



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:22 AM UTC

PDB ID : 5EML / pdb_00005eml
Title : Crystal structure of PRMT5:MEP50 with Compound 10 and SAM
Authors : Boriack-Sjodin, P.A.; Jin, L.
Deposited on : 2015-11-06
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

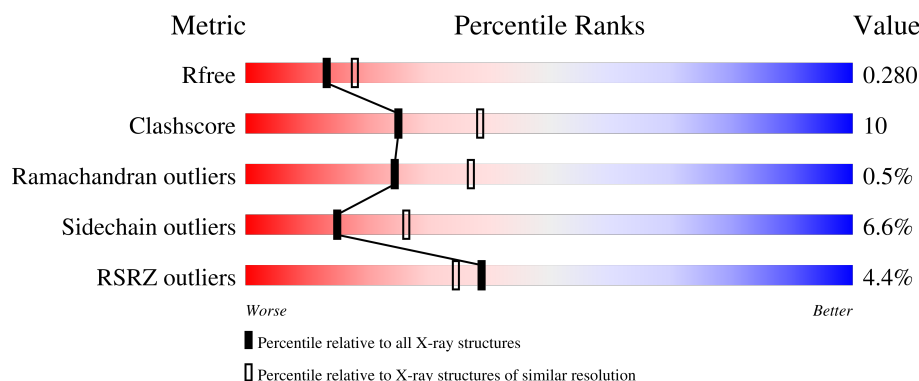
1 Overall quality at a glance

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	645	4% 75% 20% • •
2	B	350	5% 66% 21% • 11%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7699 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 Protein arginine N-methyltransferase 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	625	5137	3285	881	946	25	0	10	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	initiating methionine	UNP O14744
A	-6	ASP	-	expression tag	UNP O14744
A	-5	TYR	-	expression tag	UNP O14744
A	-4	LYS	-	expression tag	UNP O14744
A	-3	ASP	-	expression tag	UNP O14744
A	-2	ASP	-	expression tag	UNP O14744
A	-1	ASP	-	expression tag	UNP O14744
A	0	ASP	-	expression tag	UNP O14744
A	1	LYS	-	expression tag	UNP O14744

- Molecule 2 is a protein called Methylosome protein 50.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	310	2354	1476	402	461	15	0	2	0

There are 9 discrepancies between the modelled and reference sequences:

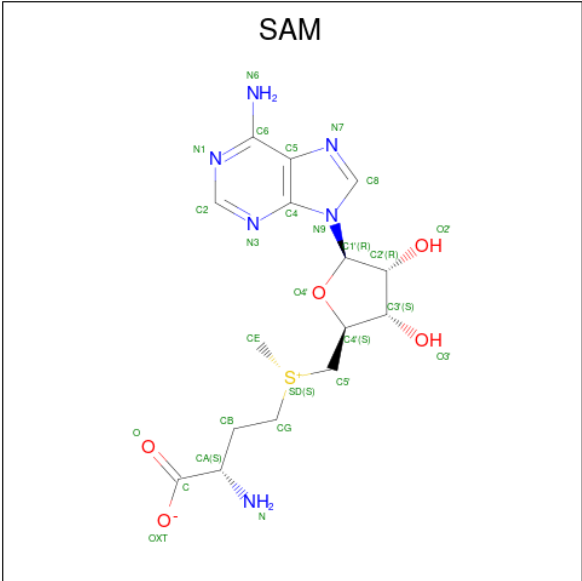
Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	MET	-	initiating methionine	UNP Q9BQA1
B	-6	HIS	-	expression tag	UNP Q9BQA1
B	-5	HIS	-	expression tag	UNP Q9BQA1
B	-4	HIS	-	expression tag	UNP Q9BQA1
B	-3	HIS	-	expression tag	UNP Q9BQA1
B	-2	HIS	-	expression tag	UNP Q9BQA1
B	-1	HIS	-	expression tag	UNP Q9BQA1

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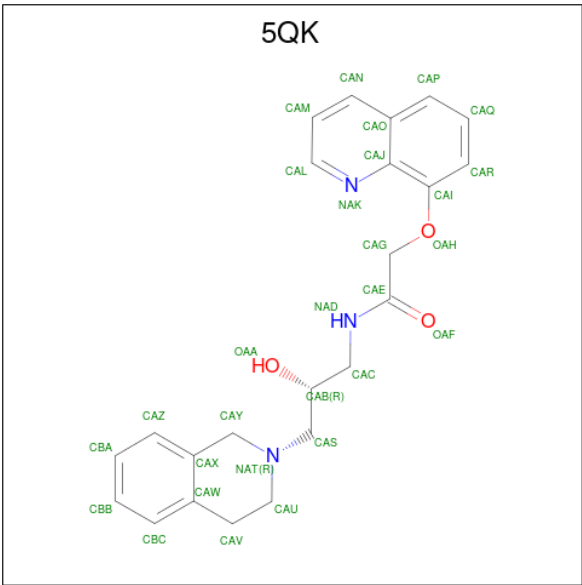
Chain	Residue	Modelled	Actual	Comment	Reference
B	0	HIS	-	expression tag	UNP Q9BQA1
B	1	HIS	-	expression tag	UNP Q9BQA1

- Molecule 3 is S-ADENOSYLMETHIONINE (CCD ID: SAM) (formula: C₁₅H₂₂N₆O₅S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			27	15	6	5	1		

- Molecule 4 is {N}-[(2 {R})-3-(3,4-dihydro-1 {H}-isoquinolin-2-yl)-2-oxidanyl-propyl]-2-quinolin-8-yloxy-ethanamide (CCD ID: 5QK) (formula: C₂₃H₂₅N₃O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			29	23	3	3		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		

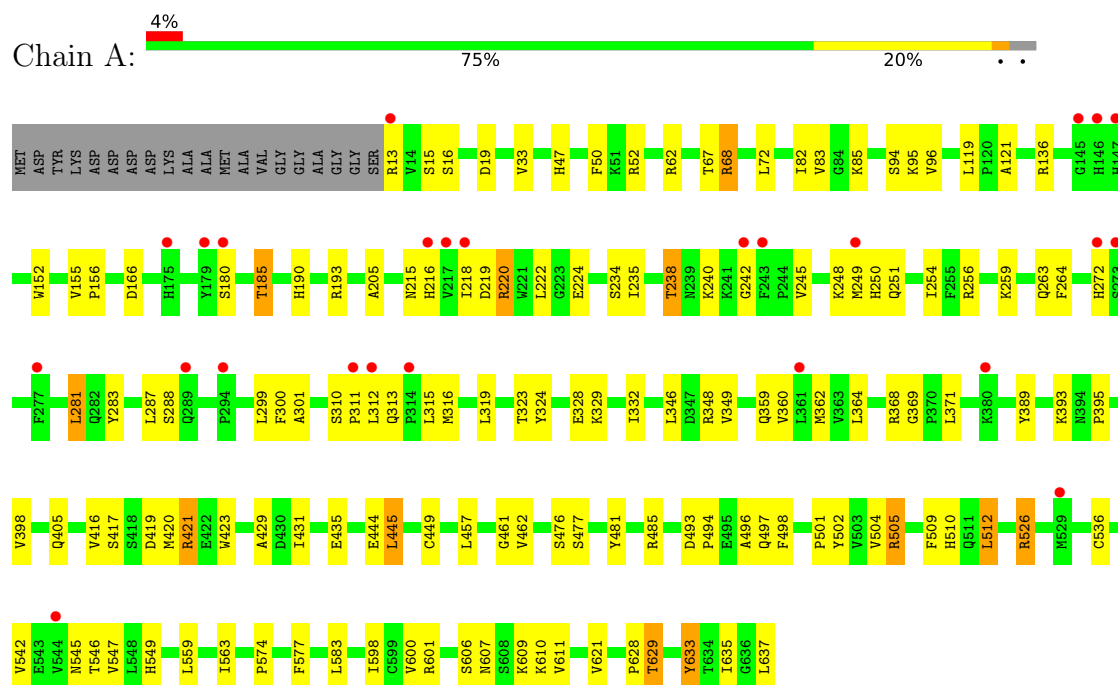
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	112	Total	O	0	0
			112	112		
6	B	28	Total	O	0	0
			28	28		

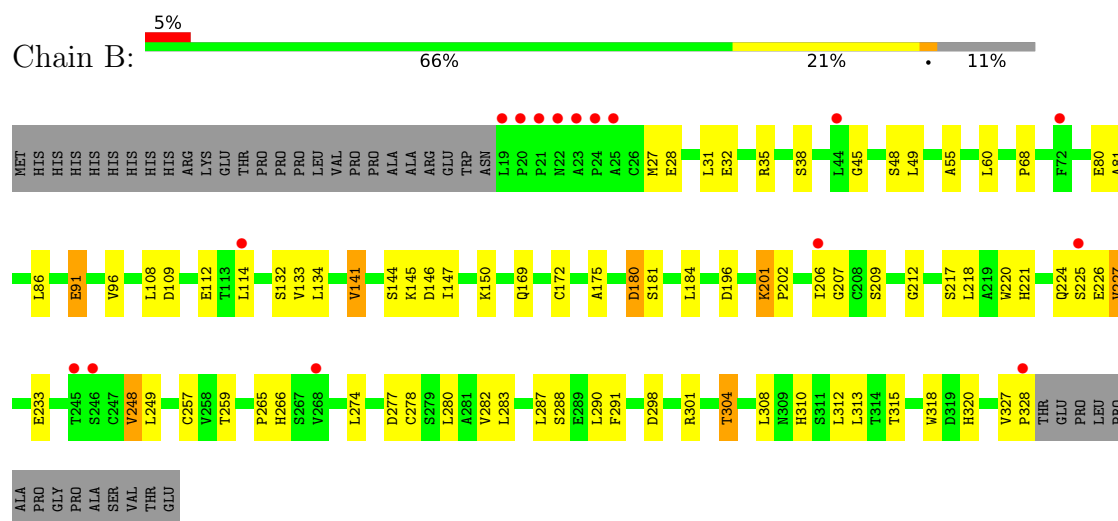
3 Residue-property plots [i](#)

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: Protein arginine N-methyltransferase 5



- Molecule 2: Methylosome protein 50



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	103.40Å 138.13Å 178.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.37 – 2.39 41.37 – 2.39	Depositor EDS
% Data completeness (in resolution range)	99.6 (41.37-2.39) 99.6 (41.37-2.39)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.08 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.208 , 0.282 0.208 , 0.280	Depositor DCC
R_{free} test set	2579 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	47.7	Xtriage
Anisotropy	0.256	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 32.3	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	0.95	EDS
Total number of atoms	7699	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.23% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SAM, 5QK

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	0.68	0/5281	0.96	7/7187 (0.1%)
2	B	0.59	0/2411	0.93	2/3295 (0.1%)
All	All	0.65	0/7692	0.95	9/10482 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	201	LYS	CA-C-N	8.24	128.20	119.05
2	B	201	LYS	C-N-CA	8.24	128.20	119.05
1	A	606	SER	N-CA-C	6.96	119.43	108.79
1	A	349	VAL	CA-C-N	5.68	125.59	119.85
1	A	349	VAL	C-N-CA	5.68	125.59	119.85
1	A	119	LEU	CA-C-N	5.24	124.75	119.19
1	A	119	LEU	C-N-CA	5.24	124.75	119.19
1	A	369	GLY	N-CA-C	5.19	122.93	112.34
1	A	82	ILE	N-CA-C	5.16	115.33	108.11

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	5137	0	5024	97	1
2	B	2354	0	2265	52	0
3	A	27	0	22	0	0
4	A	29	0	0	0	0
5	A	6	0	8	0	0
5	B	6	0	8	1	0
6	A	112	0	0	3	0
6	B	28	0	0	1	0
All	All	7699	0	7327	144	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 10.

All (144) 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:A:312[B]:LEU:H	1:A:312[B]:LEU:CD2	1.53	1.21
1:A:312[B]:LEU:H	1:A:312[B]:LEU:HD23	1.05	1.12
1:A:312[B]:LEU:CD2	1:A:312[B]:LEU:N	2.11	1.05
1:A:219:ASP:O	1:A:222:LEU:HG	1.60	1.01
1:A:215[B]:ASN:C	1:A:215[B]:ASN:HD22	1.73	0.96
1:A:312[B]:LEU:N	1:A:312[B]:LEU:HD22	1.86	0.90
1:A:222:LEU:HA	6:A:821:HOH:O	1.73	0.88
1:A:312[B]:LEU:HD23	1:A:312[B]:LEU:N	1.75	0.88
1:A:156:PRO:O	1:A:185:THR:HG21	1.81	0.80
1:A:312[B]:LEU:H	1:A:312[B]:LEU:HD22	1.46	0.77
1:A:215[B]:ASN:C	1:A:215[B]:ASN:ND2	2.42	0.77
1:A:215[B]:ASN:ND2	1:A:215[B]:ASN:O	2.20	0.74
1:A:68:ARG:HG3	1:A:72:LEU:HD12	1.72	0.71
1:A:628:PRO:O	1:A:629:THR:HB	1.93	0.69
1:A:607:ASN:HB3	1:A:609:LYS:H	1.57	0.68
1:A:419:ASP:OD1	1:A:421:ARG:HD3	1.96	0.66
2:B:35:ARG:HA	2:B:304:THR:HG21	1.77	0.65
2:B:259:THR:HG21	2:B:301:ARG:NH2	2.12	0.64
1:A:222:LEU:HD12	1:A:510:HIS:CD2	2.33	0.63
1:A:313[B]:GLN:HB3	1:A:316:MET:HB3	1.80	0.63
2:B:109:ASP:HB3	2:B:112:GLU:H	1.64	0.62
1:A:319:LEU:HD22	1:A:323:THR:HG21	1.80	0.62
1:A:250:HIS:O	1:A:254:ILE:HG12	2.00	0.62
1:A:485:ARG:HG3	1:A:498:PHE:HZ	1.63	0.62
2:B:144:SER:HB3	2:B:146:ASP:OD1	2.00	0.61
1:A:248:LYS:HA	1:A:251:GLN:HE21	1.66	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:27:MET:HE1	2:B:68:PRO:HB2	1.82	0.60
2:B:27:MET:HE1	2:B:68:PRO:CB	2.33	0.59
1:A:328:GLU:OE2	1:A:368:ARG:HD3	2.02	0.59
2:B:32:GLU:OE2	2:B:301:ARG:HD3	2.03	0.59
1:A:324:TYR:CB	1:A:368:ARG:HD2	2.33	0.59
1:A:542:VAL:HG21	1:A:598:ILE:HD12	1.85	0.57
1:A:190:HIS:ND1	1:A:193:ARG:NH2	2.46	0.57
2:B:86:LEU:HD11	2:B:96:VAL:HG22	1.86	0.57
1:A:216:HIS:O	1:A:219:ASP:HB2	2.04	0.57
2:B:172[B]:CYS:SG	2:B:217:SER:HA	2.45	0.57
1:A:16:SER:OG	1:A:281:LEU:HD21	2.05	0.57
1:A:193:ARG:NH1	1:A:224:GLU:OE2	2.38	0.56
2:B:109:ASP:HB3	2:B:112:GLU:N	2.20	0.56
2:B:221:HIS:HB2	2:B:227:VAL:CG2	2.36	0.56
1:A:362:MET:HG2	1:A:389:TYR:HB2	1.89	0.55
1:A:546:THR:OG1	1:A:547:VAL:N	2.40	0.55
2:B:27:MET:CE	2:B:68:PRO:HB2	2.37	0.54
1:A:348:ARG:HE	1:A:359:GLN:NE2	2.05	0.54
1:A:215[B]:ASN:ND2	1:A:219:ASP:OD1	2.41	0.54
1:A:444:GLU:O	1:A:445:LEU:HB2	2.07	0.54
2:B:27:MET:HE3	2:B:320:HIS:CE1	2.43	0.54
1:A:393:LYS:O	1:A:395:PRO:HD3	2.07	0.53
1:A:19:ASP:OD1	1:A:85:LYS:HE3	2.08	0.53
2:B:91[A]:GLU:CD	2:B:91[A]:GLU:N	2.67	0.52
1:A:481:TYR:HB2	1:A:549:HIS:CD2	2.44	0.52
1:A:249:MET:HA	1:A:249:MET:HE2	1.91	0.52
2:B:301:ARG:HD2	2:B:318:TRP:NE1	2.25	0.52
1:A:47:HIS:HB3	1:A:50:PHE:HB2	1.92	0.51
1:A:222:LEU:HB2	1:A:510:HIS:HB2	1.91	0.51
1:A:16:SER:HA	1:A:264:PHE:O	2.10	0.51
2:B:277:ASP:O	2:B:278:CYS:HB2	2.10	0.50
1:A:238:THR:HG23	1:A:272:HIS:HD2	1.77	0.49
2:B:27:MET:HE3	2:B:31:LEU:HD21	1.94	0.49
1:A:476:SER:HB2	1:A:512:LEU:HD11	1.95	0.49
1:A:364:LEU:HB3	1:A:420:MET:CE	2.43	0.49
1:A:633:TYR:OH	1:A:635:ILE:HD11	2.13	0.49
1:A:477:SER:HB2	1:A:509:PHE:HA	1.94	0.48
1:A:501:PRO:HA	1:A:583:LEU:O	2.14	0.48
2:B:248:VAL:CG2	2:B:249:LEU:N	2.76	0.48
2:B:109:ASP:HB2	2:B:114:LEU:O	2.14	0.48
2:B:206:ILE:HG13	2:B:207:GLY:H	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:445:LEU:HD22	1:A:635:ILE:HG22	1.95	0.48
2:B:109:ASP:HB2	2:B:114:LEU:H	1.79	0.48
1:A:68:ARG:HG3	1:A:72:LEU:CD1	2.43	0.48
1:A:445:LEU:HD11	1:A:611:VAL:HB	1.96	0.48
2:B:86:LEU:CD1	2:B:96:VAL:HG22	2.44	0.48
1:A:364:LEU:HB3	1:A:420:MET:HE2	1.95	0.48
2:B:86:LEU:HA	2:B:86:LEU:HD12	1.50	0.47
1:A:332:ILE:HG13	1:A:574:PRO:O	2.15	0.47
1:A:315[B]:LEU:HD21	1:A:449:CYS:SG	2.55	0.47
1:A:67:THR:OG1	1:A:68:ARG:N	2.48	0.47
1:A:526:ARG:HA	6:A:888:HOH:O	2.13	0.47
1:A:310[B]:SER:OG	1:A:311[B]:PRO:HD2	2.15	0.47
2:B:259:THR:HG21	2:B:301:ARG:HH22	1.77	0.47
1:A:313[A]:GLN:HG2	6:A:859:HOH:O	2.15	0.46
1:A:362:MET:SD	1:A:429:ALA:HB2	2.56	0.46
2:B:145:LYS:HA	2:B:169:GLN:CB	2.46	0.46
1:A:47:HIS:CE1	2:B:49:LEU:HD22	2.51	0.46
2:B:327:VAL:HA	2:B:328:PRO:HD2	1.84	0.46
2:B:80:GLU:HG3	2:B:81:ALA:H	1.81	0.46
2:B:145:LYS:HA	2:B:169:GLN:HB2	1.97	0.45
1:A:485:ARG:HG3	1:A:498:PHE:CZ	2.48	0.45
2:B:48:SER:HB3	2:B:55:ALA:H	1.81	0.45
1:A:245:VAL:HG12	1:A:283:TYR:CE1	2.52	0.45
1:A:301:ALA:HB1	1:A:505:ARG:HG3	1.99	0.45
1:A:502:TYR:O	1:A:583:LEU:N	2.44	0.45
1:A:600:VAL:C	1:A:601:ARG:HG2	2.42	0.45
2:B:221:HIS:HB2	2:B:227:VAL:HG22	1.98	0.45
2:B:221:HIS:CB	2:B:227:VAL:HG22	2.46	0.45
2:B:181:SER:O	2:B:196:ASP:HA	2.17	0.44
2:B:212:GLY:H	5:B:401:GOL:H31	1.82	0.44
1:A:348:ARG:HE	1:A:359:GLN:HE22	1.65	0.44
2:B:45:GLY:HA3	2:B:86:LEU:HD22	2.00	0.44
1:A:607:ASN:ND2	1:A:610:LYS:HE2	2.32	0.44
1:A:328:GLU:OE1	1:A:368:ARG:NH1	2.51	0.43
2:B:265:PRO:HB2	2:B:310:HIS:CE1	2.53	0.43
1:A:96:VAL:HG21	2:B:233:GLU:HG3	1.99	0.43
1:A:248:LYS:HD2	1:A:251:GLN:NE2	2.34	0.43
1:A:220:ARG:NH1	1:A:546:THR:HA	2.34	0.43
1:A:62:ARG:CD	2:B:298:ASP:OD2	2.67	0.43
1:A:83:VAL:HG22	1:A:121:ALA:HB3	2.01	0.43
2:B:60:LEU:HD22	2:B:108:LEU:HD21	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:ARG:HD3	2:B:298:ASP:OD2	2.18	0.42
2:B:280:LEU:HD22	2:B:315:THR:HG21	2.02	0.42
1:A:324:TYR:HB2	1:A:368:ARG:HD2	2.01	0.42
1:A:50:PHE:O	1:A:52:ARG:HD3	2.19	0.42
2:B:304:THR:O	2:B:313:LEU:HD12	2.20	0.42
1:A:152:TRP:HB3	1:A:205:ALA:HB2	2.01	0.42
1:A:512:LEU:HG	1:A:546:THR:HG21	2.02	0.42
1:A:15:SER:HB3	1:A:263:GLN:HG2	2.01	0.42
1:A:312[B]:LEU:HD23	1:A:312[B]:LEU:O	2.20	0.42
1:A:421:ARG:HH12	1:A:637:LEU:C	2.28	0.42
2:B:109:ASP:HB3	2:B:112:GLU:CA	2.50	0.42
1:A:481:TYR:HB2	1:A:549:HIS:HD2	1.81	0.42
1:A:600:VAL:HG22	1:A:601:ARG:N	2.35	0.42
1:A:398:VAL:HG11	1:A:417:SER:HB2	2.02	0.41
1:A:238:THR:HG22	1:A:242:GLY:HA2	2.02	0.41
1:A:300:PHE:CD2	1:A:577:PHE:HB2	2.55	0.41
2:B:180:ASP:HB2	6:B:521:HOH:O	2.20	0.41
1:A:166:ASP:O	2:B:201:LYS:HE2	2.20	0.41
2:B:221:HIS:HB2	2:B:227:VAL:HG23	2.01	0.41
1:A:493:ASP:HB3	1:A:496:ALA:HB2	2.03	0.41
1:A:493:ASP:HA	1:A:494:PRO:HD3	1.88	0.41
1:A:600:VAL:HG22	1:A:601:ARG:H	1.85	0.41
2:B:141:VAL:HA	2:B:150:LYS:O	2.21	0.41
2:B:184:LEU:HG	2:B:220:TRP:CZ2	2.56	0.41
2:B:28:GLU:HG3	2:B:48:SER:HB2	2.02	0.41
1:A:220:ARG:NH2	1:A:545:ASN:HB2	2.35	0.40
1:A:346:LEU:HD23	1:A:346:LEU:HA	1.87	0.40
1:A:416:VAL:HG11	1:A:423:TRP:CD1	2.57	0.40
2:B:134:LEU:HD23	2:B:175:ALA:HB1	2.03	0.40
1:A:218:ILE:HG22	1:A:218:ILE:O	2.21	0.40
1:A:563:ILE:HD12	1:A:563:ILE:N	2.37	0.40
2:B:201:LYS:HA	2:B:202:PRO:HD2	1.82	0.40
2:B:282:VAL:HB	2:B:291:PHE:HB3	2.04	0.40
1:A:235:ILE:HD12	1:A:250:HIS:CD2	2.57	0.40
1:A:431:ILE:HA	1:A:462:VAL:O	2.22	0.40
1:A:457:LEU:HG	1:A:461:GLY:HA3	2.02	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:A:240:LYS:NZ	1:A:240:LYS:NZ[2_645]	2.15	0.05

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	633/645 (98%)	600 (95%)	32 (5%)	1 (0%)	43	58
2	B	310/350 (89%)	293 (94%)	13 (4%)	4 (1%)	9	14
All	All	943/995 (95%)	893 (95%)	45 (5%)	5 (0%)	24	37

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	147	ILE
1	A	234	SER
2	B	266	HIS
2	B	209	SER
2	B	224	GLN

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	568/570 (100%)	534 (94%)	34 (6%)	17	31
2	B	265/298 (89%)	244 (92%)	21 (8%)	11	19
All	All	833/868 (96%)	778 (93%)	55 (7%)	15	26

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ARG
1	A	33	VAL
1	A	68	ARG
1	A	94	SER
1	A	95	LYS
1	A	136	ARG
1	A	155	VAL
1	A	180	SER
1	A	185	THR
1	A	220	ARG
1	A	238	THR
1	A	256	ARG
1	A	259	LYS
1	A	281	LEU
1	A	287	LEU
1	A	288	SER
1	A	299	LEU
1	A	329	LYS
1	A	360	VAL
1	A	371	LEU
1	A	405	GLN
1	A	421	ARG
1	A	435	GLU
1	A	445	LEU
1	A	497	GLN
1	A	504	VAL
1	A	505	ARG
1	A	512	LEU
1	A	526	ARG
1	A	536	CYS
1	A	559	LEU
1	A	621	VAL
1	A	629	THR
1	A	633	TYR
2	B	38	SER
2	B	91[A]	GLU
2	B	91[B]	GLU
2	B	132	SER
2	B	133	VAL
2	B	141	VAL
2	B	180	ASP
2	B	218	LEU

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Mol	Chain	Res	Type
2	B	225	SER
2	B	226	GLU
2	B	227	VAL
2	B	248	VAL
2	B	257	CYS
2	B	274	LEU
2	B	283	LEU
2	B	287	LEU
2	B	288	SER
2	B	290	LEU
2	B	304	THR
2	B	308	LEU
2	B	312	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	56	GLN
1	A	140	ASN
1	A	251	GLN
1	A	263	GLN
1	A	270	ASN
1	A	309	GLN
1	A	359	GLN
1	A	396	ASN
1	A	508	ASN
1	A	511	GLN
1	A	525	ASN
2	B	310	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	$\# Z > 2$
5	GOL	A	703	-	5,5,5	0.27	0	5,5,5	1.00	0
4	5QK	A	702	-	32,32,32	1.96	4 (12%)	43,43,43	1.31	6 (13%)
3	SAM	A	701	-	27,29,29	1.55	3 (11%)	34,42,42	2.09	9 (26%)
5	GOL	B	401	-	5,5,5	0.29	0	5,5,5	0.91	0

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
5	GOL	A	703	-	-	4/4/4/4	-
4	5QK	A	702	-	-	8/14/23/23	0/4/4/4
3	SAM	A	701	-	-	2/17/33/33	0/3/3/3
5	GOL	B	401	-	-	4/4/4/4	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	702	5QK	CAV-CAW	-6.16	1.40	1.51
4	A	702	5QK	CAL-NAK	5.20	1.42	1.32
4	A	702	5QK	CAY-CAX	-4.93	1.40	1.51
3	A	701	SAM	C5-C4	4.76	1.47	1.39
3	A	701	SAM	C5-N7	-3.20	1.33	1.39
3	A	701	SAM	C8-N7	2.31	1.36	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	702	5QK	CAE-NAD	2.15	1.38	1.33

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	701	SAM	C5-C4-N3	-6.29	118.06	126.72
3	A	701	SAM	N3-C4-N9	5.15	135.92	127.17
3	A	701	SAM	C2-N3-C4	3.31	119.91	111.83
3	A	701	SAM	CG-SD-C5'	3.30	111.51	103.43
3	A	701	SAM	C4-C5-N7	-3.18	106.95	110.58
4	A	702	5QK	CAG-OAH-CAI	-3.15	109.98	116.68
4	A	702	5QK	CAX-CAY-NAT	3.11	115.72	112.14
3	A	701	SAM	N3-C2-N1	-2.85	124.27	128.58
4	A	702	5QK	CAL-NAK-CAJ	2.75	120.55	117.31
4	A	702	5QK	CAO-CAJ-NAK	-2.73	119.75	122.63
3	A	701	SAM	C5-N7-C8	2.64	107.59	103.45
4	A	702	5QK	CAM-CAL-NAK	-2.58	120.18	123.97
3	A	701	SAM	O4'-C1'-C2'	-2.54	101.17	106.62
3	A	701	SAM	C4-N9-C8	2.26	108.11	105.74
4	A	702	5QK	CAI-CAJ-NAK	2.09	121.36	118.89

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	702	5QK	CAS-CAB-CAC-NAD
4	A	702	5QK	OAA-CAB-CAC-NAD
5	B	401	GOL	C1-C2-C3-O3
5	A	703	GOL	O1-C1-C2-C3
5	B	401	GOL	O1-C1-C2-C3
5	A	703	GOL	O1-C1-C2-O2
5	B	401	GOL	O2-C2-C3-O3
4	A	702	5QK	CAR-CAI-OAH-CAG
5	B	401	GOL	O1-C1-C2-O2
4	A	702	5QK	CAJ-CAI-OAH-CAG
4	A	702	5QK	CAC-CAB-CAS-NAT
4	A	702	5QK	OAA-CAB-CAS-NAT
5	A	703	GOL	C1-C2-C3-O3
4	A	702	5QK	CAB-CAS-NAT-CAY
4	A	702	5QK	CAB-CAS-NAT-CAU
5	A	703	GOL	O2-C2-C3-O3
3	A	701	SAM	OXT-C-CA-N

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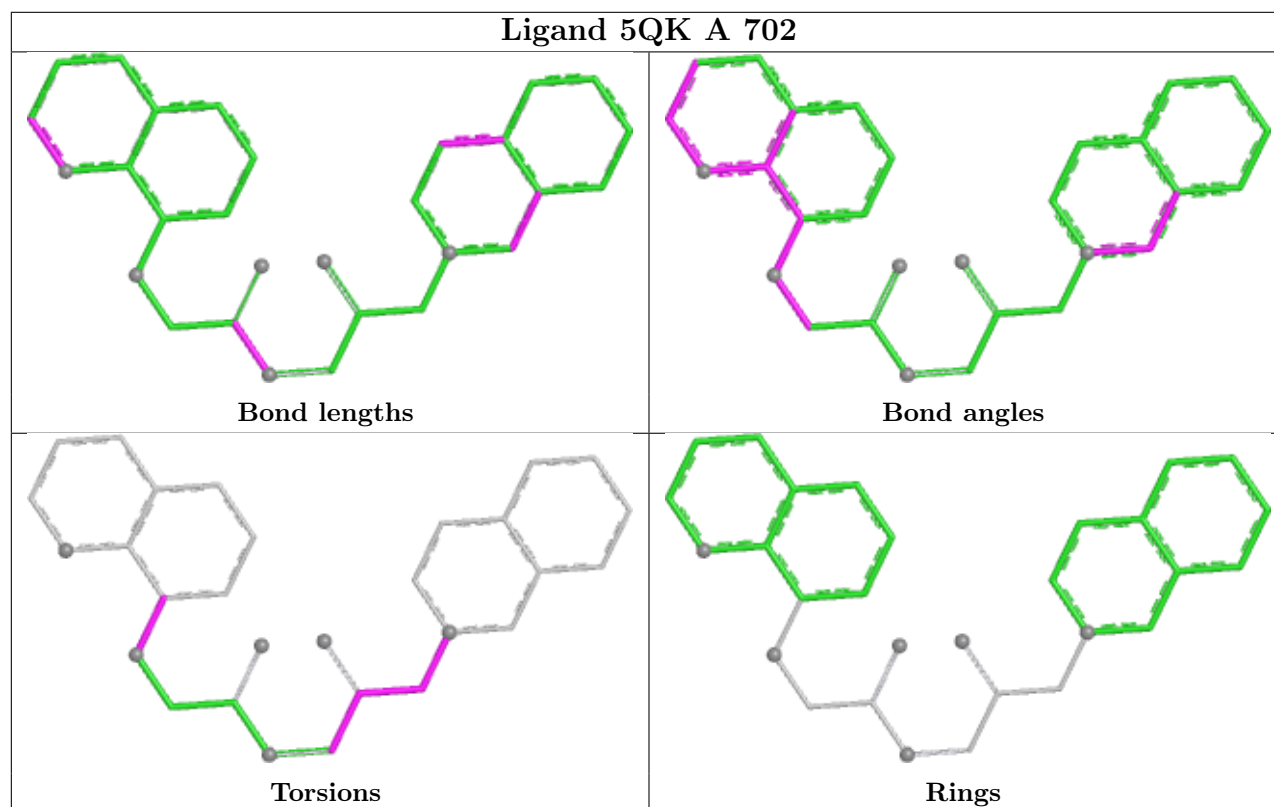
Mol	Chain	Res	Type	Atoms
3	A	701	SAM	CB-CG-SD-C5'

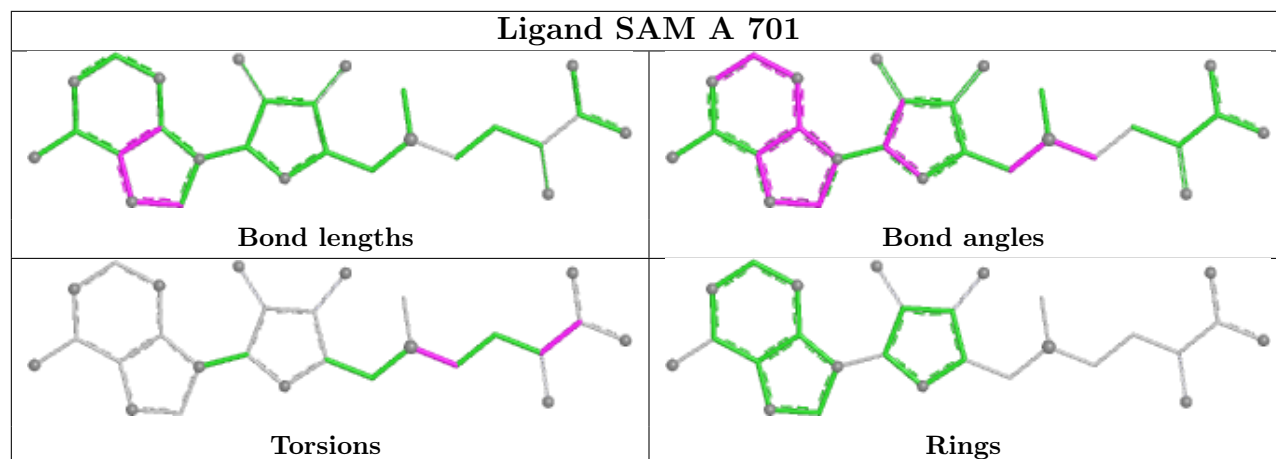
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	401	GOL	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

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	625/645 (96%)	0.03	25 (4%) 42 38	16, 44, 84, 134	10 (1%)
2	B	310/350 (88%)	0.41	16 (5%) 33 29	26, 56, 96, 142	2 (0%)
All	All	935/995 (93%)	0.15	41 (4%) 39 34	16, 49, 90, 142	12 (1%)

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	312[A]	LEU	4.8
2	B	19	LEU	4.7
2	B	21	PRO	4.7
1	A	311[A]	PRO	4.6
1	A	243	PHE	4.2
1	A	277	PHE	4.1
1	A	217	VAL	3.7
1	A	529	MET	3.6
1	A	146	HIS	3.3
2	B	25	ALA	3.3
1	A	314[A]	PRO	3.2
2	B	22	ASN	3.2
1	A	544	VAL	3.1
2	B	72	PHE	3.1
2	B	225	SER	3.0
1	A	145	GLY	2.7
2	B	246	SER	2.7
1	A	249	MET	2.7
1	A	272	HIS	2.5
2	B	268	VAL	2.5
1	A	380	LYS	2.5
1	A	180	SER	2.4
1	A	179	TYR	2.4
1	A	218	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	206	ILE	2.3
2	B	20	PRO	2.3
1	A	361	LEU	2.3
2	B	328	PRO	2.3
2	B	23	ALA	2.3
1	A	294	PRO	2.2
1	A	147	HIS	2.2
2	B	245	THR	2.2
2	B	24	PRO	2.2
1	A	242	GLY	2.2
1	A	216	HIS	2.2
1	A	175	HIS	2.1
1	A	289	GLN	2.1
2	B	114	LEU	2.1
2	B	44	LEU	2.1
1	A	273	SER	2.1
1	A	13	ARG	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

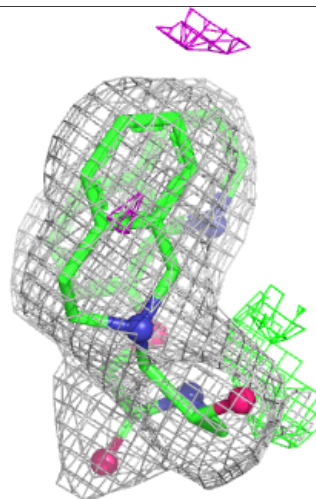
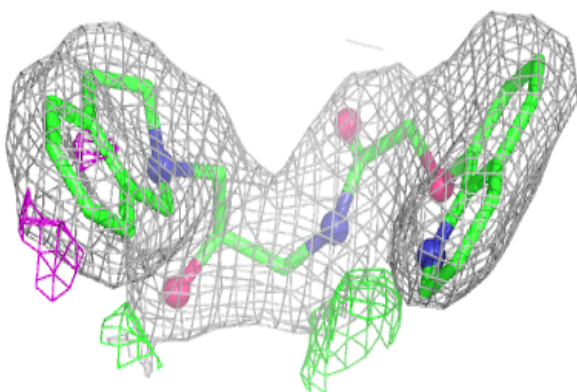
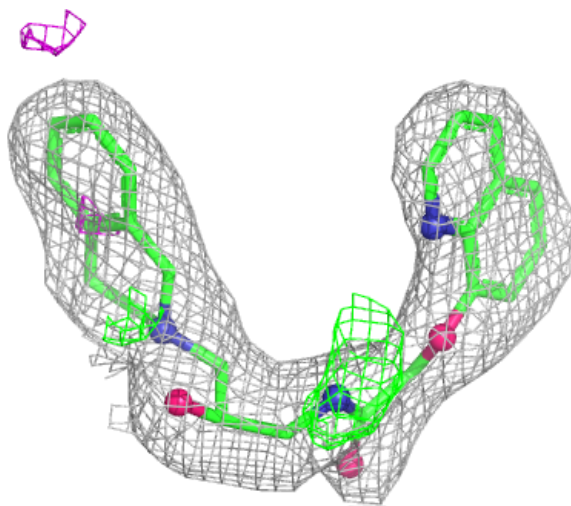
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	GOL	B	401	6/6	0.85	0.12	60,61,66,67	0
5	GOL	A	703	6/6	0.89	0.10	47,51,55,57	0
4	5QK	A	702	29/29	0.90	0.11	44,51,67,69	0
3	SAM	A	701	27/27	0.97	0.07	31,34,46,51	0

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.

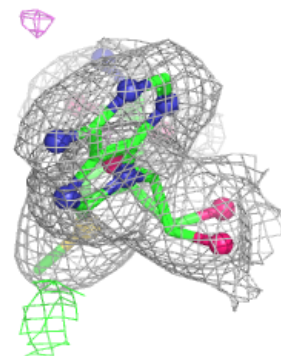
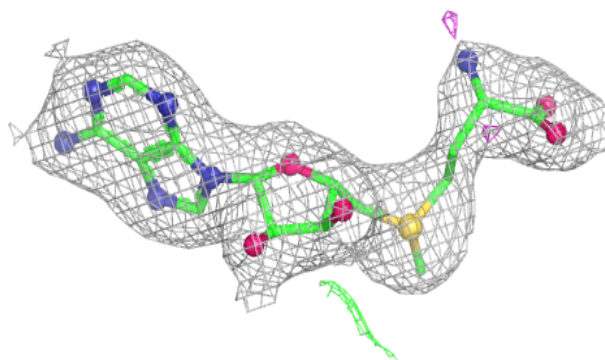
Electron density around 5QK A 702:

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



Electron density around SAM A 701:

$2mF_o-DF_c$ (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.