



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 21, 2016 – 08:55 PM EDT

PDB ID : 5EML
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027790
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027790

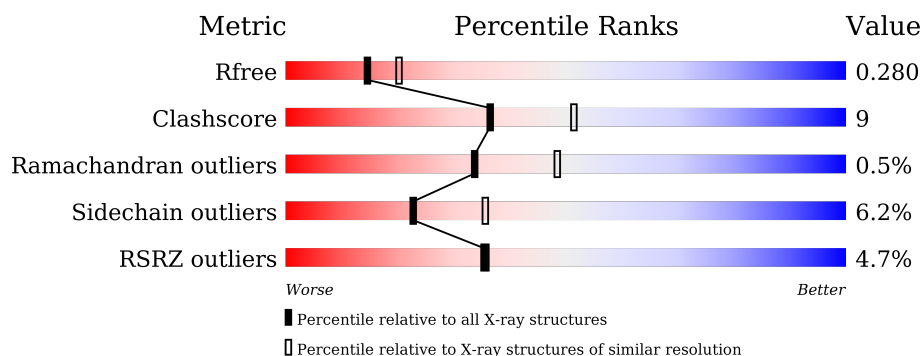
1 Overall quality at a glance

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}	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	<div> <div>3%</div> <div>76%</div> <div>19%</div> <div>• •</div> </div>
2	B	350	<div> <div>7%</div> <div>66%</div> <div>21%</div> <div>• 11%</div> </div>

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
5	GOL	A	703	-	-	-	X

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:

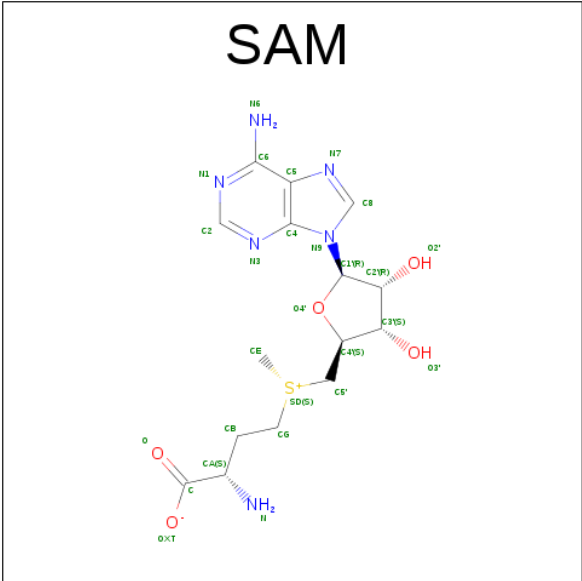
Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	MET	-	initiating methionine	UNP Q9BQA1
B	-6	HIS	-	expression tag	UNP Q9BQA1
B	-5	HIS	-	expression tag	UNP Q9BQA1
B	-4	HIS	-	expression tag	UNP Q9BQA1
B	-3	HIS	-	expression tag	UNP Q9BQA1
B	-2	HIS	-	expression tag	UNP Q9BQA1
B	-1	HIS	-	expression tag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

Continued on next page...

Continued from previous page...

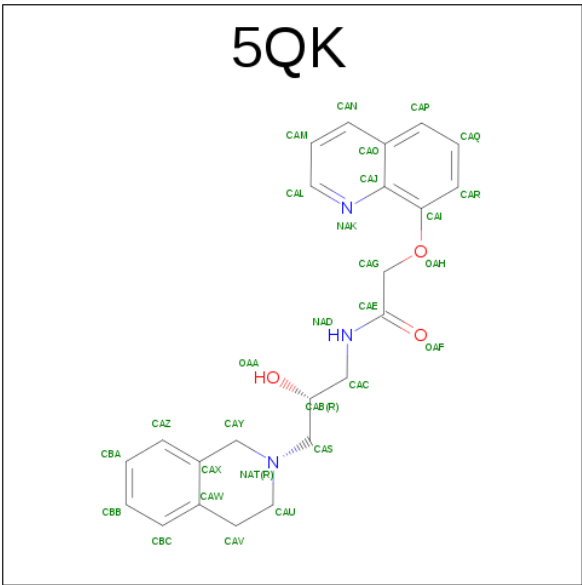
Chain	Residue	Modelled	Actual	Comment	Reference
B	0	HIS	-	expression tag	UNP Q9BQA1
B	1	HIS	-	expression tag	UNP Q9BQA1

- Molecule 3 is S-ADENOSYLMETHIONINE (three-letter code: SAM) (formula: C₁₅H₂₂N₆O₅S).



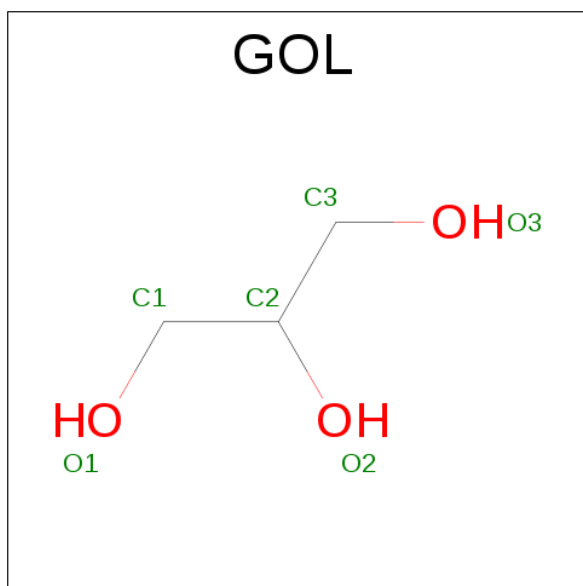
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
3	A	1	27	15	6	5	1	0	0

- Molecule 4 is {N}-[(2 {R})-3-(3,4-dihydro-1 {H}-isoquinolin-2-yl)-2-oxidanyl-propyl]-2-quinolin-8-yloxy-ethanamide (three-letter code: 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 (three-letter code: 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		

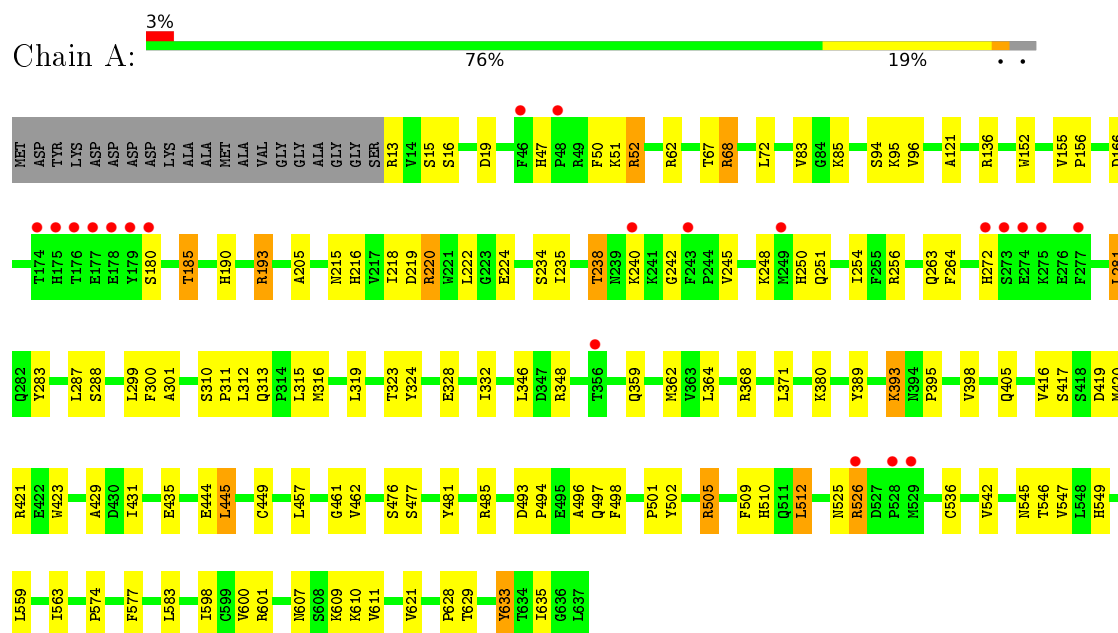
- Molecule 6 is water.

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

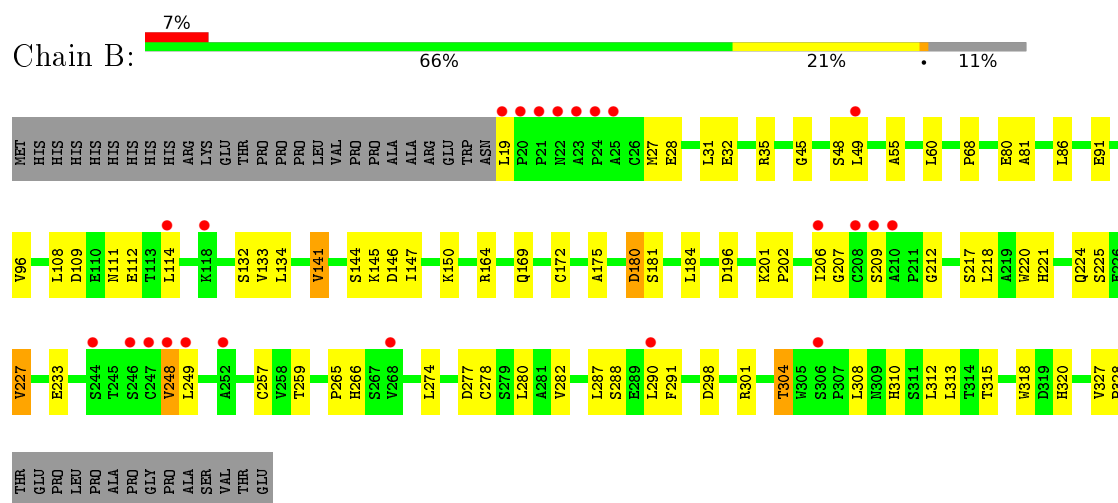
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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	103.40 Å 138.13 Å 178.92 Å 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 2.39 Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.208 , 0.282 0.208 , 0.280	Depositor DCC
R_{free} test set	2579 reflections (5.35%)	DCC
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 , 37.0	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, 5QK, SAM

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	0.53	0/5281	0.70	1/7187 (0.0%)
2	B	0.44	0/2411	0.66	0/3295
All	All	0.50	0/7692	0.69	1/10482 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	193	ARG	NE-CZ-NH1	-5.03	117.78	120.30

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	94	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

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	7699	0	7327	141	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 (141) 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:N	1:A:312[B]:LEU:CD2	2.11	1.05
1:A:219:ASP:O	1:A:222:LEU:HG	1.60	1.01
1:A:215[B]:ASN:HD22	1:A:215[B]:ASN:C	1.73	0.91
1:A:312[B]:LEU:HD22	1:A:312[B]:LEU:N	1.86	0.90
1:A:222:LEU:HA	6:A:821:HOH:O	1.73	0.88
1:A:312[B]:LEU:N	1:A:312[B]:LEU:HD23	1.75	0.88
1:A:156:PRO:O	1:A:185:THR:HG21	1.81	0.80
1:A:312[B]:LEU:HD22	1:A:312[B]:LEU:H	1.46	0.77
1:A:215[B]:ASN:O	1:A:215[B]:ASN:ND2	2.20	0.74
1:A:215[B]:ASN:C	1:A:215[B]:ASN:ND2	2.42	0.72
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
2:B:27:MET:HE1	2:B:68:PRO:HB2	1.77	0.66
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
2:B:27:MET:HE1	2:B:68:PRO:CB	2.30	0.61
1:A:248:LYS:HA	1:A:251:GLN:HE21	1.66	0.61
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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
2:B:27:MET:HE3	2:B:31:LEU:HD21	1.88	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
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
1:A:481:TYR:HB2	1:A:549:HIS:CD2	2.44	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
2:B:248:VAL:CG2	2:B:249:LEU:N	2.76	0.49
1:A:238:THR:HG23	1:A:272:HIS:HD2	1.77	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: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
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:445:LEU:HD11	1:A:611:VAL:HB	1.96	0.48
1:A:68:ARG:HG3	1:A:72:LEU:CD1	2.43	0.48
2:B:86:LEU:CD1	2:B:96:VAL:HG22	2.44	0.48
2:B:91[A]:GLU:CD	2:B:91[A]:GLU:N	2.67	0.48
2:B:86:LEU:HD12	2:B:86:LEU:HA	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:526:ARG:HA	6:A:888:HOH:O	2.13	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:THR:OG1	1:A:68:ARG:N	2.48	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
2:B:27:MET:HE3	2:B:320:HIS:CE1	2.50	0.46
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
2:B:327:VAL:HA	2:B:328:PRO:HD2	1.84	0.46
1:A:47:HIS:CE1	2:B:49:LEU:HD22	2.51	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
2:B:221:HIS:CB	2:B:227:VAL:HG22	2.46	0.45
2:B:221:HIS:HB2	2:B:227:VAL:HG22	1.98	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:248:LYS:HD2	1:A:251:GLN:NE2	2.34	0.43
1:A:96:VAL:HG21	2:B:233:GLU:HG3	1.99	0.43
1:A:220:ARG:NH1	1:A:546:THR:HA	2.34	0.43
1:A:83:VAL:HG22	1:A:121:ALA:HB3	2.01	0.43
1:A:62:ARG:CD	2:B:298:ASP:OD2	2.67	0.43
2:B:60:LEU:HD22	2:B:108:LEU:HD21	2.01	0.43
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:O	1:A:312[B]:LEU:HD23	2.20	0.42
2:B:109:ASP:HB3	2:B:112:GLU:CA	2.50	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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
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: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:HA	1:A:346:LEU:HD23	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:364:LEU:HB3	1:A:420:MET:HE1	2.04	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%)	52	69
2	B	310/350 (89%)	293 (94%)	13 (4%)	4 (1%)	15	21
All	All	943/995 (95%)	893 (95%)	45 (5%)	5 (0%)	34	48

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 ⓘ

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%)	536 (94%)	32 (6%)	26	41
2	B	265/298 (89%)	246 (93%)	19 (7%)	18	28
All	All	833/868 (96%)	782 (94%)	51 (6%)	23	36

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ARG
1	A	51	LYS
1	A	52	ARG
1	A	68	ARG
1	A	94	SER
1	A	95	LYS
1	A	136	ARG
1	A	155	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	180	SER
1	A	185	THR
1	A	220	ARG
1	A	238	THR
1	A	256	ARG
1	A	281	LEU
1	A	287	LEU
1	A	288	SER
1	A	299	LEU
1	A	371	LEU
1	A	380	LYS
1	A	393	LYS
1	A	405	GLN
1	A	435	GLU
1	A	445	LEU
1	A	497	GLN
1	A	505	ARG
1	A	512	LEU
1	A	525	ASN
1	A	526	ARG
1	A	536	CYS
1	A	559	LEU
1	A	621	VAL
1	A	633	TYR
2	B	19	LEU
2	B	111	ASN
2	B	132	SER
2	B	133	VAL
2	B	141	VAL
2	B	164	ARG
2	B	180	ASP
2	B	218	LEU
2	B	225	SER
2	B	227	VAL
2	B	248	VAL
2	B	257	CYS
2	B	274	LEU
2	B	287	LEU
2	B	288	SER
2	B	290	LEU
2	B	304	THR
2	B	308	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	B	312	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	GLN
1	A	140	ASN
1	A	251	GLN
1	A	309	GLN
1	A	336	GLN
1	A	359	GLN
1	A	510	HIS
1	A	511	GLN
1	A	525	ASN
2	B	310	HIS

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 carbohydrates 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
3	SAM	A	701	-	23,29,29	1.71	3 (13%)	15,42,42	1.79	2 (13%)
4	5QK	A	702	-	32,32,32	1.93	6 (18%)	42,43,43	1.54	6 (14%)
5	GOL	A	703	-	5,5,5	0.23	0	5,5,5	0.85	0
5	GOL	B	401	-	5,5,5	0.26	0	5,5,5	0.76	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
3	SAM	A	701	-	-	0/8/33/33	0/3/3/3
4	5QK	A	702	-	-	0/14/23/23	0/4/4/4
5	GOL	A	703	-	-	0/4/4/4	0/0/0/0
5	GOL	B	401	-	-	0/4/4/4	0/0/0/0

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	702	5QK	CAV-CAW	-6.02	1.40	1.51
3	A	701	SAM	CG-SD	-5.79	1.68	1.80
4	A	702	5QK	CAY-CAX	-4.84	1.40	1.51
3	A	701	SAM	O4'-C1'	2.05	1.44	1.41
4	A	702	5QK	CAC-CAB	2.06	1.56	1.52
4	A	702	5QK	CAE-NAD	2.10	1.38	1.33
4	A	702	5QK	CAS-CAB	2.17	1.57	1.52
3	A	701	SAM	C5-C4	3.14	1.47	1.40
4	A	702	5QK	CAL-NAK	4.98	1.42	1.32

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	702	5QK	CAG-OAH-CAI	-6.18	109.98	116.88
3	A	701	SAM	N3-C2-N1	-5.85	124.27	128.87
4	A	702	5QK	CAO-CAJ-NAK	-2.52	119.75	122.54
4	A	702	5QK	CAM-CAL-NAK	-2.37	120.18	123.94
3	A	701	SAM	N6-C6-N1	2.42	122.57	118.52
4	A	702	5QK	CAX-CAY-NAT	2.44	115.72	111.83
4	A	702	5QK	CAL-NAK-CAJ	2.77	120.55	117.36
4	A	702	5QK	CAI-CAJ-NAK	2.93	121.36	118.86

There are no chirality outliers.

There are no torsion outliers.

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

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%)	0.01	21 (3%) 49 49	19, 45, 84, 134	0
2	B	310/350 (88%)	0.35	23 (7%) 17 17	37, 56, 96, 142	0
All	All	935/995 (93%)	0.12	44 (4%) 35 36	19, 49, 90, 142	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	25	ALA	5.9
1	A	176	THR	5.8
1	A	175	HIS	5.8
2	B	21	PRO	5.3
2	B	19	LEU	5.2
1	A	177	GLU	4.4
2	B	22	ASN	4.3
1	A	272	HIS	4.1
1	A	240	LYS	4.1
1	A	174	THR	3.9
1	A	243	PHE	3.8
2	B	114	LEU	3.6
2	B	23	ALA	3.4
1	A	529	MET	3.2
2	B	244	SER	3.1
2	B	118	LYS	3.1
1	A	179	TYR	3.0
1	A	277	PHE	3.0
2	B	290	LEU	2.9
1	A	180	SER	2.9
1	A	178	GLU	2.9
2	B	210	ALA	2.9
2	B	209	SER	2.9
2	B	248	VAL	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	273	SER	2.7
1	A	274	GLU	2.6
1	A	249	MET	2.6
2	B	306	SER	2.6
1	A	48	PRO	2.5
2	B	208	CYS	2.5
2	B	20	PRO	2.4
2	B	249	LEU	2.4
2	B	268	VAL	2.4
1	A	275	LYS	2.2
2	B	247	CYS	2.2
2	B	206	ILE	2.1
2	B	24	PRO	2.1
1	A	356	THR	2.1
1	A	528	PRO	2.1
1	A	526	ARG	2.1
2	B	252	ALA	2.1
1	A	46	PHE	2.1
2	B	246	SER	2.0
2	B	49	LEU	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
5	GOL	A	703	6/6	0.89	0.20	2.30	47,51,55,57	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	5QK	A	702	29/29	0.89	0.15	0.73	44,51,67,69	0
3	SAM	A	701	27/27	0.97	0.12	-0.19	31,34,46,51	0
5	GOL	B	401	6/6	0.89	0.16	-0.44	60,61,66,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.