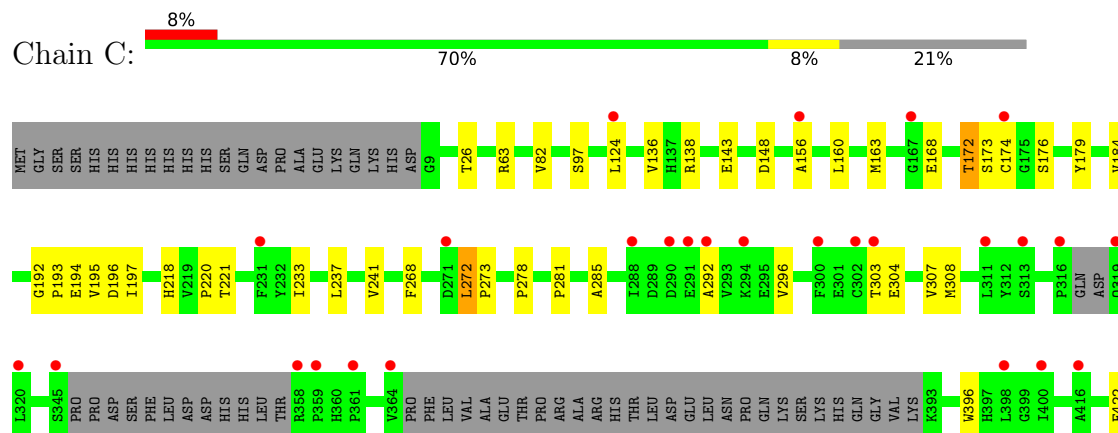
3 Residue-property plots

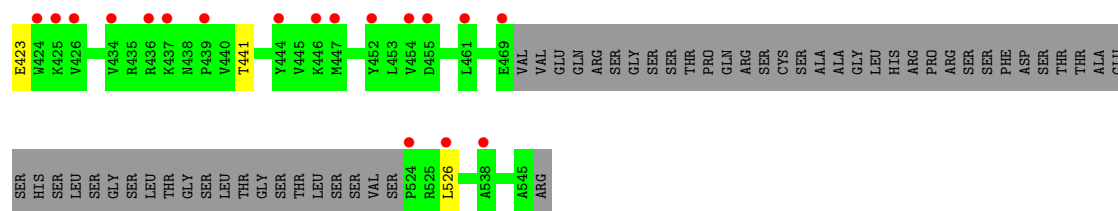
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: 5'-AMP-activated protein kinase catalytic subunit alpha-2/alpha-1 RIM SWAP chimera

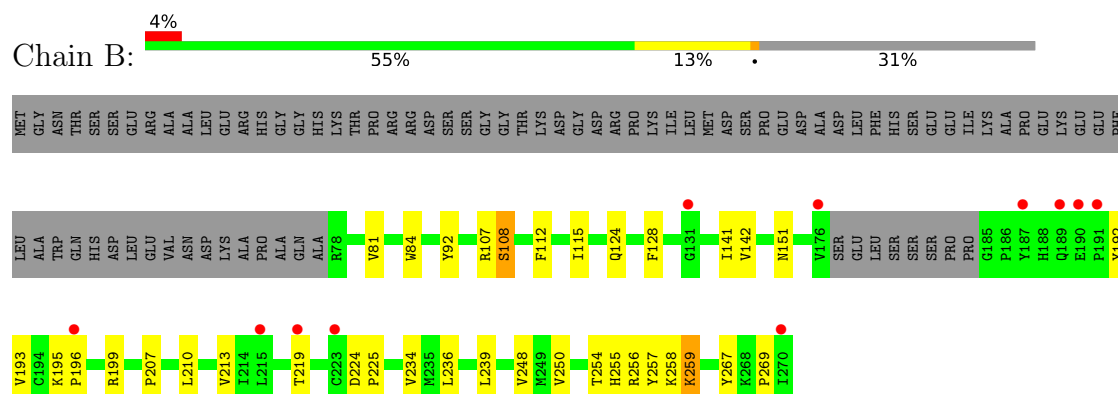


- Molecule 1: 5'-AMP-activated protein kinase catalytic subunit alpha-2/alpha-1 RIM SWAP chimera

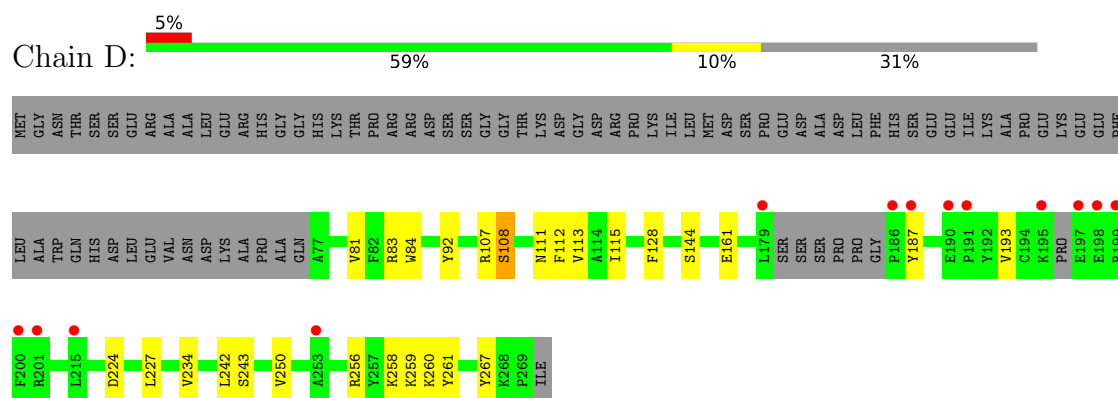




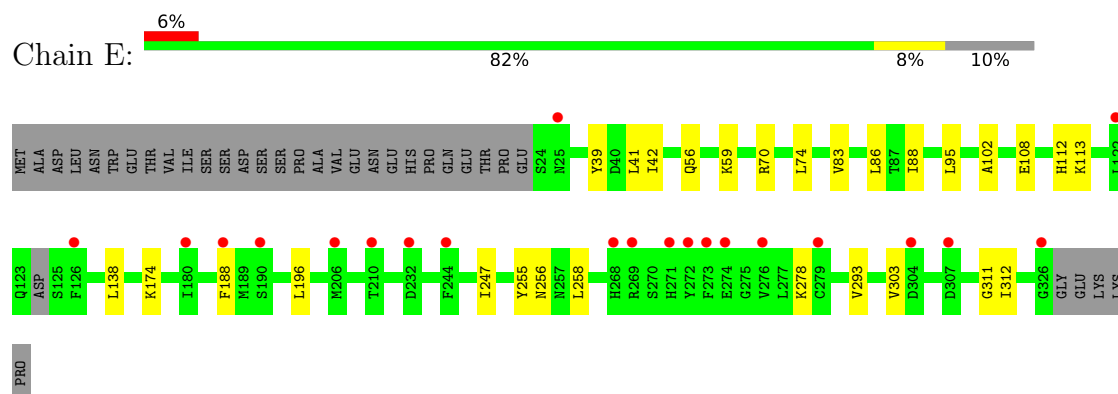
• Molecule 2: 5'-AMP-activated protein kinase subunit beta-1



• Molecule 2: 5'-AMP-activated protein kinase subunit beta-1



• Molecule 3: 5'-AMP-activated protein kinase subunit gamma-1



• Molecule 3: 5'-AMP-activated protein kinase subunit gamma-1





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.61Å 134.11Å 141.29Å 90.00° 93.16° 90.00°	Depositor
Resolution (Å)	49.30 – 2.99 49.30 – 2.99	Depositor EDS
% Data completeness (in resolution range)	99.3 (49.30-2.99) 99.3 (49.30-2.99)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.66 (at 3.01Å)	Xtriage
Refinement program	BUSTER-TNT 1.10.0	Depositor
R, R_{free}	0.224 , 0.245 0.241 , 0.262	Depositor DCC
R_{free} test set	2867 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	67.5	Xtriage
Anisotropy	0.203	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 44.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	14682	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.87% 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: C2Z, SEP, STU, C1V, TPO

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	0.72	0/3528	1.30	5/4778 (0.1%)
1	C	0.71	0/3485	1.27	7/4724 (0.1%)
2	B	0.70	0/1431	1.13	2/1964 (0.1%)
2	D	0.71	0/1452	1.14	2/1989 (0.1%)
3	E	0.73	0/2411	1.30	2/3281 (0.1%)
3	F	0.72	0/2402	1.27	4/3269 (0.1%)
All	All	0.72	0/14709	1.26	22/20005 (0.1%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	156	ALA	N-CA-C	7.38	122.01	112.12
3	E	255	TYR	N-CA-C	6.91	119.84	111.82
1	C	156	ALA	N-CA-C	6.38	120.84	111.34
1	A	218	HIS	CA-C-N	6.06	125.18	120.33
1	A	218	HIS	C-N-CA	6.06	125.18	120.33
1	C	278	PRO	N-CA-C	5.80	122.39	113.75
3	F	305	GLU	N-CA-C	-5.76	105.00	111.28
1	C	422	PHE	CA-CB-CG	5.72	119.52	113.80
2	B	259	LYS	N-CA-CB	-5.71	101.78	111.20
1	C	281	PRO	N-CA-C	5.58	122.81	113.78
1	C	97	SER	N-CA-C	5.54	117.32	111.28
3	F	276	VAL	N-CA-C	5.46	116.19	110.62
1	A	97	SER	N-CA-C	5.45	117.22	111.28
2	B	219	THR	N-CA-C	5.35	118.03	111.82
3	F	57	VAL	N-CA-C	5.26	117.63	111.05
1	C	148	ASP	CA-CB-CG	5.24	117.84	112.60
1	A	426	VAL	N-CA-C	5.24	111.81	106.21
1	C	82	VAL	CB-CA-C	5.19	116.33	110.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	83	VAL	N-CA-C	5.17	116.26	111.81
3	F	265	ALA	N-CA-C	5.10	116.84	111.28
2	D	161	GLU	CA-C-N	5.05	127.11	120.60
2	D	161	GLU	C-N-CA	5.05	127.11	120.60

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	3462	0	3311	35	0
1	C	3419	0	3255	42	1
2	B	1400	0	1308	30	0
2	D	1422	0	1317	22	1
3	E	2362	0	2379	18	0
3	F	2351	0	2368	41	0
4	A	35	0	26	2	0
4	C	35	0	26	3	0
5	A	26	0	12	0	0
5	C	26	0	12	1	0
6	E	30	0	0	0	0
6	F	30	0	0	0	0
7	A	32	0	0	0	0
7	B	8	0	0	0	0
7	C	25	0	0	0	0
7	D	7	0	0	0	0
7	E	7	0	0	0	0
7	F	5	0	0	0	0
All	All	14682	0	14014	175	1

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 6.

All (175) 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:292:ALA:O	1:C:296:VAL:HG23	1.55	1.06
1:A:138:ARG:NH2	1:A:172:TPO:O2P	1.89	1.05
3:E:256:ASN:O	3:E:256:ASN:ND2	1.92	1.03
3:F:262:VAL:O	3:F:266:LEU:HD13	1.60	1.00
1:C:296:VAL:CG1	1:C:303:THR:HA	1.94	0.97
2:D:107:ARG:HG2	2:D:108:SEP:O	1.64	0.96
2:B:224:ASP:OD1	2:B:225:PRO:HD2	1.67	0.92
1:C:237:LEU:HD11	1:C:241:VAL:CG1	2.02	0.89
1:A:124:LEU:HD11	1:A:259:ILE:HG12	1.53	0.89
1:A:138:ARG:HH22	1:A:172:TPO:HB	1.38	0.89
2:B:224:ASP:OD1	2:B:225:PRO:CD	2.24	0.84
1:C:296:VAL:HG13	1:C:303:THR:HA	1.60	0.84
3:F:240:ILE:HG22	3:F:275:GLY:HA3	1.57	0.84
1:C:296:VAL:HG11	1:C:303:THR:HA	1.59	0.81
1:C:396:TRP:CH2	2:D:250:VAL:HG12	2.15	0.81
1:A:138:ARG:NH2	1:A:172:TPO:HB	1.96	0.80
2:D:256:ARG:HG2	2:D:256:ARG:HH11	1.44	0.80
1:C:272:LEU:C	1:C:272:LEU:HD13	2.08	0.78
3:E:256:ASN:HD22	3:E:256:ASN:C	1.91	0.78
2:D:256:ARG:HG2	2:D:256:ARG:NH1	1.99	0.77
3:F:239:ASP:OD1	3:F:240:ILE:N	2.19	0.75
1:C:237:LEU:HD11	1:C:241:VAL:HG13	1.70	0.72
1:C:138:ARG:NH1	1:C:172:TPO:O2P	2.20	0.72
1:C:441:THR:HG22	1:C:526:LEU:HD22	1.71	0.72
1:C:192:GLY:O	1:C:195:VAL:HG22	1.90	0.71
2:B:256:ARG:HG2	2:B:257:TYR:O	1.91	0.70
1:C:396:TRP:HH2	2:D:250:VAL:HG12	1.57	0.69
1:C:304:GLU:O	1:C:307:VAL:HG12	1.92	0.69
1:A:124:LEU:HD23	1:A:268:PHE:CZ	2.31	0.66
4:C:601:STU:H261	4:C:601:STU:H16	1.78	0.65
3:F:57:VAL:HG22	3:F:61:PHE:CE2	2.32	0.65
3:F:241:TYR:HD1	3:F:266:LEU:HD11	1.62	0.64
1:A:422:PHE:CE2	1:A:436:ARG:HD3	2.33	0.64
1:A:422:PHE:HE2	1:A:436:ARG:HD3	1.62	0.64
4:A:601:STU:H261	4:A:601:STU:H16	1.79	0.64
1:A:138:ARG:CZ	1:A:172:TPO:O2P	2.47	0.63
1:A:396:TRP:CH2	2:B:250:VAL:HG12	2.33	0.63
3:E:41:LEU:HD13	3:E:138:LEU:HD21	1.82	0.62
3:F:239:ASP:CG	3:F:240:ILE:H	2.06	0.62
1:A:168:GLU:O	2:B:234:VAL:HG12	1.98	0.62
1:C:233:ILE:HG23	1:C:237:LEU:HD23	1.81	0.62
3:F:241:TYR:CD1	3:F:266:LEU:HD11	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:259:LYS:HB2	3:E:39:TYR:OH	2.01	0.61
1:C:237:LEU:HD11	1:C:241:VAL:HG11	1.81	0.60
3:F:41:LEU:HD11	3:F:167:LEU:HD13	1.83	0.60
1:C:272:LEU:C	1:C:272:LEU:CD1	2.74	0.60
1:C:272:LEU:HD13	1:C:273:PRO:N	2.16	0.59
1:C:237:LEU:CD1	1:C:241:VAL:CG1	2.80	0.59
2:D:83:ARG:NE	2:D:111:ASN:OD1	2.33	0.58
2:B:195:LYS:CB	2:B:196:PRO:HD2	2.32	0.58
1:C:143:GLU:HB3	4:C:601:STU:H281	1.84	0.58
2:D:107:ARG:HD2	2:D:112:PHE:CZ	2.39	0.57
1:A:143:GLU:HB3	4:A:601:STU:H281	1.86	0.57
1:A:413:VAL:O	1:A:417:MET:HG3	2.05	0.57
3:F:266:LEU:HD12	3:F:266:LEU:N	2.19	0.57
1:C:168:GLU:O	2:D:234:VAL:HG12	2.03	0.57
1:C:138:ARG:NH2	1:C:172:TPO:O2P	2.38	0.56
1:A:31:LYS:NZ	2:B:108:SEP:O2P	2.30	0.55
2:B:195:LYS:CB	2:B:199:ARG:CB	2.85	0.54
1:C:173:SER:HA	1:C:184:VAL:CG2	2.37	0.54
1:A:413:VAL:HG12	1:A:417:MET:HE2	1.88	0.54
3:F:240:ILE:HG22	3:F:275:GLY:CA	2.36	0.54
1:A:428:ASN:HB3	1:A:431:HIS:HB3	1.89	0.54
1:C:176:SER:HB3	1:C:179:TYR:HD2	1.73	0.53
2:B:224:ASP:OD1	2:B:225:PRO:N	2.42	0.52
1:C:193:PRO:O	1:C:197:ILE:HG12	2.10	0.51
3:E:88:ILE:HG23	3:E:247:ILE:HG23	1.93	0.51
3:F:266:LEU:O	3:F:270:SER:N	2.43	0.51
2:B:250:VAL:HG22	2:B:267:TYR:CD2	2.45	0.51
1:A:420:LEU:O	1:A:421:ASP:HB3	2.10	0.51
1:C:173:SER:HA	1:C:184:VAL:HG22	1.91	0.51
1:C:396:TRP:CH2	2:D:250:VAL:CG1	2.89	0.51
1:C:218:HIS:CE1	1:C:220:PRO:HD2	2.46	0.51
1:A:396:TRP:HH2	2:B:250:VAL:HG12	1.74	0.51
3:F:39:TYR:HA	3:F:42:ILE:HD12	1.93	0.51
1:A:234:PRO:HD2	1:A:237:LEU:HD12	1.93	0.50
3:F:266:LEU:N	3:F:266:LEU:CD1	2.74	0.50
1:A:362:GLU:HG2	3:E:70:ARG:HH21	1.76	0.50
1:A:396:TRP:HB2	2:B:213:VAL:HG11	1.94	0.50
1:A:47:LEU:HB2	1:A:89:PHE:HB2	1.93	0.50
2:B:124:GLN:HE21	2:B:142:VAL:HG11	1.77	0.49
3:F:32:PHE:HD2	3:F:33:MET:HE2	1.76	0.49
2:B:107:ARG:HB2	2:B:112:PHE:CE2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:41:LEU:HD11	3:F:167:LEU:CD1	2.41	0.49
3:E:174:LYS:HG2	3:E:293:VAL:HG13	1.94	0.49
1:A:138:ARG:NH1	1:A:172:TPO:O2P	2.45	0.49
3:E:256:ASN:ND2	3:E:256:ASN:C	2.61	0.49
2:D:258:LYS:NZ	3:F:37:ARG:HH12	2.11	0.48
2:B:84:TRP:HB3	2:B:112:PHE:HB2	1.94	0.48
2:D:242:LEU:HD23	2:D:243:SER:O	2.13	0.48
1:C:296:VAL:HG11	1:C:303:THR:CA	2.37	0.48
3:F:61:PHE:HA	3:F:64:LEU:HD12	1.96	0.48
2:D:260:LYS:NZ	3:F:39:TYR:O	2.47	0.48
2:B:92:TYR:HB2	2:B:128:PHE:HB3	1.93	0.48
3:E:108:GLU:O	3:E:112:HIS:HB2	2.14	0.48
3:F:262:VAL:O	3:F:266:LEU:CD1	2.48	0.48
2:B:195:LYS:CB	2:B:196:PRO:CD	2.92	0.48
1:C:307:VAL:HG13	1:C:308:MET:N	2.28	0.47
1:A:530:THR:HG23	2:B:255:HIS:CE1	2.49	0.47
1:A:288:ILE:HG12	1:A:308:MET:HE1	1.95	0.47
3:E:56:GLN:HB2	3:E:59:LYS:HE3	1.96	0.47
2:B:248:VAL:HA	2:B:269:PRO:HA	1.97	0.47
3:E:74:LEU:HD21	3:E:86:LEU:HB2	1.97	0.47
3:F:25:ASN:O	3:F:28:VAL:HG23	2.15	0.47
3:E:56:GLN:HA	3:E:113:LYS:HA	1.96	0.47
2:B:124:GLN:NE2	2:B:142:VAL:HG11	2.30	0.47
2:D:250:VAL:HG22	2:D:267:TYR:CD2	2.50	0.46
1:A:212:LEU:HB2	1:A:215:ASP:HB2	1.97	0.46
1:A:463:PHE:HB2	2:B:239:LEU:HB3	1.97	0.46
2:D:92:TYR:HB2	2:D:128:PHE:HB3	1.98	0.46
3:F:303:VAL:HG13	3:F:307:ASP:O	2.16	0.46
1:C:441:THR:HG22	1:C:526:LEU:CD2	2.42	0.46
1:A:177:PRO:HB3	1:A:223:PHE:HE1	1.81	0.46
1:C:272:LEU:HD13	1:C:273:PRO:O	2.15	0.46
1:A:421:ASP:O	1:A:421:ASP:CG	2.59	0.46
1:A:422:PHE:HB3	1:A:434:VAL:CG2	2.46	0.46
2:B:207:PRO:HD2	2:B:210:LEU:HD12	1.98	0.46
3:F:74:LEU:HD22	3:F:114:ILE:HG21	1.97	0.46
3:F:239:ASP:CG	3:F:240:ILE:N	2.72	0.46
1:C:423:GLU:HB3	2:D:187:TYR:HB3	1.98	0.45
1:C:26:THR:HG21	1:C:160:LEU:HG	1.99	0.45
3:F:41:LEU:HD13	3:F:171:ARG:CG	2.47	0.45
1:C:63:ARG:HE	1:C:163:MET:HG3	1.82	0.45
1:C:138:ARG:CZ	1:C:172:TPO:O2P	2.65	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:81:VAL:HG22	2:B:115:ILE:HG12	1.97	0.45
2:B:192:TYR:CE1	2:B:193:VAL:O	2.70	0.45
1:A:172:TPO:HG23	1:A:174:CYS:SG	2.57	0.45
1:A:208:LEU:HB3	1:A:237:LEU:HD21	1.99	0.45
3:F:174:LYS:HG3	3:F:293:VAL:HG13	1.99	0.45
1:C:307:VAL:CG1	1:C:308:MET:N	2.80	0.44
3:F:25:ASN:HA	3:F:28:VAL:HG23	1.99	0.44
3:F:88:ILE:HG23	3:F:247:ILE:HG23	2.00	0.44
1:C:124:LEU:HD23	1:C:268:PHE:CZ	2.53	0.44
2:D:256:ARG:HH11	2:D:256:ARG:CG	2.15	0.44
1:C:218:HIS:NE2	1:C:220:PRO:HD2	2.33	0.43
2:B:192:TYR:CD1	2:B:193:VAL:O	2.70	0.43
3:E:311:GLY:O	3:E:312:ILE:HD13	2.17	0.43
1:C:160:LEU:HD13	1:C:174:CYS:HB3	1.99	0.43
3:F:74:LEU:HD21	3:F:86:LEU:HB2	2.01	0.43
5:C:602:C1V:HAM	2:D:113:VAL:HG12	2.00	0.43
2:D:256:ARG:HD3	2:D:261:TYR:HE1	1.84	0.43
3:F:188:PHE:HB2	3:F:196:LEU:HD21	2.00	0.43
3:E:188:PHE:HB2	3:E:196:LEU:HD21	2.02	0.42
3:F:208:ARG:HB2	3:F:211:THR:HG23	2.00	0.42
3:F:231:VAL:HG22	3:F:237:VAL:HG22	2.02	0.42
3:F:303:VAL:HG12	3:F:304:ASP:O	2.19	0.42
2:B:250:VAL:HG21	2:B:267:TYR:CE2	2.54	0.42
3:E:39:TYR:HA	3:E:42:ILE:HD12	2.02	0.42
3:F:177:LYS:HG2	3:F:293:VAL:HG11	2.01	0.42
3:F:303:VAL:CG1	3:F:304:ASP:N	2.83	0.42
2:D:81:VAL:HG22	2:D:115:ILE:HG12	2.00	0.41
3:E:102:ALA:HA	3:E:256:ASN:ND2	2.35	0.41
3:F:25:ASN:C	3:F:28:VAL:HG23	2.44	0.41
2:D:224:ASP:HB3	2:D:227:LEU:HD12	2.01	0.41
3:E:278:LYS:HD2	3:E:303:VAL:HG21	2.01	0.41
3:F:280:TYR:HB2	3:F:283:GLU:HG3	2.02	0.41
3:F:95:LEU:HD22	3:F:258:LEU:HD11	2.03	0.41
1:A:315:ASP:HA	1:A:316:PRO:HD3	1.98	0.41
3:F:193:LEU:HD13	3:F:288:ILE:HD13	2.02	0.41
1:A:32:ILE:HD11	1:A:41:LYS:HD3	2.02	0.41
1:C:136:VAL:HG23	1:C:196:ASP:OD2	2.21	0.41
2:B:141:ILE:HG22	2:B:151:ASN:ND2	2.35	0.41
2:B:236:LEU:HD23	2:B:254:THR:HG22	2.03	0.41
2:B:257:TYR:O	2:B:258:LYS:C	2.63	0.41
4:C:601:STU:H261	4:C:601:STU:C16	2.48	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:41:LEU:HD13	3:F:171:ARG:HG2	2.03	0.41
1:A:176:SER:HB2	1:A:179:TYR:HD2	1.85	0.41
2:D:84:TRP:HB3	2:D:112:PHE:HB2	2.03	0.40
2:D:107:ARG:CG	2:D:108:SEP:O	2.51	0.40
1:A:417:MET:HB3	1:A:422:PHE:HB2	2.03	0.40
1:C:194:GLU:HG2	1:C:195:VAL:N	2.36	0.40
1:C:218:HIS:ND1	1:C:221:THR:HG23	2.36	0.40
3:F:303:VAL:HG12	3:F:304:ASP:N	2.36	0.40
3:E:95:LEU:HD22	3:E:258:LEU:HD11	2.03	0.40
3:F:32:PHE:CE1	3:F:36:HIS:CD2	3.10	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:285:ALA:O	2:D:144:SER:OG[2_946]	2.16	0.04

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	433/560 (77%)	422 (98%)	11 (2%)	0	100	100
1	C	430/560 (77%)	413 (96%)	17 (4%)	0	100	100
2	B	180/270 (67%)	177 (98%)	3 (2%)	0	100	100
2	D	179/270 (66%)	174 (97%)	4 (2%)	1 (1%)	21	56
3	E	298/336 (89%)	293 (98%)	5 (2%)	0	100	100
3	F	299/336 (89%)	293 (98%)	6 (2%)	0	100	100
All	All	1819/2332 (78%)	1772 (97%)	46 (2%)	1 (0%)	48	80

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	193	VAL

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	358/497 (72%)	358 (100%)	0	100	100
1	C	349/497 (70%)	348 (100%)	1 (0%)	86	91
2	B	145/239 (61%)	145 (100%)	0	100	100
2	D	149/239 (62%)	148 (99%)	1 (1%)	76	86
3	E	259/308 (84%)	259 (100%)	0	100	100
3	F	255/308 (83%)	253 (99%)	2 (1%)	73	86
All	All	1515/2088 (73%)	1511 (100%)	4 (0%)	86	91

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	272	LEU
2	D	259	LYS
3	F	147	ARG
3	F	269	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	ASN
1	C	66	GLN
1	C	397	HIS
2	D	135	HIS
2	D	150	ASN
2	D	212	GLN
2	D	237	ASN
3	E	67	ASN
3	E	80	GLN

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Mol	Chain	Res	Type
3	E	93	ASN
3	E	197	GLN
3	E	203	ASN
3	E	268	HIS
3	E	290	ASN
3	E	320	GLN
3	F	56	GLN
3	F	67	ASN
3	F	135	ASN
3	F	148	ASN
3	F	257	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 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	TPO	C	172	1	8,10,11	0.82	0	10,14,16	1.32	1 (10%)
1	TPO	A	172	1	8,10,11	0.87	0	10,14,16	1.25	0
2	SEP	D	108	2	8,9,10	0.95	0	7,12,14	1.25	1 (14%)
2	SEP	B	108	2	8,9,10	0.94	0	7,12,14	1.45	1 (14%)

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	TPO	C	172	1	-	2/9/11/13	-
1	TPO	A	172	1	-	1/9/11/13	-
2	SEP	D	108	2	-	2/6/8/10	-
2	SEP	B	108	2	-	4/6/8/10	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	108	SEP	OG-CB-CA	2.89	110.96	108.14
1	C	172	TPO	O-C-CA	-2.09	119.39	124.77
2	D	108	SEP	O3P-P-O2P	2.07	115.57	107.80

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	172	TPO	O-C-CA-CB
2	B	108	SEP	C-CA-CB-OG
2	B	108	SEP	CB-OG-P-O1P
2	B	108	SEP	CB-OG-P-O2P
2	B	108	SEP	CB-OG-P-O3P
1	C	172	TPO	O-C-CA-CB
2	D	108	SEP	CB-OG-P-O2P
1	C	172	TPO	CB-OG1-P-O1P
2	D	108	SEP	N-CA-CB-OG

There are no ring outliers.

4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	172	TPO	3	0
1	A	172	TPO	6	0
2	D	108	SEP	2	0
2	B	108	SEP	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

8 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
6	C2Z	E	401	-	14,16,16	3.55	8 (57%)	13,24,24	2.08	4 (30%)
5	C1V	A	602	-	28,29,29	3.22	9 (32%)	32,42,42	5.18	7 (21%)
5	C1V	C	602	-	28,29,29	3.24	9 (32%)	32,42,42	5.25	6 (18%)
6	C2Z	F	402	-	14,16,16	3.53	8 (57%)	13,24,24	2.12	4 (30%)
4	STU	C	601	-	39,42,42	2.69	12 (30%)	50,68,68	4.10	24 (48%)
4	STU	A	601	-	39,42,42	2.70	13 (33%)	50,68,68	4.01	26 (52%)
6	C2Z	F	401	-	14,16,16	3.40	8 (57%)	13,24,24	2.16	5 (38%)
6	C2Z	E	402	-	14,16,16	3.04	7 (50%)	13,24,24	2.08	3 (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
6	C2Z	E	401	-	-	0/4/10/10	0/2/2/2
5	C1V	A	602	-	-	0/8/10/10	0/4/4/4
5	C1V	C	602	-	-	0/8/10/10	0/4/4/4
6	C2Z	F	402	-	-	0/4/10/10	0/2/2/2
4	STU	C	601	-	-	2/4/42/42	-
4	STU	A	601	-	-	2/4/42/42	-
6	C2Z	F	401	-	-	0/4/10/10	0/2/2/2
6	C2Z	E	402	-	-	0/4/10/10	0/2/2/2

All (74) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	602	C1V	CAE-NAA	9.86	1.32	1.14
5	A	602	C1V	CAE-NAA	9.81	1.32	1.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	602	C1V	OAB-CAX	9.78	1.42	1.23
5	A	602	C1V	OAB-CAX	9.77	1.42	1.23
6	E	402	C2Z	OAI-CAK	-6.81	1.35	1.41
6	E	401	C2Z	OAJ-CAN	-6.74	1.32	1.37
6	F	402	C2Z	OAI-CAK	-6.64	1.35	1.41
6	F	401	C2Z	PAO-OAC	6.28	1.61	1.49
4	C	601	STU	C9-C10	-6.28	1.45	1.50
4	A	601	STU	C9-C10	-6.10	1.45	1.50
6	F	402	C2Z	PAO-OAC	6.02	1.60	1.49
6	E	401	C2Z	PAO-OAC	5.92	1.60	1.49
6	E	401	C2Z	CAM-CAL	-5.62	1.33	1.44
6	F	401	C2Z	CAM-CAL	-5.58	1.34	1.44
6	F	401	C2Z	OAI-CAK	-5.56	1.36	1.41
6	F	402	C2Z	CAM-CAL	-5.49	1.34	1.44
6	E	402	C2Z	CAM-CAL	-5.47	1.34	1.44
6	E	401	C2Z	OAI-CAK	-5.47	1.36	1.41
6	F	401	C2Z	OAJ-CAN	-5.42	1.33	1.37
4	A	601	STU	C12-C17	5.29	1.49	1.41
4	A	601	STU	C5-C20	5.25	1.49	1.41
6	F	402	C2Z	OAJ-CAN	-5.23	1.33	1.37
4	A	601	STU	C7-C10	5.21	1.48	1.39
4	C	601	STU	C5-C20	5.20	1.49	1.41
4	C	601	STU	C12-C17	5.19	1.48	1.41
4	A	601	STU	C11-C10	5.08	1.50	1.40
4	C	601	STU	C7-C10	5.06	1.48	1.39
4	C	601	STU	C11-C10	5.00	1.50	1.40
4	A	601	STU	C6-C7	4.90	1.49	1.40
5	C	602	C1V	CAS-CAW	4.86	1.56	1.48
4	C	601	STU	C6-C7	4.83	1.49	1.40
5	A	602	C1V	CAS-CAW	4.82	1.56	1.48
5	A	602	C1V	CAE-CAU	4.78	1.53	1.43
5	C	602	C1V	CAE-CAU	4.74	1.53	1.43
4	C	601	STU	C25-N3	4.64	1.49	1.45
4	A	601	STU	C19-C18	4.59	1.50	1.40
4	A	601	STU	C25-N3	4.54	1.49	1.45
4	C	601	STU	C19-C18	4.54	1.50	1.40
4	A	601	STU	C11-C18	4.46	1.49	1.41
4	C	601	STU	C11-C18	4.41	1.49	1.41
6	E	402	C2Z	OAJ-CAN	-4.24	1.34	1.37
4	A	601	STU	C6-C19	3.91	1.48	1.41
4	C	601	STU	C6-C19	3.88	1.48	1.41
6	E	401	C2Z	CAG-CAL	-3.16	1.33	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	F	401	C2Z	CAG-CAL	-3.16	1.33	1.41
6	E	402	C2Z	CAG-CAL	-3.11	1.33	1.41
6	F	402	C2Z	CAG-CAL	-3.06	1.33	1.41
5	A	602	C1V	CAX-NAO	-3.01	1.33	1.38
5	C	602	C1V	CAX-NAO	-3.01	1.33	1.38
6	F	402	C2Z	PAO-OAB	2.66	1.61	1.54
6	E	401	C2Z	PAO-OAB	2.65	1.61	1.54
6	E	402	C2Z	PAO-OAD	2.64	1.61	1.54
5	C	602	C1V	OAC-CAQ	2.58	1.41	1.36
5	A	602	C1V	OAC-CAQ	2.56	1.41	1.36
6	E	401	C2Z	OAJ-CAM	-2.52	1.34	1.38
6	E	402	C2Z	PAO-OAB	2.34	1.60	1.54
5	C	602	C1V	CAV-CAR	2.32	1.53	1.49
6	F	402	C2Z	PAO-OAD	-2.29	1.48	1.54
6	F	402	C2Z	OAJ-CAM	-2.28	1.34	1.38
6	F	401	C2Z	PAO-OAD	-2.28	1.48	1.54
6	E	401	C2Z	PAO-OAD	-2.24	1.48	1.54
4	C	601	STU	O5-C8	2.23	1.28	1.23
4	A	601	STU	O5-C8	2.23	1.28	1.23
6	F	401	C2Z	PAO-OAB	2.19	1.60	1.54
4	A	601	STU	C18-N2	-2.13	1.37	1.40
4	C	601	STU	C18-N2	-2.12	1.37	1.40
5	A	602	C1V	CAV-CAR	2.12	1.53	1.49
5	C	602	C1V	CAY-NAO	-2.12	1.35	1.38
6	F	401	C2Z	OAJ-CAM	-2.12	1.35	1.38
5	C	602	C1V	CAW-CAZ	2.10	1.49	1.44
5	A	602	C1V	CAW-CAZ	2.10	1.49	1.44
5	A	602	C1V	CAY-NAO	-2.10	1.35	1.38
4	A	601	STU	C9-N1	2.08	1.47	1.45
6	E	402	C2Z	OAJ-CAM	-2.04	1.35	1.38

All (79) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	602	C1V	CAU-CAE-NAA	-25.21	120.06	177.39
5	A	602	C1V	CAU-CAE-NAA	-24.64	121.36	177.39
5	A	602	C1V	CAN-SAP-CAY	12.60	99.32	90.33
5	C	602	C1V	CAN-SAP-CAY	12.59	99.32	90.33
4	C	601	STU	O5-C8-N1	-12.36	111.60	125.35
4	A	601	STU	O5-C8-N1	-11.91	112.10	125.35
4	C	601	STU	C19-C18-N2	-9.61	123.21	131.38
4	C	601	STU	O5-C8-C7	-9.58	114.58	128.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	601	STU	O5-C8-C7	-9.30	115.00	128.47
4	A	601	STU	C19-C18-N2	-9.09	123.65	131.38
4	C	601	STU	C10-C7-C8	-8.84	100.70	108.09
4	A	601	STU	C10-C7-C8	-8.84	100.70	108.09
4	C	601	STU	C6-C7-C8	-7.63	119.76	130.53
4	A	601	STU	C6-C7-C8	-7.50	119.94	130.53
4	C	601	STU	C18-C19-N3	-7.22	123.29	129.42
4	A	601	STU	C12-C17-N2	-7.17	104.22	108.94
4	C	601	STU	C12-C17-N2	-6.98	104.35	108.94
4	A	601	STU	C18-C19-N3	-6.74	123.70	129.42
4	A	601	STU	C9-N1-C8	-6.45	107.98	113.86
5	C	602	C1V	CAU-CAX-NAO	6.40	120.59	114.81
5	A	602	C1V	CAU-CAX-NAO	6.28	120.49	114.81
4	C	601	STU	C9-N1-C8	-6.16	108.24	113.86
4	A	601	STU	C16-C17-N2	5.41	138.35	130.22
4	C	601	STU	C16-C17-N2	5.36	138.26	130.22
6	E	401	C2Z	OAA-CAK-CAG	-4.82	123.34	132.71
6	F	401	C2Z	OAA-CAK-CAG	-4.73	123.50	132.71
6	E	402	C2Z	OAA-CAK-CAG	-4.63	123.71	132.71
6	F	402	C2Z	OAA-CAK-CAG	-4.53	123.89	132.71
4	C	601	STU	C6-C7-C10	-4.47	117.65	120.76
4	C	601	STU	C28-N4-C23	4.30	119.53	114.39
4	A	601	STU	C28-N4-C23	4.22	119.42	114.39
4	C	601	STU	C16-C17-C12	-4.14	116.75	121.59
4	A	601	STU	C16-C17-C12	-4.13	116.76	121.59
4	A	601	STU	C11-C10-C7	-3.90	118.68	121.25
4	C	601	STU	C1-C20-C5	-3.89	117.04	121.59
4	A	601	STU	C6-C7-C10	-3.80	118.11	120.76
4	A	601	STU	C1-C20-C5	-3.73	117.22	121.59
6	F	401	C2Z	OAB-PAO-OAC	-3.72	101.38	112.52
5	A	602	C1V	CAE-CAU-CAX	3.72	121.02	115.79
6	F	402	C2Z	OAB-PAO-OAC	-3.63	101.66	112.52
4	C	601	STU	C11-C10-C7	-3.42	119.00	121.25
4	C	601	STU	C9-C10-C11	-3.40	126.14	128.73
6	E	401	C2Z	OAB-PAO-OAC	-3.40	102.35	112.52
6	E	402	C2Z	CAN-OAJ-CAM	-3.31	104.93	106.57
4	A	601	STU	C18-N2-C17	3.28	111.33	107.37
5	C	602	C1V	CAZ-CAY-SAP	-3.26	108.13	112.74
5	A	602	C1V	CAZ-CAY-SAP	-3.24	108.15	112.74
4	C	601	STU	C1-C20-N3	3.24	135.24	129.73
4	C	601	STU	C18-N2-C17	3.23	111.27	107.37
4	A	601	STU	C1-C20-N3	3.16	135.11	129.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	601	STU	C11-C18-N2	-3.03	106.10	109.38
4	A	601	STU	C11-C18-N2	-3.01	106.13	109.38
4	A	601	STU	C9-C10-C11	-3.00	126.45	128.73
4	A	601	STU	C5-C6-C19	-2.98	102.25	106.44
5	C	602	C1V	CAE-CAU-CAX	2.93	119.91	115.79
6	F	401	C2Z	CAN-OAJ-CAM	-2.93	105.12	106.57
4	C	601	STU	C12-C11-C10	-2.82	127.21	132.33
4	C	601	STU	C5-C6-C19	-2.75	102.58	106.44
4	C	601	STU	C4-C5-C20	2.62	122.27	119.24
4	C	601	STU	C13-C12-C11	-2.58	127.35	133.49
4	A	601	STU	C13-C12-C11	-2.58	127.36	133.49
6	F	402	C2Z	CAN-OAJ-CAM	-2.56	105.30	106.57
4	A	601	STU	C12-C11-C10	-2.48	127.83	132.33
5	A	602	C1V	CAW-CAN-SAP	-2.41	107.19	114.23
6	E	401	C2Z	CAN-OAJ-CAM	-2.37	105.40	106.57
5	C	602	C1V	CAW-CAN-SAP	-2.36	107.32	114.23
4	C	601	STU	C6-C19-C18	-2.33	115.67	121.25
4	A	601	STU	C6-C19-C18	-2.26	115.83	121.25
6	E	402	C2Z	CAG-CAL-NAH	-2.23	110.25	112.11
4	A	601	STU	C4-C5-C20	2.21	121.79	119.24
4	A	601	STU	C20-N3-C25	-2.14	121.97	127.45
6	F	402	C2Z	CAG-CAL-NAH	-2.11	110.35	112.11
6	F	401	C2Z	CAE-CAM-CAL	-2.09	125.94	132.29
4	A	601	STU	C13-C12-C17	2.09	121.65	119.24
4	C	601	STU	C20-N3-C25	-2.04	122.22	127.45
4	A	601	STU	C4-C5-C6	-2.04	128.63	133.49
6	E	401	C2Z	OAI-CAK-CAG	-2.01	109.77	118.17
5	A	602	C1V	OAD-CAT-CAU	-2.01	119.49	122.81
6	F	401	C2Z	OAI-CAK-CAG	-2.01	109.79	118.17

There are no chirality outliers.

All (4) torsion outliers are listed below:

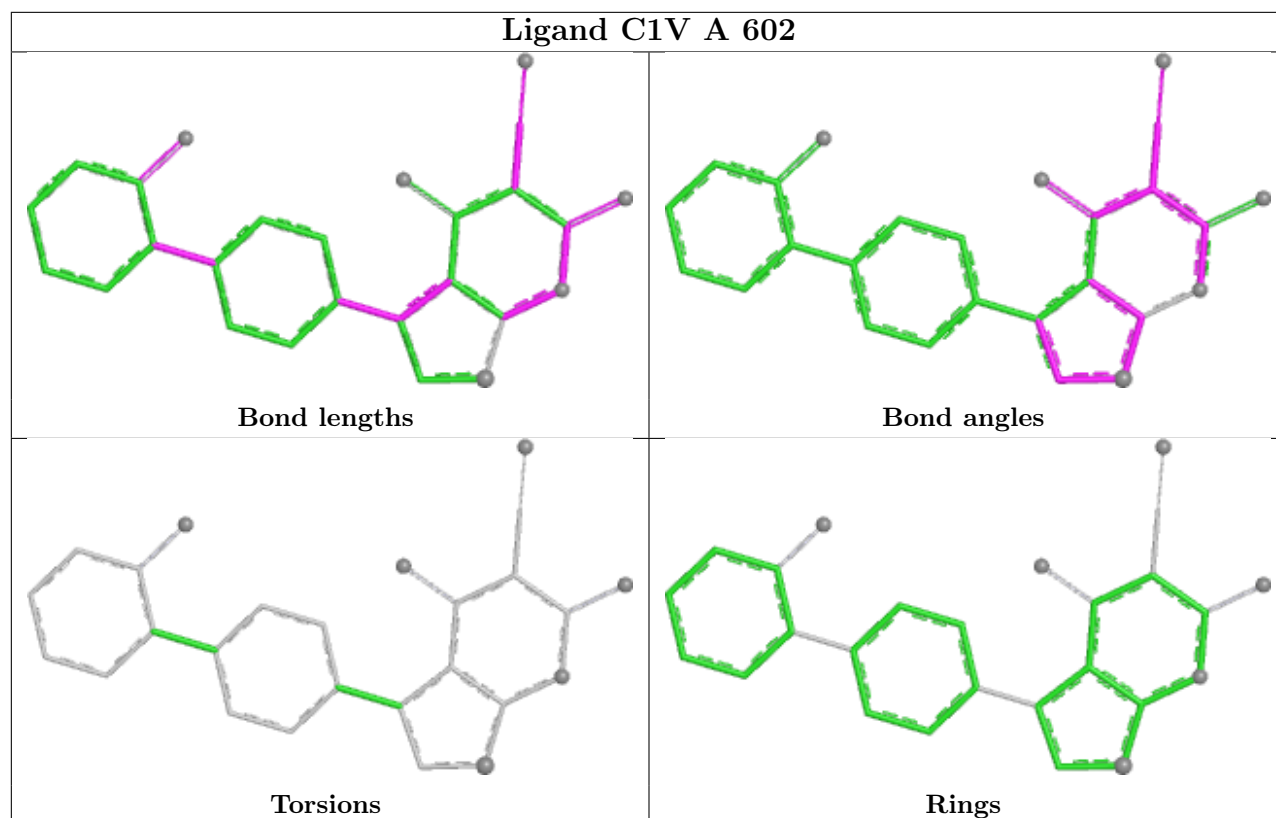
Mol	Chain	Res	Type	Atoms
4	C	601	STU	C22-C23-N4-C28
4	A	601	STU	C24-C23-N4-C28
4	C	601	STU	C24-C23-N4-C28
4	A	601	STU	C22-C23-N4-C28

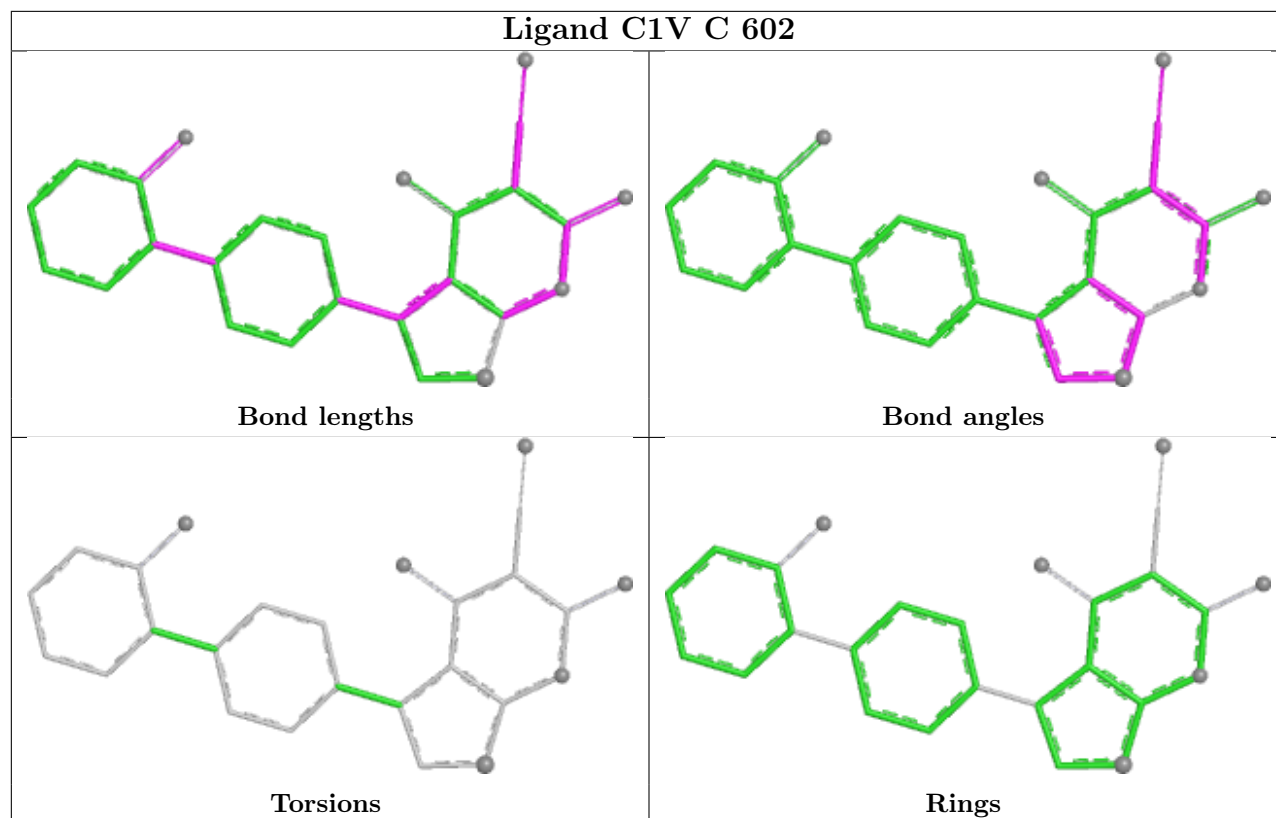
There are no ring outliers.

3 monomers are involved in 6 short contacts:

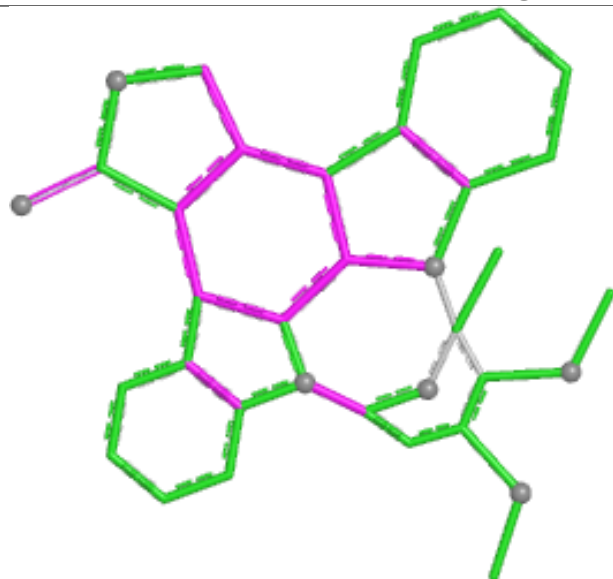
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	602	C1V	1	0
4	C	601	STU	3	0
4	A	601	STU	2	0

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

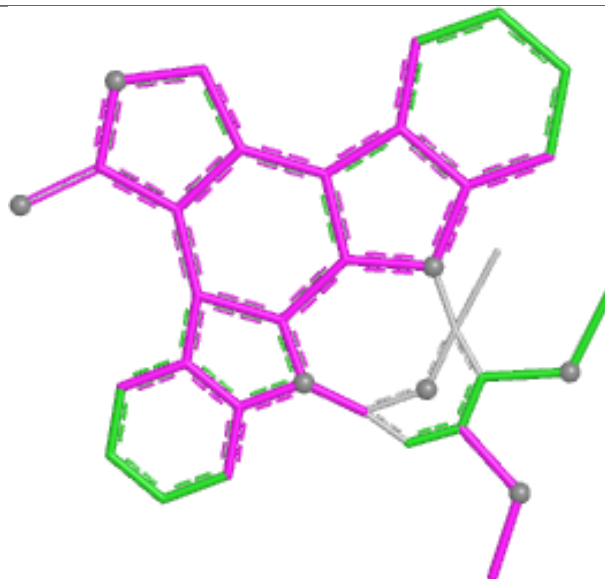




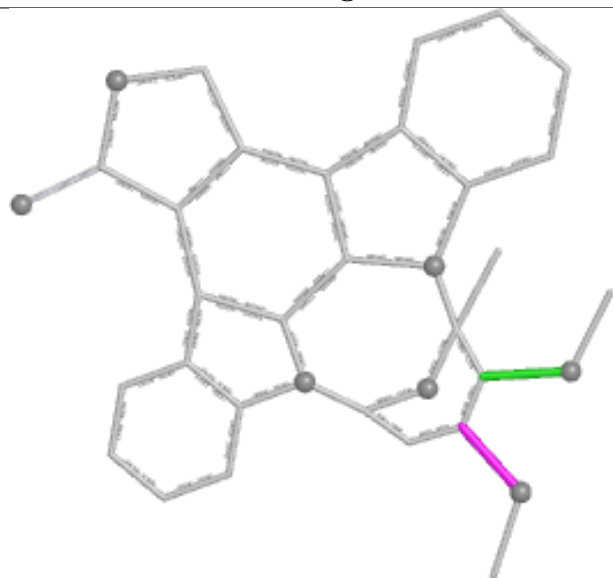
Ligand STU C 601



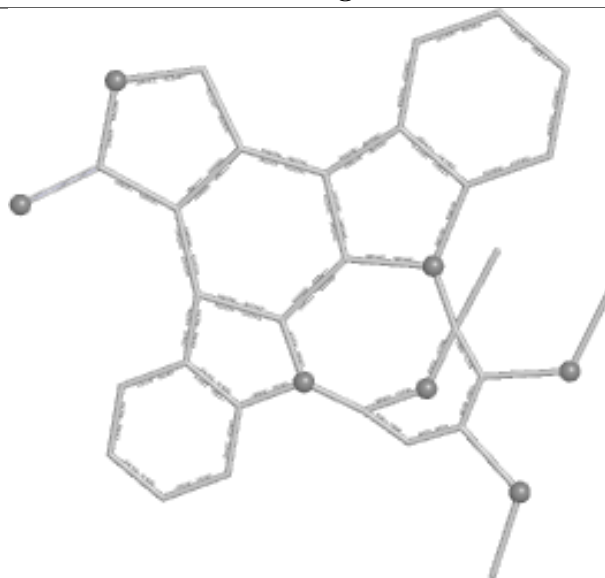
Bond lengths



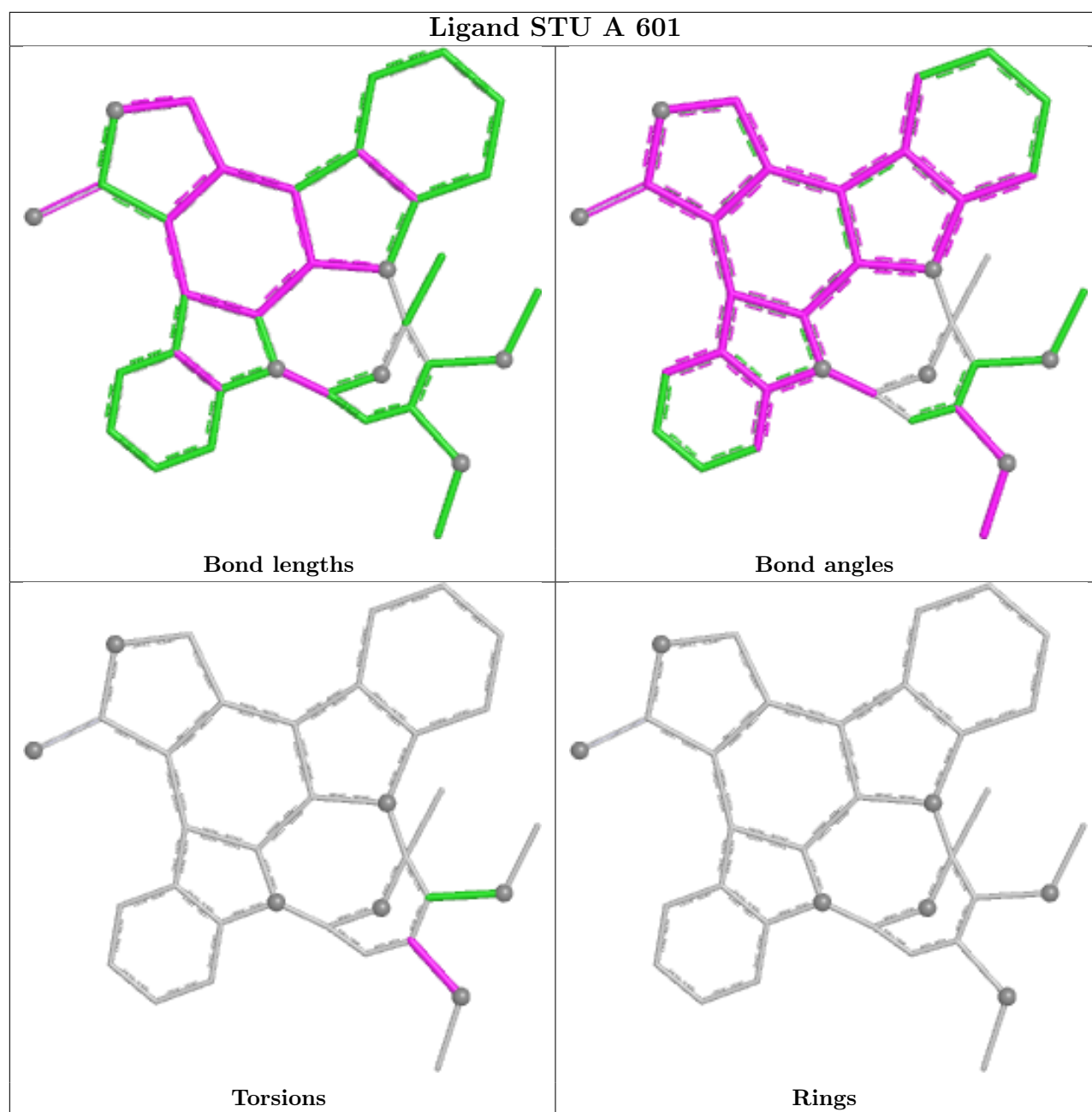
Bond angles



Torsions



Rings



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2		OWAB(Å ²)	Q < 0.9
1	A	441/560 (78%)	0.76	36 (8%)	17 9	40, 63, 98, 142	0
1	C	440/560 (78%)	0.84	45 (10%)	12 6	38, 67, 105, 126	0
2	B	184/270 (68%)	0.92	11 (5%)	27 14	42, 71, 109, 131	0
2	D	185/270 (68%)	0.88	13 (7%)	22 12	51, 71, 115, 146	0
3	E	302/336 (89%)	0.75	21 (6%)	22 12	40, 65, 93, 113	0
3	F	300/336 (89%)	0.69	21 (7%)	22 12	43, 68, 93, 109	1 (0%)
All	All	1852/2332 (79%)	0.79	147 (7%)	18 10	38, 67, 103, 146	1 (0%)

All (147) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	314	GLY	8.9
3	F	324	LEU	5.9
2	B	176	VAL	5.0
1	A	315	ASP	4.9
1	C	364	VAL	4.6
3	F	304	ASP	4.6
1	C	319	GLN	4.5
1	C	316	PRO	4.4
3	E	326	GLY	4.4
1	A	361	PRO	4.3
1	A	366	PHE	4.2
1	A	410	MET	4.1
1	A	156	ALA	3.9
2	D	179	LEU	3.9
1	A	300	PHE	3.8
1	A	468	ASP	3.7
1	A	344	SER	3.6
1	A	316	PRO	3.6
2	D	195	LYS	3.6

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Mol	Chain	Res	Type	RSRZ
3	E	274	GLU	3.4
3	E	276	VAL	3.4
1	C	361	PRO	3.4
1	A	359	PRO	3.3
1	A	365	PRO	3.3
1	C	469	GLU	3.3
3	F	25	ASN	3.2
3	E	272	TYR	3.2
1	A	304	GLU	3.2
1	A	426	VAL	3.2
3	E	268	HIS	3.1
1	A	364	VAL	3.1
1	C	300	PHE	3.1
3	F	273	PHE	3.1
1	C	292	ALA	3.1
2	D	186	PRO	3.1
3	E	273	PHE	3.0
1	A	362	GLU	3.0
2	B	196	PRO	3.0
1	C	174	CYS	2.9
1	A	320	LEU	2.9
1	C	320	LEU	2.9
3	F	117	TRP	2.9
1	C	359	PRO	2.9
1	C	294	LYS	2.9
1	A	323	ALA	2.9
1	C	271	ASP	2.8
3	E	304	ASP	2.8
2	B	219	THR	2.8
2	D	201	ARG	2.8
3	F	306	ASN	2.8
2	D	198	GLU	2.7
2	D	200	PHE	2.7
1	C	288	ILE	2.7
1	C	461	LEU	2.7
2	B	270	ILE	2.7
2	B	190	GLU	2.7
1	A	363	ARG	2.7
1	C	156	ALA	2.7
3	E	279	CYS	2.7
3	E	25	ASN	2.7
3	E	190	SER	2.7

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Mol	Chain	Res	Type	RSRZ
2	B	191	PRO	2.6
1	A	299	LYS	2.6
1	C	446	LYS	2.6
3	F	277	LEU	2.6
3	E	271	HIS	2.6
3	F	125	SER	2.6
1	A	360	HIS	2.6
1	A	13	ILE	2.6
1	A	290	ASP	2.6
1	A	312	TYR	2.6
2	D	253	ALA	2.6
1	C	398	LEU	2.5
3	E	244	PHE	2.5
3	E	206	MET	2.5
3	F	322	LEU	2.5
1	C	426	VAL	2.5
3	F	303	VAL	2.5
3	F	26	ASN	2.5
1	A	427	VAL	2.5
3	E	210	THR	2.5
1	C	302	CYS	2.5
2	B	189	GLN	2.4
1	C	454	VAL	2.4
2	D	187	TYR	2.4
1	C	526	LEU	2.4
1	A	415	ARG	2.4
1	C	436	ARG	2.4
1	C	358	ARG	2.3
1	C	439	PRO	2.3
2	D	199	ARG	2.3
1	A	397	HIS	2.3
1	A	192	GLY	2.3
1	C	167	GLY	2.3
2	B	131	GLY	2.3
1	C	447	MET	2.3
1	C	400	ILE	2.3
3	F	187	GLU	2.3
2	B	223	CYS	2.3
1	C	444	TYR	2.3
3	F	122	LEU	2.2
1	C	524	PRO	2.2
3	F	276	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	215	LEU	2.2
1	C	455	ASP	2.2
1	C	311	LEU	2.2
1	C	416	ALA	2.2
1	C	437	LYS	2.2
3	F	127	LYS	2.2
3	E	122	LEU	2.2
3	E	269	ARG	2.2
1	A	317	GLN	2.2
1	C	424	TRP	2.2
1	C	425	LYS	2.2
3	F	244	PHE	2.2
1	A	394	ALA	2.2
1	C	538	ALA	2.2
3	E	307	ASP	2.1
2	B	187	TYR	2.1
2	D	197	GLU	2.1
1	C	434	VAL	2.1
3	F	309	VAL	2.1
1	C	290	ASP	2.1
3	E	232	ASP	2.1
3	E	126	PHE	2.1
1	A	545	ALA	2.1
1	C	345	SER	2.1
3	E	180	ILE	2.1
1	C	124	LEU	2.1
1	A	275	TYR	2.1
1	C	452	TYR	2.1
3	F	54	SER	2.1
3	F	323	VAL	2.1
1	A	212	LEU	2.1
3	E	188	PHE	2.1
3	F	118	ARG	2.1
3	F	75	TRP	2.1
1	A	434	VAL	2.0
1	C	231	PHE	2.0
1	C	303	THR	2.0
2	D	190	GLU	2.0
1	C	313	SER	2.0
2	D	191	PRO	2.0
2	D	215	LEU	2.0
1	A	393	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	423	GLU	2.0
1	C	291	GLU	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	TPO	A	172	11/12	0.86	0.12	56,57,60,61	1
2	SEP	D	108	10/11	0.89	0.10	65,65,69,69	0
1	TPO	C	172	11/12	0.90	0.09	69,69,73,73	2
2	SEP	B	108	10/11	0.93	0.09	64,64,67,67	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

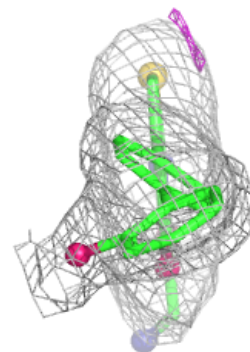
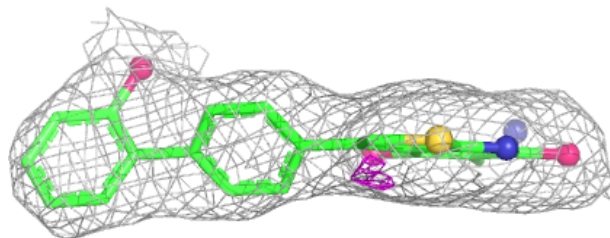
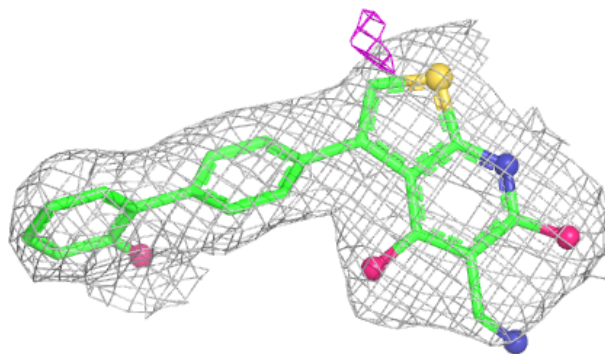
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	C2Z	E	401	15/15	0.80	0.17	100,100,101,101	0
6	C2Z	E	402	15/15	0.80	0.17	92,92,93,93	0
6	C2Z	F	401	15/15	0.84	0.14	92,93,94,94	0
6	C2Z	F	402	15/15	0.84	0.15	88,88,88,88	0
5	C1V	A	602	26/26	0.92	0.11	44,45,47,48	0
4	STU	C	601	35/35	0.94	0.10	39,40,42,43	0
4	STU	A	601	35/35	0.95	0.09	38,39,40,40	0
5	C1V	C	602	26/26	0.95	0.08	42,44,45,45	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.

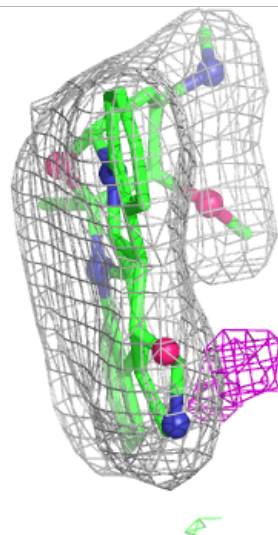
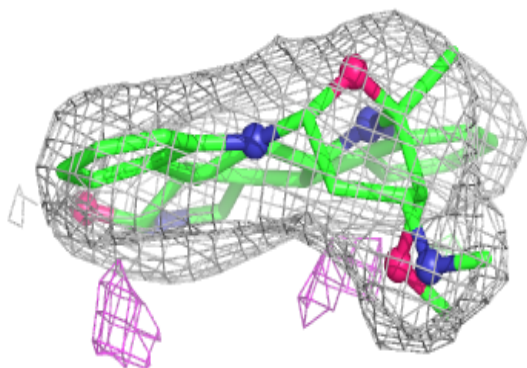
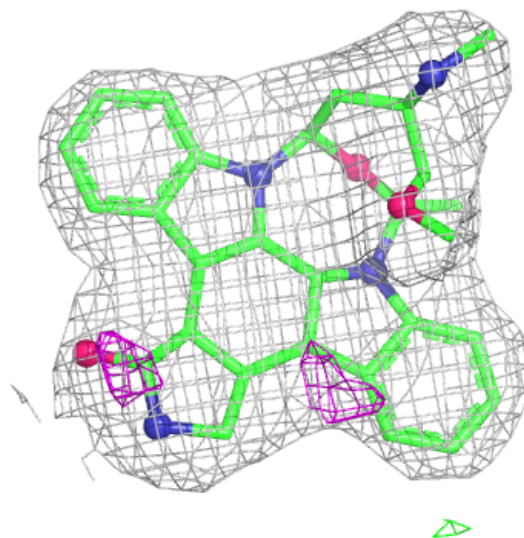
Electron density around C1V A 602:

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



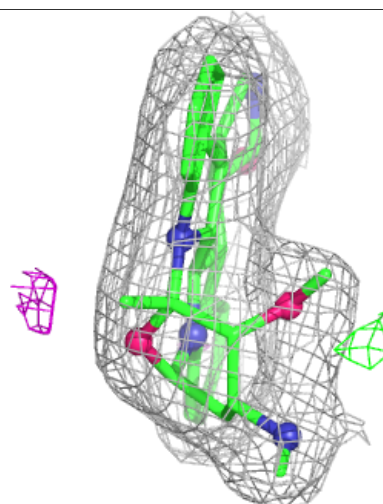
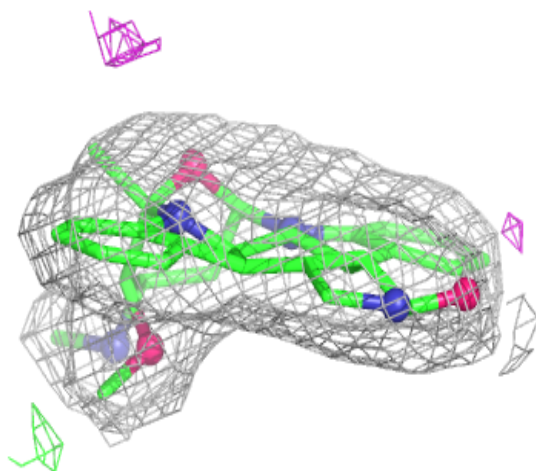
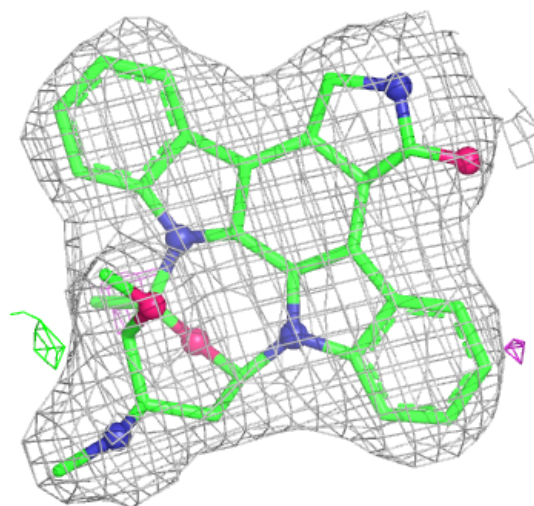
Electron density around STU C 601:

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



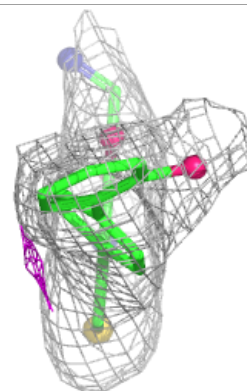
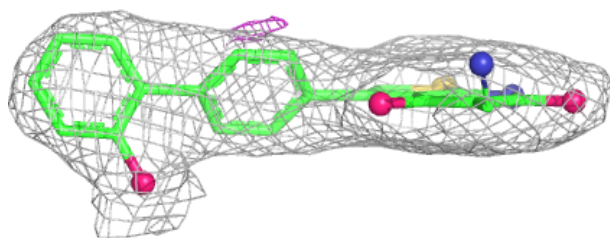
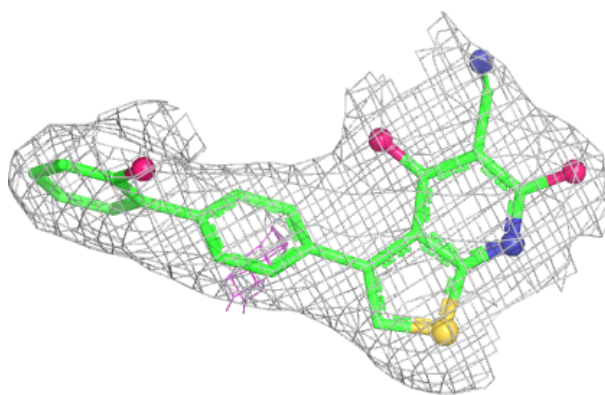
Electron density around STU A 601:

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



Electron density around C1V C 602:

$2mF_o-DF_c$ (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.