



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 08:22 PM UTC

PDB ID : 9J2Z / pdb_00009j2z
Title : mouse cGAS catalytic domain bound with RU.521
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Deposited on : 2024-08-07
Resolution : 2.39 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

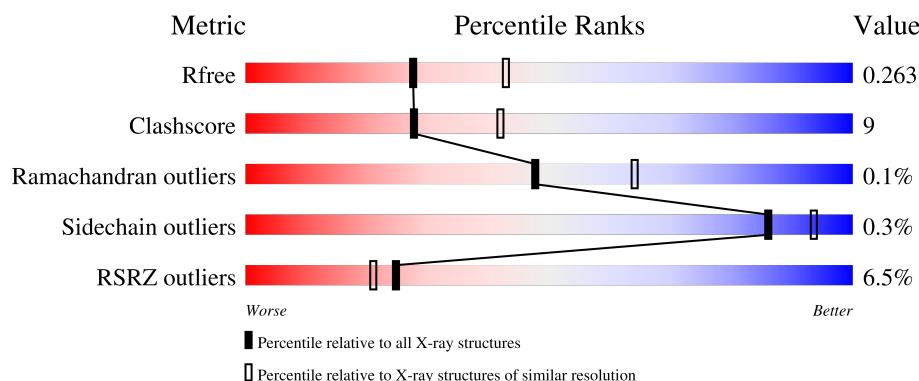
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.39 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	362	3% 86% 13% .
1	B	362	6% 82% 15% .
1	C	362	7% 81% 17% .
1	D	362	10% 75% 22% .

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
3	A1EAM	D	601	X	-	-	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11317 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 Cyclic GMP-AMP synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	360	Total	C	N	O	S	0	0	0
			2907	1869	487	538	13			
1	B	352	Total	C	N	O	S	0	0	0
			2822	1813	469	527	13			
1	C	356	Total	C	N	O	S	0	0	0
			2803	1794	478	518	13			
1	D	351	Total	C	N	O	S	0	0	0
			2697	1730	447	508	12			

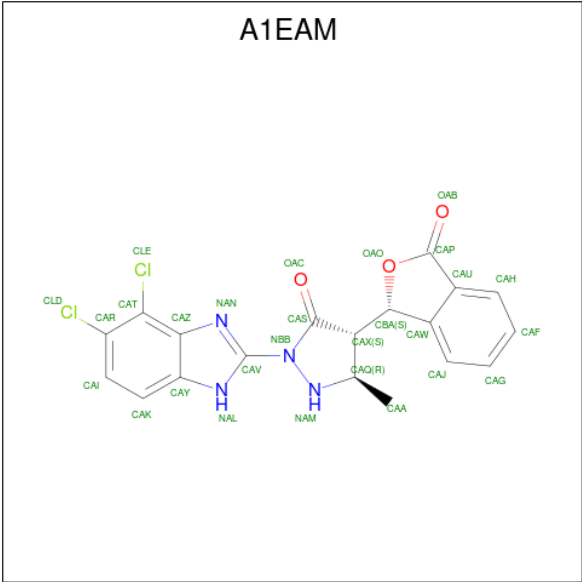
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	146	SER	GLU	conflict	UNP Q8C6L5
B	146	SER	GLU	conflict	UNP Q8C6L5
C	146	SER	GLU	conflict	UNP Q8C6L5
D	146	SER	GLU	conflict	UNP Q8C6L5

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		

- Molecule 3 is (4 {S},5 {R})-2-[4,5-bis(chloranyl)-1 {H}-benzimidazol-2-yl]-5-methyl-4-[(1 {S})-3-oxidanylidene-1 {H}-2-benzofuran-1-yl]pyrazolidin-3-one (CCD ID: A1EAM) (formula: C₁₉H₁₄Cl₂N₄O₃) (labeled as "Ligand of Interest" by depositor).

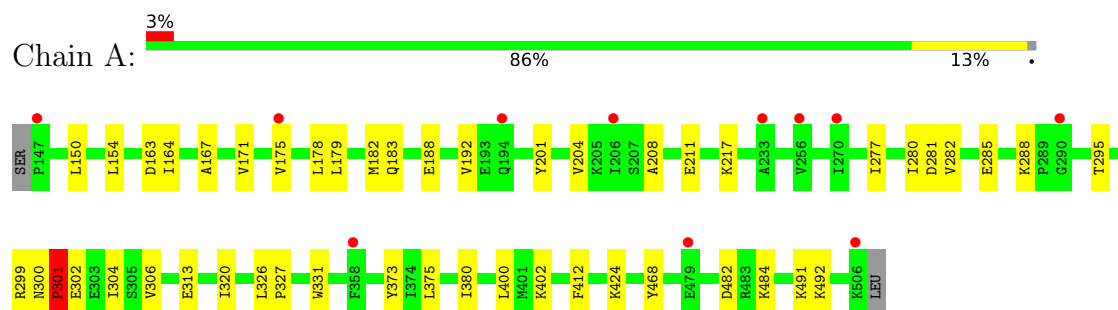


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	Cl	N	O	0	0
			28	19	2	4	3		
3	C	1	Total	C	Cl	N	O	0	0
			28	19	2	4	3		
3	D	1	Total	C	Cl	N	O	0	0
			28	19	2	4	3		

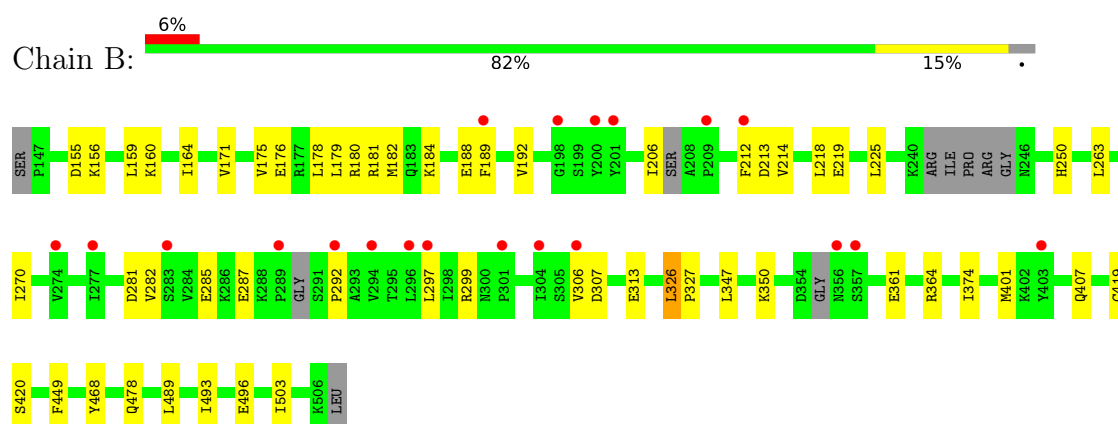
3 Residue-property plots [i](#)

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.

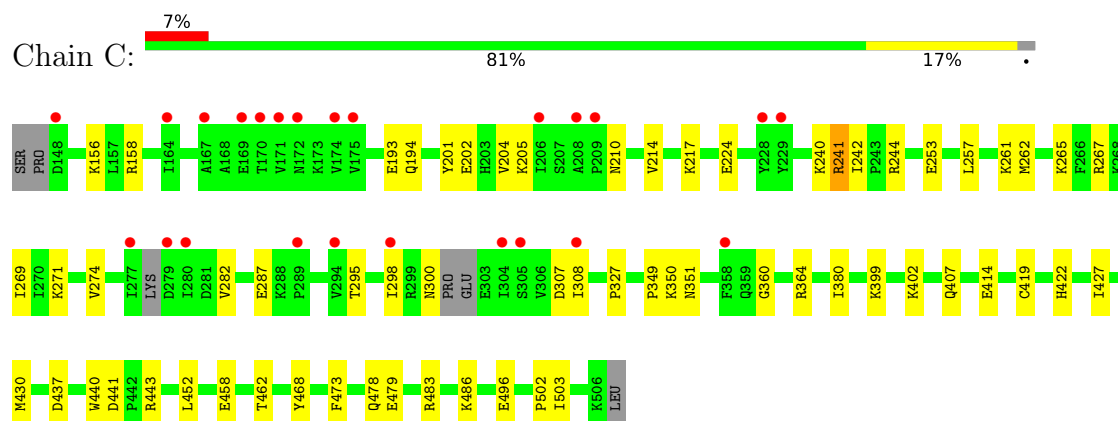
• Molecule 1: Cyclic GMP-AMP synthase



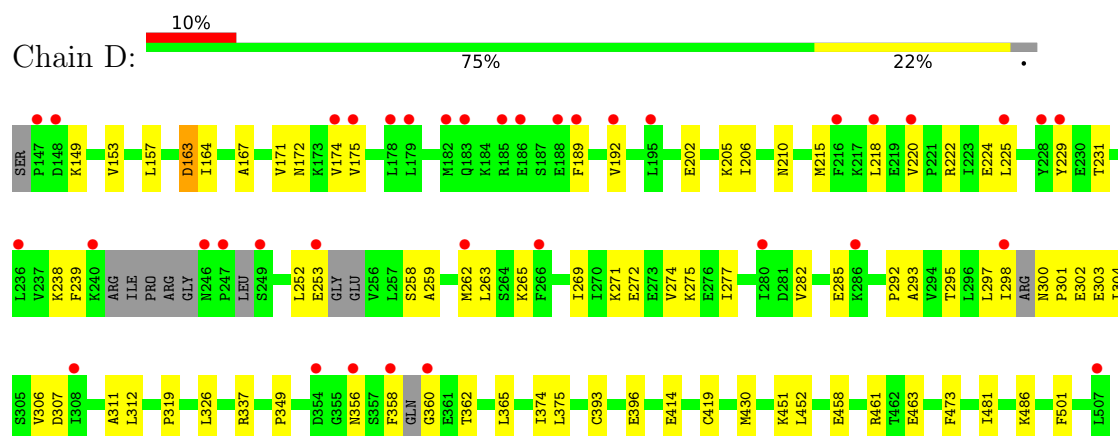
• Molecule 1: Cyclic GMP-AMP synthase



• Molecule 1: Cyclic GMP-AMP synthase



● Molecule 1: Cyclic GMP-AMP synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	86.75Å 84.38Å 124.68Å 90.00° 92.52° 90.00°	Depositor
Resolution (Å)	44.06 – 2.39 44.06 – 2.39	Depositor EDS
% Data completeness (in resolution range)	97.3 (44.06-2.39) 97.3 (44.06-2.39)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 2.39Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.222 , 0.263 0.222 , 0.263	Depositor DCC
R_{free} test set	3332 reflections (4.68%)	wwPDB-VP
Wilson B-factor (Å ²)	55.6	Xtriage
Anisotropy	0.236	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 29.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.009 for k,h,-l 0.004 for -k,-h,-l 0.021 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11317	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 4.48% 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: A1EAM, ZN

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	0.74	2/2972 (0.1%)	0.51	4/4004 (0.1%)
1	B	0.32	1/2881 (0.0%)	0.45	0/3883
1	C	0.40	1/2861 (0.0%)	0.51	1/3855 (0.0%)
1	D	0.44	3/2753 (0.1%)	0.57	1/3726 (0.0%)
All	All	0.50	7/11467 (0.1%)	0.51	6/15468 (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	C	0	2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	326	LEU	C-O	-36.52	1.05	1.23
1	C	419	CYS	CB-SG	-10.18	1.47	1.81
1	D	262	MET	SD-CE	-8.67	1.57	1.79
1	B	326	LEU	C-O	7.24	1.27	1.23
1	A	326	LEU	CA-C	7.04	1.60	1.53
1	D	501	PHE	C-O	5.87	1.26	1.23
1	D	419	CYS	CB-SG	-5.02	1.64	1.81

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	301	PRO	N-CA-CB	-12.09	90.55	103.25
1	A	326	LEU	O-C-N	8.18	125.29	121.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	419	CYS	CA-CB-SG	7.28	131.15	114.40
1	A	326	LEU	CB-CA-C	-6.58	106.53	111.20
1	A	301	PRO	N-CD-CG	-6.10	94.05	103.20
1	D	356	ASN	N-CA-C	-5.71	105.42	112.90

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	158	ARG	Sidechain
1	C	241	ARG	Sidechain

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	2907	0	2872	39	0
1	B	2822	0	2740	41	1
1	C	2803	0	2698	53	1
1	D	2697	0	2513	62	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	B	28	0	0	4	0
3	C	28	0	0	5	0
3	D	28	0	0	3	0
All	All	11317	0	10823	199	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 9.

All (199) 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)
3:B:601:A1EAM:CBA	3:B:601:A1EAM:OAO	1.64	1.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:601:A1EAM:OAO	3:D:601:A1EAM:CBA	1.65	1.36
1:D:430:MET:HE1	1:D:451:LYS:HG3	1.29	1.14
1:B:478:GLN:N	1:B:478:GLN:OE1	2.01	0.92
1:C:257:LEU:HD21	1:C:262:MET:HE1	1.54	0.90
1:A:277:ILE:HD11	1:A:282:VAL:HB	1.55	0.87
1:C:473:PHE:HE2	3:C:601:A1EAM:CLE	1.96	0.85
1:C:253:GLU:OE2	1:C:261:LYS:HE2	1.77	0.84
1:C:473:PHE:CE2	3:C:601:A1EAM:CLE	2.68	0.84
1:A:484:LYS:N	1:A:484:LYS:HD3	1.93	0.83
1:C:414:GLU:OE1	1:C:414:GLU:N	2.10	0.82
1:C:224:GLU:HB2	1:C:240:LYS:HG3	1.64	0.79
1:D:172:ASN:HA	1:D:175:VAL:CG1	2.13	0.79
1:D:285:GLU:HB3	1:D:297:LEU:HD11	1.65	0.79
1:D:430:MET:CE	1:D:451:LYS:HG3	2.10	0.79
1:B:188:GLU:OE2	1:B:250:HIS:NE2	2.16	0.77
1:C:458:GLU:O	1:C:462:THR:HG22	1.85	0.76
1:C:265:LYS:O	1:C:269:ILE:HG13	1.85	0.75
1:D:167:ALA:O	1:D:171:VAL:HG23	1.86	0.75
1:D:302:GLU:N	1:D:302:GLU:OE1	2.20	0.75
1:D:263:LEU:HD21	1:D:349:PRO:HG2	1.70	0.73
1:C:479:GLU:OE1	1:C:479:GLU:N	2.16	0.73
1:C:349:PRO:O	1:C:350:LYS:HG2	1.89	0.72
1:D:414:GLU:OE1	1:D:414:GLU:N	2.18	0.72
1:B:287:GLU:O	1:B:287:GLU:CG	2.40	0.70
1:A:182:MET:HE2	1:A:192:VAL:HG11	1.72	0.69
1:A:208:ALA:HB3	1:A:211:GLU:HG3	1.75	0.69
1:B:287:GLU:O	1:B:287:GLU:HG2	1.92	0.69
1:C:201:TYR:HA	1:C:204:VAL:CG1	2.24	0.68
1:D:300:ASN:HB3	1:D:301:PRO:HD3	1.73	0.68
1:C:351:ASN:OD1	1:C:360:GLY:N	2.22	0.67
3:D:601:A1EAM:OAO	3:D:601:A1EAM:CAX	2.42	0.66
1:A:179:LEU:O	1:A:183:GLN:HG3	1.96	0.66
1:A:484:LYS:HD3	1:A:484:LYS:H	1.59	0.66
1:C:496:GLU:OE1	1:C:502:PRO:HD2	1.96	0.66
1:D:463:GLU:OE1	1:D:486:LYS:NZ	2.29	0.65
1:D:297:LEU:HD23	1:D:303:GLU:OE1	1.97	0.64
1:D:224:GLU:C	1:D:225:LEU:HD12	2.22	0.64
1:B:178:LEU:HD12	1:B:178:LEU:O	1.96	0.63
1:D:172:ASN:HA	1:D:175:VAL:HG12	1.80	0.63
1:D:171:VAL:O	1:D:175:VAL:HG12	1.98	0.63
3:B:601:A1EAM:OAO	3:B:601:A1EAM:CAX	2.41	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:430:MET:HE1	1:D:451:LYS:CG	2.19	0.63
1:A:285:GLU:CD	1:A:288:LYS:HZ3	2.07	0.63
1:B:178:LEU:HD12	1:B:182:MET:HG3	1.80	0.63
1:C:201:TYR:HA	1:C:204:VAL:HG12	1.81	0.63
1:A:171:VAL:O	1:A:175:VAL:HG13	1.99	0.62
1:D:171:VAL:HG13	1:D:306:VAL:HG21	1.82	0.62
1:D:189:PHE:O	1:D:192:VAL:HG13	2.00	0.61
1:B:285:GLU:OE1	1:B:297:LEU:HD21	2.00	0.61
1:D:215:MET:HE3	1:D:311:ALA:HB2	1.82	0.61
1:C:156:LYS:N	1:C:156:LYS:HD3	2.16	0.61
1:C:271:LYS:O	1:C:274:VAL:HG12	2.00	0.61
1:C:430:MET:HE3	1:C:440:TRP:HH2	1.65	0.61
1:D:153:VAL:O	1:D:157:LEU:HG	2.01	0.61
1:C:204:VAL:HG22	1:C:205:LYS:O	2.00	0.60
1:D:265:LYS:O	1:D:269:ILE:HG13	2.02	0.60
1:B:181:ARG:HA	1:B:184:LYS:HE2	1.84	0.60
1:B:178:LEU:HD21	1:B:270:ILE:HG23	1.85	0.59
1:C:422:HIS:CE1	3:C:601:A1EAM:CAA	2.86	0.59
1:C:267:ARG:NH2	1:C:287:GLU:HB2	2.18	0.59
1:A:204:VAL:O	1:A:402:LYS:HE3	2.03	0.58
1:A:285:GLU:CD	1:A:288:LYS:NZ	2.62	0.58
1:D:210:ASN:HA	1:D:304:ILE:HG23	1.85	0.58
1:D:149:LYS:O	1:D:153:VAL:HG23	2.02	0.58
1:A:178:LEU:O	1:A:182:MET:HG3	2.04	0.58
1:D:220:VAL:HG21	1:D:312:LEU:HD22	1.86	0.58
1:C:430:MET:HE2	1:C:452:LEU:HA	1.86	0.57
1:D:218:LEU:HD22	1:D:220:VAL:HG22	1.86	0.57
1:D:358:PHE:C	1:D:360:GLY:N	2.63	0.57
1:D:271:LYS:O	1:D:274:VAL:HG12	2.02	0.57
1:B:160:LYS:O	1:B:164:ILE:HG12	2.05	0.57
1:B:178:LEU:O	1:B:182:MET:HG3	2.04	0.57
1:D:300:ASN:CB	1:D:301:PRO:HD3	2.36	0.56
1:A:277:ILE:HD11	1:A:282:VAL:CB	2.34	0.56
1:A:320:ILE:HD12	1:A:320:ILE:N	2.21	0.56
1:D:430:MET:HE2	1:D:452:LEU:HA	1.88	0.55
1:A:299:ARG:O	1:A:300:ASN:C	2.49	0.55
1:C:194:GLN:OE1	1:C:214:VAL:HG21	2.07	0.55
1:B:263:LEU:HD13	1:B:292:PRO:HB3	1.88	0.55
1:D:285:GLU:HG2	1:D:295:THR:HB	1.88	0.55
1:B:282:VAL:HA	1:B:297:LEU:O	2.05	0.55
1:D:302:GLU:H	1:D:302:GLU:CD	2.15	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:176:GLU:O	1:B:176:GLU:OE1	2.26	0.54
1:B:175:VAL:O	1:B:179:LEU:HD12	2.07	0.54
1:C:253:GLU:CD	1:C:261:LYS:HE2	2.33	0.54
1:D:229:TYR:HB3	1:D:231:THR:HG23	1.90	0.54
1:A:201:TYR:CE1	1:A:402:LYS:NZ	2.76	0.54
1:C:478:GLN:HE22	1:C:483:ARG:HB2	1.73	0.53
1:D:271:LYS:C	1:D:274:VAL:HG12	2.33	0.53
1:A:217:LYS:NZ	1:A:313:GLU:OE1	2.40	0.53
1:A:182:MET:CE	1:A:192:VAL:HG11	2.38	0.53
1:C:295:THR:HG22	1:C:307:ASP:OD1	2.08	0.53
1:B:419:CYS:HB3	3:B:601:A1EAM:CAJ	2.38	0.53
1:C:257:LEU:CD2	1:C:262:MET:HE1	2.34	0.53
1:C:201:TYR:O	1:C:402:LYS:NZ	2.25	0.52
1:A:280:ILE:HD12	1:A:281:ASP:H	1.75	0.52
1:D:481:ILE:HG22	1:D:486:LYS:HE3	1.91	0.52
1:B:181:ARG:O	1:B:184:LYS:HG2	2.10	0.51
1:D:301:PRO:HD2	1:D:302:GLU:OE1	2.09	0.51
1:A:320:ILE:HD12	1:A:320:ILE:H	1.76	0.51
1:D:302:GLU:N	1:D:302:GLU:CD	2.69	0.51
1:A:484:LYS:N	1:A:484:LYS:CD	2.72	0.51
1:C:242:ILE:HG22	1:C:244:ARG:HG3	1.93	0.51
1:D:461:ARG:NH1	1:D:461:ARG:HG2	2.25	0.51
1:D:473:PHE:CE2	3:D:601:A1EAM:CLE	3.02	0.50
1:C:214:VAL:CG1	1:C:308:ILE:HD12	2.41	0.50
1:C:210:ASN:OD1	1:C:210:ASN:N	2.45	0.49
1:D:274:VAL:CG2	1:D:282:VAL:HG12	2.42	0.49
1:A:300:ASN:O	1:A:302:GLU:N	2.46	0.49
1:D:297:LEU:HD23	1:D:303:GLU:CD	2.37	0.49
1:D:297:LEU:HD12	1:D:297:LEU:N	2.28	0.49
1:C:253:GLU:HG2	1:C:261:LYS:CE	2.42	0.49
1:B:175:VAL:HG21	1:B:214:VAL:HG21	1.95	0.49
1:A:320:ILE:H	1:A:320:ILE:CD1	2.25	0.48
1:C:282:VAL:HG13	1:C:298:ILE:HG12	1.95	0.48
1:B:401:MET:HG2	1:B:449:PHE:HE1	1.79	0.48
1:B:327:PRO:HD2	1:B:468:TYR:CZ	2.47	0.48
1:B:407:GLN:HB3	1:B:503:ILE:HG13	1.95	0.48
1:D:258:SER:O	1:D:259:ALA:C	2.56	0.48
1:D:252:LEU:O	1:D:253:GLU:C	2.57	0.48
1:B:212:PHE:HB2	1:B:306:VAL:HG22	1.96	0.47
1:A:331:TRP:O	1:A:380:ILE:HD12	2.13	0.47
1:C:267:ARG:HH11	1:C:267:ARG:HG3	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:327:PRO:HD2	1:A:468:TYR:CZ	2.49	0.47
1:B:176:GLU:OE1	1:B:180:ARG:NE	2.47	0.47
1:A:167:ALA:HB1	1:A:304:ILE:HD13	1.96	0.47
1:C:479:GLU:H	1:C:479:GLU:CD	2.14	0.47
1:C:253:GLU:CG	1:C:261:LYS:HE2	2.45	0.47
1:B:281:ASP:HB3	1:B:299:ARG:HB2	1.96	0.46
1:D:461:ARG:HG2	1:D:461:ARG:HH11	1.80	0.46
1:A:484:LYS:H	1:A:484:LYS:CD	2.28	0.46
1:C:430:MET:HE3	1:C:440:TRP:CH2	2.47	0.46
1:C:483:ARG:HA	1:C:486:LYS:HE3	1.98	0.46
1:B:206:ILE:HG13	1:B:420:SER:HB3	1.97	0.46
1:C:262:MET:HE2	1:C:262:MET:HB2	1.81	0.46
1:C:441:ASP:OD1	1:C:443:ARG:HG2	2.16	0.46
1:D:319:PRO:HG2	1:D:365:LEU:HD21	1.99	0.45
1:A:154:LEU:HG	1:A:400:LEU:HD21	1.97	0.45
1:B:263:LEU:CD1	1:B:292:PRO:HB3	2.45	0.45
1:A:300:ASN:O	1:A:301:PRO:C	2.59	0.45
1:C:201:TYR:CA	1:C:204:VAL:HG12	2.45	0.45
1:B:225:LEU:HD21	1:B:347:LEU:HD11	1.97	0.45
1:B:178:LEU:HD12	1:B:178:LEU:C	2.42	0.45
1:D:326:LEU:HD11	1:D:374:ILE:HD11	1.99	0.45
1:D:393:CYS:HA	1:D:396:GLU:OE1	2.17	0.45
1:D:300:ASN:HB3	1:D:301:PRO:CD	2.43	0.45
1:D:458:GLU:OE2	1:D:461:ARG:NH2	2.50	0.44
1:B:496:GLU:CD	1:B:503:ILE:HB	2.42	0.44
1:A:491:LYS:HB3	1:A:491:LYS:HE3	1.72	0.44
1:B:326:LEU:HD11	1:B:374:ILE:HD11	1.99	0.44
1:D:202:GLU:CD	1:D:375:LEU:HD12	2.41	0.44
1:A:412:PHE:CZ	1:A:492:LYS:HE3	2.52	0.44
1:A:183:GLN:NE2	1:A:192:VAL:O	2.45	0.44
1:B:213:ASP:OD1	1:B:307:ASP:HB2	2.18	0.44
1:C:202:GLU:OE1	1:C:399:LYS:HE2	2.18	0.44
1:C:327:PRO:HD2	1:C:468:TYR:CZ	2.53	0.44
1:B:171:VAL:HG22	1:B:306:VAL:HG21	1.99	0.44
1:B:219:GLU:HA	1:B:313:GLU:HB3	1.99	0.44
1:D:163:ASP:O	1:D:164:ILE:C	2.60	0.44
1:B:189:PHE:O	1:B:192:VAL:HG13	2.18	0.43
1:D:269:ILE:O	1:D:272:GLU:HB3	2.17	0.43
1:D:206:ILE:HG22	1:D:206:ILE:O	2.18	0.43
1:B:192:VAL:HG12	1:B:218:LEU:HB2	2.01	0.43
1:A:295:THR:HA	1:A:306:VAL:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:441:ASP:CG	1:C:443:ARG:HE	2.27	0.43
1:C:380:ILE:HG22	1:C:437:ASP:CG	2.44	0.43
1:C:364:ARG:HD2	3:C:601:A1EAM:CLE	2.56	0.43
1:D:220:VAL:CG2	1:D:312:LEU:HD22	2.47	0.43
1:D:326:LEU:HD23	1:D:337:ARG:HG3	2.00	0.43
1:D:358:PHE:O	1:D:360:GLY:N	2.52	0.43
1:B:350:LYS:HD2	1:B:364:ARG:NH1	2.33	0.42
1:D:174:VAL:HG21	1:D:298:ILE:HD11	2.00	0.42
1:D:292:PRO:O	1:D:293:ALA:C	2.62	0.42
1:B:364:ARG:HD2	3:B:601:A1EAM:CLE	2.56	0.42
1:C:193:GLU:HB2	1:C:217:LYS:HE3	2.00	0.42
1:C:201:TYR:CE2	1:C:402:LYS:HD2	2.54	0.42
1:C:364:ARG:HD2	3:C:601:A1EAM:CAT	2.50	0.42
1:C:267:ARG:HG3	1:C:267:ARG:NH1	2.34	0.42
1:C:267:ARG:CZ	1:C:287:GLU:HB2	2.49	0.42
1:A:150:LEU:HD22	1:A:400:LEU:HD13	2.00	0.41
1:A:373:TYR:CD1	1:A:373:TYR:C	2.97	0.41
1:D:293:ALA:HB1	1:D:307:ASP:HB3	2.03	0.41
1:B:155:ASP:O	1:B:159:LEU:HG	2.20	0.41
1:A:163:ASP:C	1:A:164:ILE:HD13	2.45	0.41
1:C:407:GLN:HB3	1:C:503:ILE:HG12	2.01	0.41
1:D:238:LYS:O	1:D:239:PHE:CD1	2.74	0.41
1:A:280:ILE:HD12	1:A:281:ASP:N	2.36	0.41
1:C:274:VAL:HG22	1:C:282:VAL:HB	2.03	0.41
1:B:164:ILE:N	1:B:164:ILE:HD13	2.36	0.41
1:C:427:ILE:HA	1:C:430:MET:HB2	2.02	0.41
1:D:205:LYS:C	1:D:206:ILE:HD13	2.45	0.41
1:D:275:LYS:C	1:D:277:ILE:N	2.74	0.41
1:A:375:LEU:HD11	1:A:424:LYS:HD2	2.03	0.40
1:D:461:ARG:HH11	1:D:461:ARG:CG	2.34	0.40
1:B:156:LYS:HA	1:B:159:LEU:HG	2.02	0.40
1:A:482:ASP:HB3	1:A:484:LYS:HE2	2.03	0.40
1:A:188:GLU:H	1:A:188:GLU:CD	2.29	0.40
1:B:489:LEU:HG	1:B:493:ILE:HD12	2.03	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:B:361:GLU:OE1	1:C:241:ARG:NH2[2_645]	1.62	0.58

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	358/362 (99%)	347 (97%)	10 (3%)	1 (0%)	36	50
1	B	342/362 (94%)	330 (96%)	12 (4%)	0	100	100
1	C	350/362 (97%)	343 (98%)	7 (2%)	0	100	100
1	D	339/362 (94%)	322 (95%)	16 (5%)	1 (0%)	36	50
All	All	1389/1448 (96%)	1342 (97%)	45 (3%)	2 (0%)	48	64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	301	PRO
1	D	222	ARG

5.3.2 Protein sidechains [i](#)

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	317/335 (95%)	316 (100%)	1 (0%)	86	93
1	B	305/335 (91%)	305 (100%)	0	100	100
1	C	293/335 (88%)	292 (100%)	1 (0%)	86	93
1	D	275/335 (82%)	273 (99%)	2 (1%)	76	88
All	All	1190/1340 (89%)	1186 (100%)	4 (0%)	86	93

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

Mol	Chain	Res	Type
1	A	301	PRO
1	C	300	ASN
1	D	163	ASP
1	D	362	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	194	GLN
1	A	478	GLN
1	B	210	ASN
1	C	203	HIS
1	C	376	ASN
1	C	478	GLN
1	D	194	GLN
1	D	351	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	A1EAM	B	601	1	30,32,32	5.45	21 (70%)	34,49,49	4.36	13 (38%)
3	A1EAM	D	601	1	30,32,32	5.66	22 (73%)	34,49,49	4.88	15 (44%)
3	A1EAM	C	601	1	30,32,32	5.13	21 (70%)	34,49,49	3.31	15 (44%)

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	A1EAM	D	601	1	1/1/6/7	2/4/36/36	0/5/5/5
3	A1EAM	B	601	1	-	2/4/36/36	0/5/5/5
3	A1EAM	C	601	1	-	3/4/36/36	0/5/5/5

All (64) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	601	A1EAM	CAX-CAQ	-22.41	1.29	1.54
3	D	601	A1EAM	CAX-CAQ	-20.57	1.31	1.54
3	B	601	A1EAM	CAX-CAQ	-20.22	1.31	1.54
3	D	601	A1EAM	OAQ-CBA	14.30	1.65	1.46
3	B	601	A1EAM	OAQ-CBA	13.44	1.64	1.46
3	D	601	A1EAM	CAW-CBA	8.33	1.64	1.51
3	B	601	A1EAM	CAW-CBA	6.48	1.61	1.51
3	B	601	A1EAM	CAY-NAL	-5.78	1.28	1.38
3	C	601	A1EAM	CAY-NAL	-5.70	1.28	1.38
3	D	601	A1EAM	CAY-NAL	-5.50	1.28	1.38
3	C	601	A1EAM	CAY-CAZ	-5.29	1.35	1.40
3	D	601	A1EAM	CAK-CAY	5.22	1.48	1.39
3	C	601	A1EAM	CAK-CAY	5.18	1.48	1.39
3	B	601	A1EAM	CAK-CAY	5.16	1.48	1.39
3	B	601	A1EAM	CAY-CAZ	-5.00	1.35	1.40
3	D	601	A1EAM	CAX-CBA	-4.90	1.45	1.54
3	D	601	A1EAM	CAY-CAZ	-4.88	1.35	1.40
3	C	601	A1EAM	CAX-CAS	-4.72	1.44	1.51
3	B	601	A1EAM	CAV-NBB	4.70	1.53	1.38
3	C	601	A1EAM	OAB-CAP	-4.69	1.09	1.22
3	C	601	A1EAM	CAV-NBB	4.66	1.53	1.38
3	D	601	A1EAM	CAV-NBB	4.61	1.53	1.38
3	B	601	A1EAM	OAB-CAP	-4.44	1.10	1.22
3	D	601	A1EAM	OAB-CAP	-4.43	1.10	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	601	A1EAM	CAX-CBA	-4.27	1.46	1.54
3	B	601	A1EAM	CAZ-NAN	-4.25	1.31	1.39
3	C	601	A1EAM	CAZ-NAN	-4.22	1.31	1.39
3	D	601	A1EAM	CAZ-NAN	-4.17	1.32	1.39
3	C	601	A1EAM	CAW-CBA	4.12	1.57	1.51
3	D	601	A1EAM	CAA-CAQ	3.91	1.60	1.52
3	C	601	A1EAM	OAQ-CBA	3.90	1.51	1.46
3	B	601	A1EAM	CAA-CAQ	3.82	1.60	1.52
3	D	601	A1EAM	OAQ-CAP	3.40	1.40	1.36
3	D	601	A1EAM	CAX-CAS	-3.30	1.46	1.51
3	B	601	A1EAM	CAH-CAU	-3.05	1.35	1.39
3	B	601	A1EAM	OAQ-CAP	2.98	1.40	1.36
3	D	601	A1EAM	CAH-CAU	-2.93	1.35	1.39
3	D	601	A1EAM	CAI-CAR	2.92	1.44	1.38
3	C	601	A1EAM	CAI-CAR	2.83	1.44	1.38
3	C	601	A1EAM	CAH-CAU	-2.80	1.35	1.39
3	B	601	A1EAM	CAI-CAR	2.78	1.44	1.38
3	C	601	A1EAM	CAS-NBB	-2.75	1.34	1.39
3	D	601	A1EAM	CAF-CAH	2.65	1.43	1.38
3	C	601	A1EAM	CAF-CAH	2.64	1.43	1.38
3	B	601	A1EAM	CAF-CAH	2.63	1.43	1.38
3	C	601	A1EAM	OAQ-CAP	2.61	1.39	1.36
3	B	601	A1EAM	CAR-CLD	2.59	1.79	1.73
3	C	601	A1EAM	CAA-CAQ	2.57	1.57	1.52
3	B	601	A1EAM	CAX-CAS	-2.54	1.47	1.51
3	D	601	A1EAM	CAR-CLD	2.52	1.79	1.73
3	C	601	A1EAM	CAR-CLD	2.44	1.79	1.73
3	B	601	A1EAM	OAC-CAS	-2.33	1.18	1.22
3	B	601	A1EAM	CAI-CAK	-2.24	1.35	1.38
3	C	601	A1EAM	CAJ-CAW	2.23	1.42	1.39
3	D	601	A1EAM	OAC-CAS	-2.23	1.18	1.22
3	B	601	A1EAM	CAS-NBB	-2.22	1.35	1.39
3	C	601	A1EAM	CAG-CAF	-2.22	1.33	1.38
3	C	601	A1EAM	OAC-CAS	-2.19	1.18	1.22
3	B	601	A1EAM	CAG-CAF	-2.15	1.33	1.38
3	C	601	A1EAM	CAI-CAK	-2.15	1.35	1.38
3	D	601	A1EAM	CAG-CAF	-2.14	1.33	1.38
3	D	601	A1EAM	CAS-NBB	-2.08	1.36	1.39
3	D	601	A1EAM	CAZ-CAT	2.05	1.43	1.40
3	D	601	A1EAM	CAI-CAK	-2.01	1.35	1.38

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	601	A1EAM	OA0-CBA-CAW	-23.98	89.50	103.69
3	B	601	A1EAM	OA0-CBA-CAW	-20.87	91.34	103.69
3	C	601	A1EAM	OA0-CBA-CAW	-10.67	97.38	103.69
3	D	601	A1EAM	OA0-CAP-OAB	7.53	127.68	121.18
3	B	601	A1EAM	OA0-CAP-OAB	7.43	127.59	121.18
3	C	601	A1EAM	OA0-CAP-OAB	6.12	126.46	121.18
3	C	601	A1EAM	CAX-CAS-NBB	-6.02	102.67	107.48
3	B	601	A1EAM	CBA-CAX-CAQ	6.01	126.21	113.69
3	D	601	A1EAM	CBA-CAX-CAQ	5.40	124.93	113.69
3	C	601	A1EAM	CAW-CAU-CAP	-5.14	106.17	108.44
3	D	601	A1EAM	OA0-CAP-CAU	-5.08	105.38	108.37
3	D	601	A1EAM	NAL-CAV-NAN	-4.83	109.61	114.73
3	C	601	A1EAM	NAL-CAV-NAN	-4.82	109.62	114.73
3	D	601	A1EAM	CAZ-CAY-NAL	4.70	107.81	105.13
3	B	601	A1EAM	CAZ-CAY-NAL	4.69	107.80	105.13
3	C	601	A1EAM	CBA-OA0-CAP	-4.63	108.30	111.00
3	B	601	A1EAM	NAL-CAV-NAN	-4.55	109.91	114.73
3	C	601	A1EAM	CAZ-CAY-NAL	4.48	107.68	105.13
3	B	601	A1EAM	OA0-CAP-CAU	-4.45	105.75	108.37
3	C	601	A1EAM	CAJ-CAW-CAU	-4.38	116.59	120.45
3	C	601	A1EAM	OA0-CAP-CAU	-3.81	106.13	108.37
3	C	601	A1EAM	OAC-CAS-CAX	3.18	131.58	127.51
3	D	601	A1EAM	CAY-CAZ-NAN	-2.93	107.77	110.31
3	C	601	A1EAM	CAT-CAR-CLD	-2.88	117.86	120.51
3	D	601	A1EAM	CBA-OA0-CAP	2.86	112.68	111.00
3	D	601	A1EAM	CAA-CAQ-NAM	-2.79	103.92	111.70
3	B	601	A1EAM	CAA-CAQ-NAM	-2.74	104.05	111.70
3	D	601	A1EAM	CAK-CAY-CAZ	-2.71	120.03	122.57
3	C	601	A1EAM	CAK-CAY-CAZ	-2.66	120.07	122.57
3	D	601	A1EAM	CAT-CAR-CLD	-2.65	118.08	120.51
3	D	601	A1EAM	CAJ-CAW-CAU	-2.62	118.14	120.45
3	B	601	A1EAM	CAK-CAY-CAZ	-2.56	120.16	122.57
3	B	601	A1EAM	CAY-CAZ-NAN	-2.56	108.10	110.31
3	C	601	A1EAM	CAY-CAZ-NAN	-2.52	108.13	110.31
3	D	601	A1EAM	CAJ-CAW-CBA	-2.46	125.52	132.14
3	B	601	A1EAM	CAJ-CAW-CAU	-2.46	118.29	120.45
3	D	601	A1EAM	CAW-CAU-CAP	-2.39	107.38	108.44
3	B	601	A1EAM	CAJ-CAW-CBA	-2.39	125.70	132.14
3	C	601	A1EAM	CAA-CAQ-NAM	-2.34	105.18	111.70
3	B	601	A1EAM	CAW-CAU-CAP	-2.31	107.42	108.44
3	C	601	A1EAM	CBA-CAX-CAQ	2.26	118.40	113.69
3	B	601	A1EAM	CAT-CAR-CLD	-2.23	118.46	120.51
3	D	601	A1EAM	CAX-CAS-NBB	-2.15	105.76	107.48

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	D	601	A1EAM	CAX

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	601	A1EAM	CAS-CAX-CBA-CAW
3	B	601	A1EAM	CAQ-CAX-CBA-CAW
3	C	601	A1EAM	CAS-CAX-CBA-CAW
3	C	601	A1EAM	CAQ-CAX-CBA-CAW
3	C	601	A1EAM	CAS-CAX-CBA-OAO
3	D	601	A1EAM	CAS-CAX-CBA-CAW
3	D	601	A1EAM	CAQ-CAX-CBA-CAW

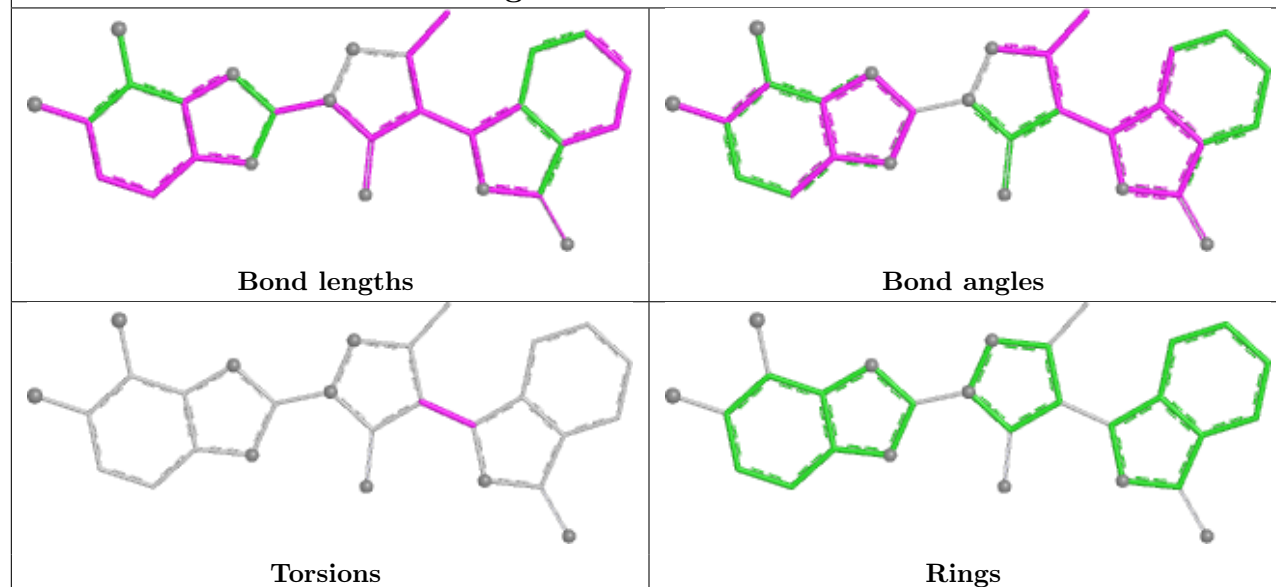
There are no ring outliers.

3 monomers are involved in 12 short contacts:

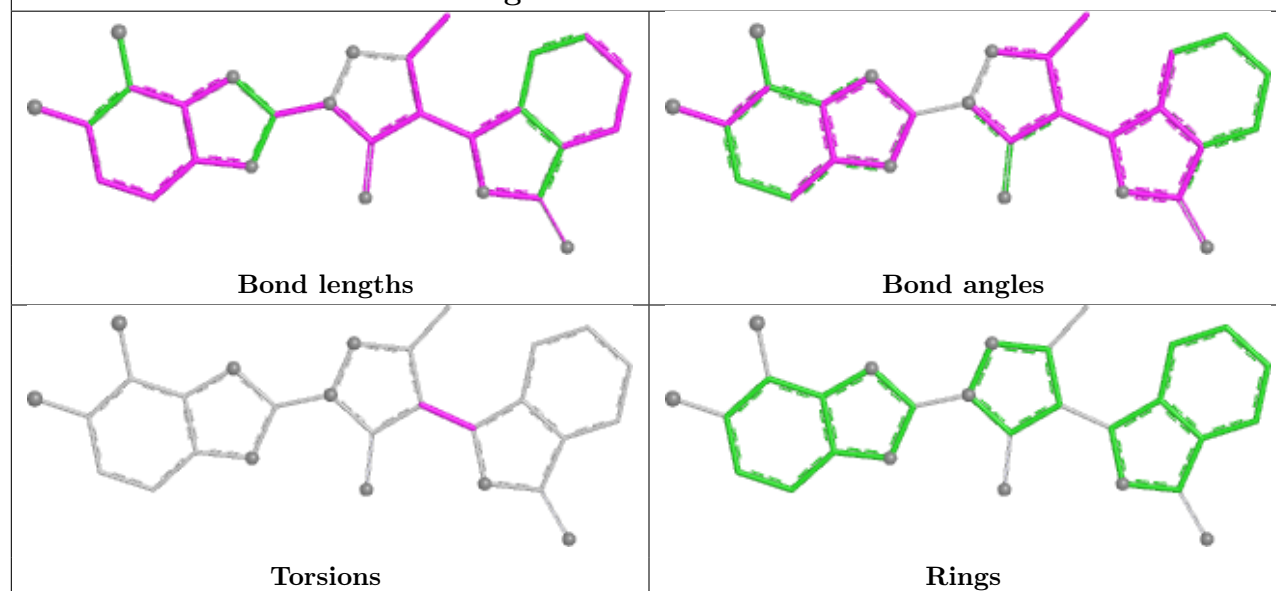
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	601	A1EAM	4	0
3	D	601	A1EAM	3	0
3	C	601	A1EAM	5	0

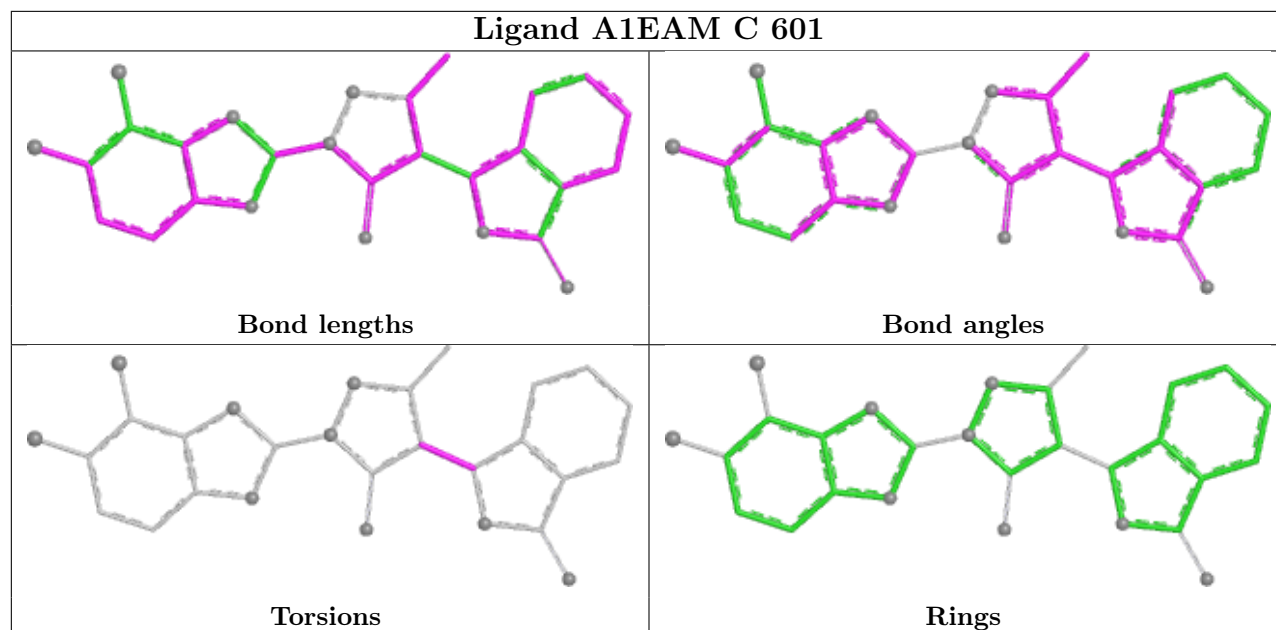
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 A1EAM B 601



Ligand A1EAM D 601





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	360/362 (99%)	0.27	11 (3%) 51 47	43, 62, 90, 103	0
1	B	352/362 (97%)	0.46	20 (5%) 29 25	43, 63, 106, 117	0
1	C	356/362 (98%)	0.53	24 (6%) 24 20	42, 66, 100, 125	0
1	D	351/362 (96%)	0.70	37 (10%) 11 8	39, 75, 111, 131	0
All	All	1419/1448 (97%)	0.49	92 (6%) 25 21	39, 66, 106, 131	0

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	148	ASP	5.5
1	C	304	ILE	5.3
1	D	216	PHE	4.5
1	D	308	ILE	4.4
1	C	279	ASP	4.4
1	D	507	LEU	3.9
1	D	189	PHE	3.8
1	C	280	ILE	3.7
1	C	174	VAL	3.7
1	C	228	TYR	3.5
1	A	506	LYS	3.4
1	B	189	PHE	3.2
1	B	289	PRO	3.2
1	B	277	ILE	3.2
1	A	147	PRO	3.0
1	D	356	ASN	3.0
1	C	289	PRO	3.0
1	B	198	GLY	3.0
1	B	274	VAL	2.9
1	D	354	ASP	2.9
1	D	178	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	218	LEU	2.9
1	C	171	VAL	2.9
1	D	147	PRO	2.8
1	D	186	GLU	2.8
1	D	240	LYS	2.8
1	C	358	PHE	2.8
1	C	209	PRO	2.7
1	B	296	LEU	2.7
1	A	233	ALA	2.7
1	A	358	PHE	2.7
1	B	209	PRO	2.7
1	D	182	MET	2.7
1	C	206	ILE	2.6
1	C	298	ILE	2.6
1	C	305	SER	2.6
1	D	220	VAL	2.6
1	D	358	PHE	2.6
1	C	170	THR	2.6
1	C	294	VAL	2.6
1	C	308	ILE	2.6
1	D	249	SER	2.6
1	B	200	TYR	2.6
1	B	357	SER	2.6
1	C	164	ILE	2.6
1	B	356	ASN	2.5
1	B	294	VAL	2.5
1	C	175	VAL	2.5
1	D	174	VAL	2.5
1	A	270	ILE	2.5
1	C	169	GLU	2.5
1	C	229	TYR	2.5
1	D	185	ARG	2.5
1	D	247	PRO	2.5
1	D	246	ASN	2.5
1	A	175	VAL	2.4
1	D	266	PHE	2.4
1	B	201	TYR	2.4
1	B	292	PRO	2.4
1	B	301	PRO	2.4
1	C	167	ALA	2.4
1	B	306	VAL	2.4
1	D	262	MET	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	194	GLN	2.4
1	C	172	ASN	2.4
1	B	283	SER	2.4
1	B	304	ILE	2.3
1	D	225	LEU	2.3
1	D	360	GLY	2.3
1	A	206	ILE	2.3
1	D	280	ILE	2.3
1	D	175	VAL	2.3
1	D	192	VAL	2.3
1	D	253	GLU	2.3
1	D	195	LEU	2.3
1	A	290	GLY	2.3
1	B	212	PHE	2.2
1	B	297	LEU	2.2
1	B	403	TYR	2.2
1	C	208	ALA	2.2
1	D	183	GLN	2.2
1	D	229	TYR	2.2
1	D	148	ASP	2.2
1	A	479	GLU	2.2
1	C	277	ILE	2.1
1	A	256	VAL	2.1
1	D	228	TYR	2.1
1	D	188	GLU	2.0
1	D	298	ILE	2.0
1	D	179	LEU	2.0
1	D	236	LEU	2.0
1	D	286	LYS	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

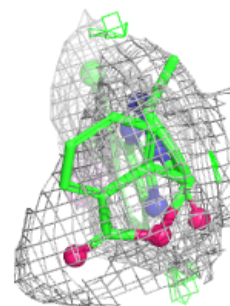
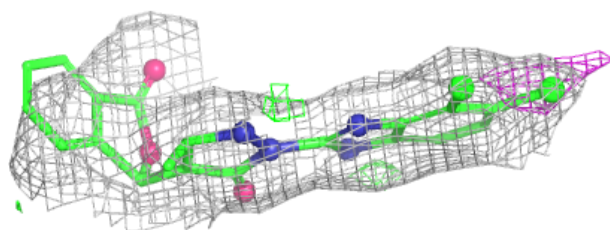
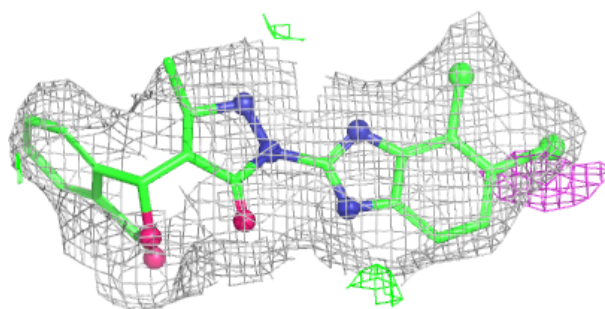
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	A1EAM	D	601	28/28	0.79	0.13	60,74,91,105	0
3	A1EAM	C	601	28/28	0.80	0.13	64,77,93,98	0
3	A1EAM	B	601	28/28	0.84	0.12	57,67,83,87	0
2	ZN	C	602	1/1	0.99	0.03	47,47,47,47	0
2	ZN	B	602	1/1	0.99	0.05	50,50,50,50	0
2	ZN	D	602	1/1	1.00	0.02	47,47,47,47	0
2	ZN	A	601	1/1	1.00	0.04	45,45,45,45	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.

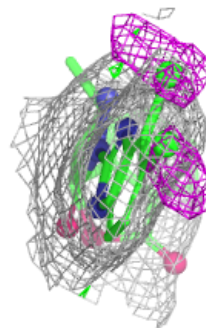
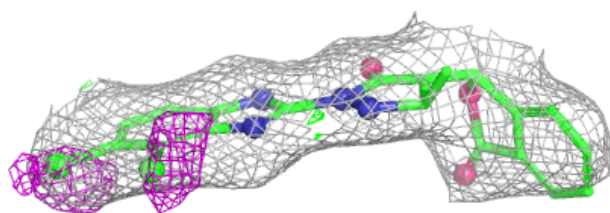
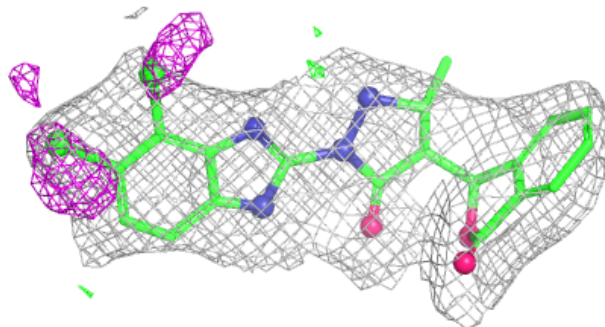
Electron density around A1EAM D 601:

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

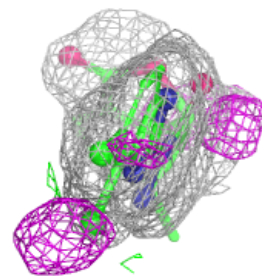
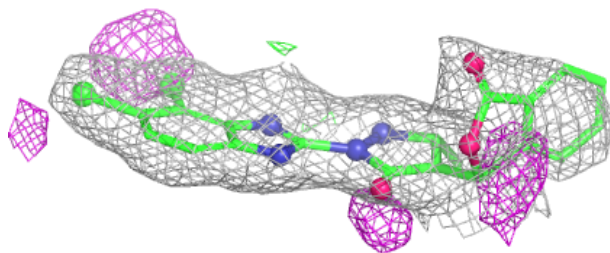
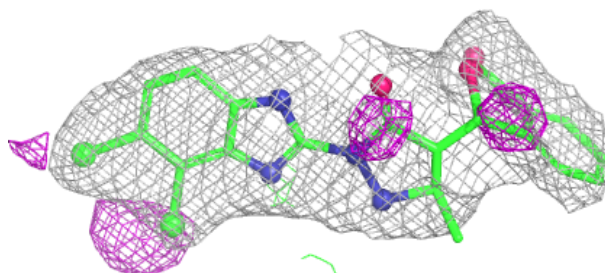


Electron density around A1EAM C 601:

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

**Electron density around A1EAM B 601:**

$2mF_o-DF_c$ (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.