



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 08:32 pm GMT

PDB ID : 3I53
Title : Crystal structure of an O-methyltransferase (NcsB1) from neocarzinostatin biosynthesis in complex with S-adenosyl-L-homocysteine (SAH)
Authors : Cooke, H.A.; Bruner, S.D.
Deposited on : 2009-07-03
Resolution : 2.08 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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

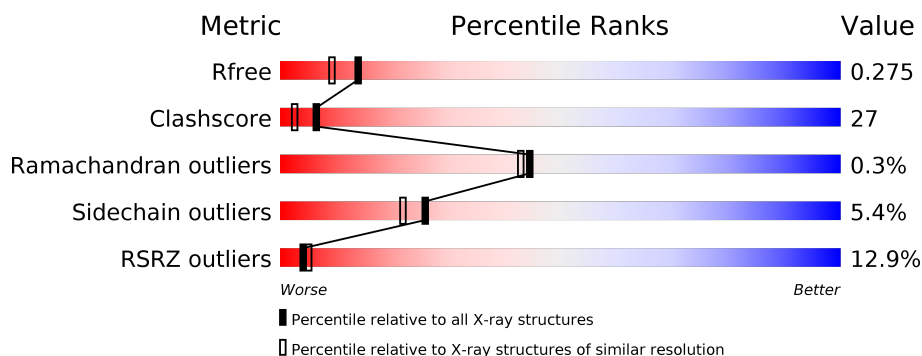
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.08 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}	100719	4955 (2.10-2.06)
Clashscore	112137	5537 (2.10-2.06)
Ramachandran outliers	110173	5483 (2.10-2.06)
Sidechain outliers	110143	5484 (2.10-2.06)
RSRZ outliers	101464	4991 (2.10-2.06)

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	332	<div> <div>3%</div> <div>73%</div> <div>25%</div> <div>.</div> </div>
1	B	332	<div> <div>21%</div> <div>44%</div> <div>42%</div> <div>5%</div> <div>9%</div> </div>

2 Entry composition [i](#)

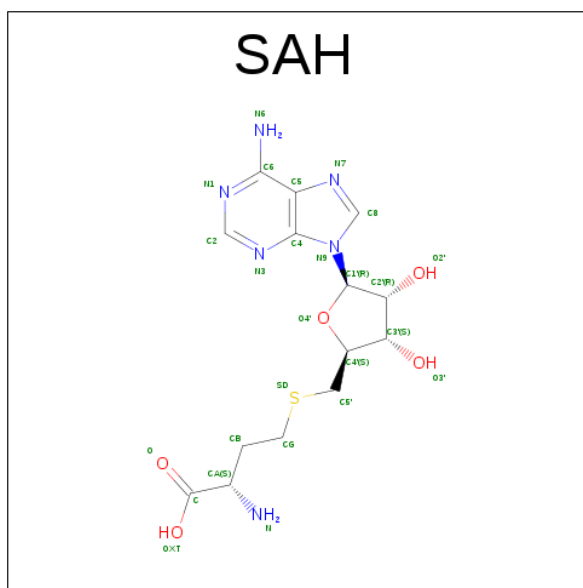
There are 4 unique types of molecules in this entry. The entry contains 4935 atoms, of which 0 are hydrogens and 0 are deuteriums.

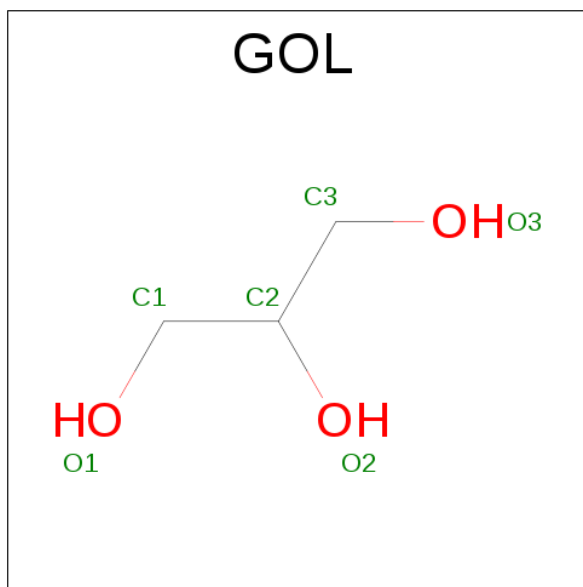
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called O-methyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	325	Total	C	N	O	S	0	0	0
			2379	1495	428	449	7			
1	B	303	Total	C	N	O	S	0	0	0
			2193	1378	396	412	7			

- Molecule 2 is S-ADENOSYL-L-HOMOCYSTEINE (three-letter code: SAH) (formula: $C_{14}H_{20}N_6O_5S$).





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

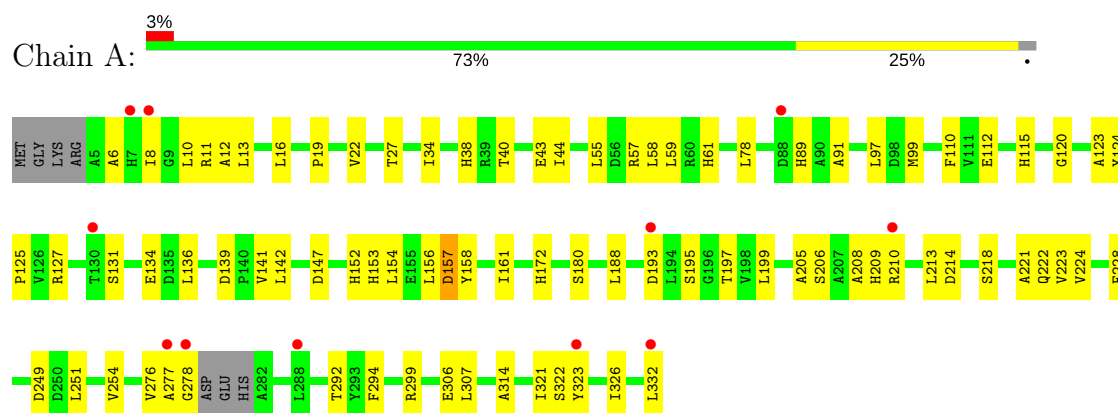
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	188	Total	O	0	0
			188	188		
4	B	117	Total	O	0	0
			117	117		

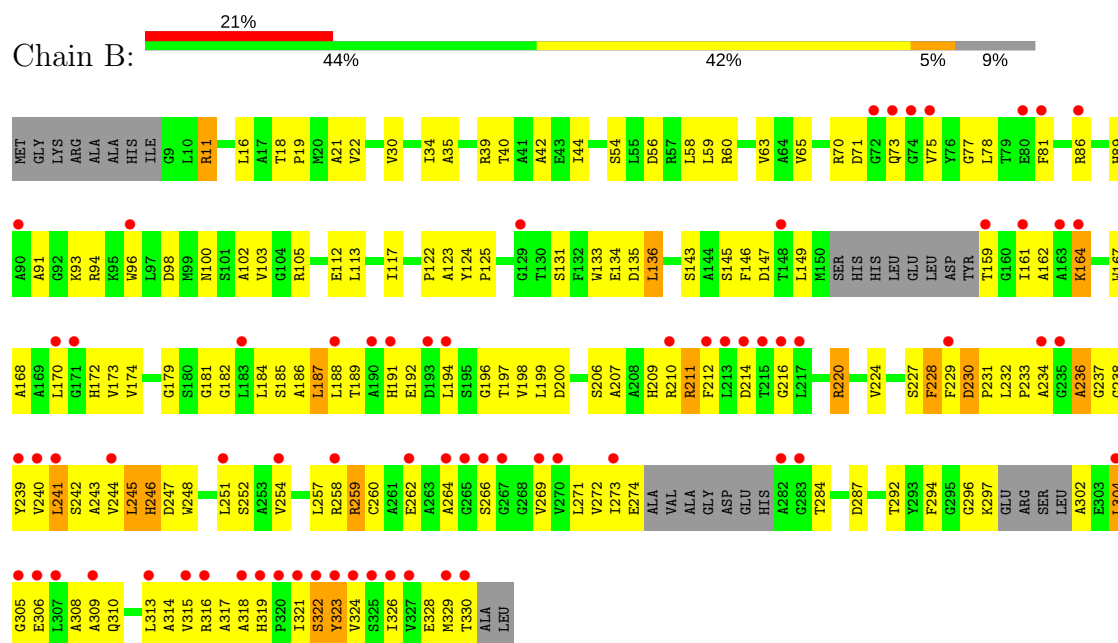
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: O-methyltransferase



• Molecule 1: O-methyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	91.25Å 161.57Å 98.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.08 31.29 – 2.08	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-2.08) 93.6 (31.29-2.08)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.84 (at 2.08Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.240 , 0.272 0.242 , 0.275	Depositor DCC
R_{free} test set	2071 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	32.5	Xtriage
Anisotropy	0.825	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 59.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.016 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.027 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4935	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% 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, SAH

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	0.36	0/2424	0.60	0/3291
1	B	0.33	0/2232	0.62	2/3029 (0.1%)
All	All	0.34	0/4656	0.61	2/6320 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	236	ALA	CB-CA-C	-5.92	101.21	110.10
1	B	266	SER	CB-CA-C	-5.44	99.77	110.10

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	2379	0	2344	70	0
1	B	2193	0	2149	180	0
2	A	26	0	19	0	0
2	B	26	0	19	4	0
3	B	6	0	8	0	0
4	A	188	0	0	7	0
4	B	117	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4935	0	4539	244	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 27.

All (244) 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:34:ILE:HD11	1:A:78:LEU:HG	1.30	1.13
1:B:164:LYS:HG2	1:B:321:ILE:HG22	1.24	1.13
1:B:164:LYS:HG2	1:B:321:ILE:CG2	1.87	1.05
1:B:316:ARG:HG3	1:B:330:THR:HG22	1.41	1.01
1:B:241:LEU:HD13	1:B:245:LEU:HD11	1.44	0.99
1:B:131:SER:OG	1:B:134:GLU:HG2	1.66	0.95
1:B:322:SER:OG	1:B:323:TYR:N	1.99	0.94
1:B:241:LEU:HD11	1:B:272:VAL:HG22	1.51	0.93
1:B:304:LEU:H	1:B:304:LEU:HD12	1.34	0.92
1:B:245:LEU:HD22	1:B:245:LEU:H	1.32	0.92
1:B:162:ALA:HA	1:B:167:TRP:HE1	1.38	0.89
1:A:11:ARG:HG2	1:A:11:ARG:HH11	1.38	0.88
1:B:257:LEU:HD13	1:B:308:ALA:HB1	1.57	0.86
1:A:161:ILE:HD11	1:A:326:ILE:HD11	1.59	0.85
1:B:254:VAL:HA	1:B:257:LEU:HD12	1.59	0.82
1:A:197:THR:HG22	1:A:222:GLN:HG2	1.64	0.80
1:A:34:ILE:HD11	1:A:78:LEU:CG	2.10	0.80
1:A:40:THR:HG23	1:A:43:GLU:H	1.46	0.80
1:B:245:LEU:N	1:B:245:LEU:HD22	1.96	0.80
1:B:246:HIS:HA	1:B:297:LYS:HE2	1.64	0.80
1:B:182:GLY:H	1:B:211:ARG:NH2	1.79	0.79
1:B:309:ALA:HA	1:B:313:LEU:O	1.85	0.75
1:B:241:LEU:HD12	1:B:241:LEU:O	1.87	0.74
1:B:168:ALA:HA	1:B:194:LEU:HD21	1.70	0.74
1:B:309:ALA:HB2	1:B:315:VAL:HG23	1.69	0.74
1:A:34:ILE:HA	1:A:38:HIS:O	1.86	0.74
1:B:322:SER:OG	1:B:323:TYR:CD2	2.40	0.73
1:B:305:GLY:O	1:B:308:ALA:HB3	1.90	0.72
1:B:316:ARG:CG	1:B:330:THR:HG22	2.19	0.71
1:B:162:ALA:HA	1:B:167:TRP:NE1	2.06	0.70
1:B:174:VAL:HG23	1:B:236:ALA:HB2	1.74	0.70
1:B:241:LEU:CD1	1:B:245:LEU:HD11	2.20	0.70
1:B:245:LEU:CD2	1:B:245:LEU:H	2.02	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:304:LEU:N	1:B:304:LEU:HD12	2.08	0.69
1:B:321:ILE:HG12	1:B:324:VAL:O	1.93	0.69
1:B:11:ARG:HG3	1:B:11:ARG:HH11	1.57	0.69
1:B:39:ARG:NH1	1:B:77:GLY:HA2	2.09	0.68
1:B:273:ILE:N	1:B:273:ILE:HD12	2.09	0.68
1:B:182:GLY:HA3	4:B:575:HOH:O	1.93	0.68
1:B:96:TRP:HA	1:B:103:VAL:CG2	2.25	0.67
1:B:321:ILE:HG13	1:B:321:ILE:O	1.93	0.67
1:B:184:LEU:HD11	1:B:196:GLY:HA3	1.78	0.66
1:A:206:SER:O	1:A:210:ARG:HG3	1.96	0.65
1:A:147:ASP:HB2	4:A:773:HOH:O	1.97	0.64
1:B:321:ILE:CG1	1:B:324:VAL:O	2.46	0.64
1:B:96:TRP:CD1	1:B:149:LEU:HD11	2.32	0.64
1:B:185:SER:O	1:B:189:THR:HG23	1.98	0.64
1:B:60:ARG:HH21	1:B:70:ARG:NH2	1.95	0.64
1:B:123:ALA:HB3	1:B:292:THR:HG23	1.79	0.64
1:B:248:TRP:HB3	1:B:252:SER:HB2	1.80	0.63
1:B:248:TRP:HB3	1:B:252:SER:CB	2.27	0.63
1:A:172:HIS:NE2	1:A:197:THR:HG23	2.12	0.63
1:B:102:ALA:CB	1:B:146:PHE:HB2	2.28	0.63
1:B:304:LEU:H	1:B:304:LEU:CD1	2.10	0.63
1:B:234:ALA:HA	1:B:239:TYR:OH	1.99	0.62
1:B:245:LEU:HD23	1:B:274:GLU:OE2	1.98	0.62
1:A:154:LEU:O	1:A:158:TYR:HB2	1.99	0.62
1:A:314:ALA:HB2	1:A:332:LEU:HD13	1.81	0.62
1:B:230:ASP:HB2	1:B:231:PRO:HD2	1.82	0.62
1:A:61:HIS:ND1	1:B:284:THR:HG23	2.15	0.61
1:B:18:THR:O	1:B:22:VAL:HG23	2.00	0.61
1:B:136:LEU:HD13	1:B:143:SER:HA	1.83	0.61
1:B:329:MET:O	1:B:330:THR:HB	1.99	0.61
1:B:316:ARG:HG3	1:B:330:THR:CG2	2.24	0.61
1:B:246:HIS:HD2	1:B:247:ASP:H	1.46	0.61
1:A:6:ALA:HB3	4:A:776:HOH:O	2.00	0.61
1:B:172:HIS:CE1	1:B:197:THR:HG23	2.36	0.60
1:A:40:THR:HG22	1:A:43:GLU:CG	2.32	0.60
1:B:174:VAL:HG23	1:B:236:ALA:CB	2.32	0.60
1:A:136:LEU:HD22	1:A:142:LEU:HB3	1.83	0.60
1:B:315:VAL:O	1:B:316:ARG:HD3	2.03	0.59
1:A:34:ILE:CD1	1:A:78:LEU:HG	2.20	0.59
1:B:242:SER:O	1:B:244:VAL:HG23	2.03	0.59
1:B:164:LYS:HB2	1:B:164:LYS:NZ	2.18	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:39:ARG:HH12	1:B:77:GLY:HA2	1.67	0.59
1:A:306:GLU:HB2	4:A:678:HOH:O	2.02	0.58
1:A:40:THR:CG2	1:A:43:GLU:H	2.16	0.58
1:B:240:VAL:HG13	1:B:273:ILE:HD13	1.84	0.58
1:B:317:ALA:HB3	1:B:328:GLU:HG2	1.86	0.58
1:B:34:ILE:HG21	1:B:78:LEU:HD13	1.85	0.57
1:B:71:ASP:OD1	1:B:73:GLN:HB3	2.05	0.57
1:B:71:ASP:OD2	1:B:75:VAL:HB	2.05	0.57
1:B:241:LEU:HD12	1:B:272:VAL:HA	1.86	0.57
1:B:230:ASP:HB2	1:B:231:PRO:CD	2.35	0.56
1:B:269:VAL:HG12	1:B:330:THR:HB	1.86	0.56
1:B:161:ILE:HG13	1:B:326:ILE:HD11	1.88	0.56
1:B:164:LYS:CG	1:B:321:ILE:HG22	2.17	0.56
1:A:11:ARG:HG2	1:A:11:ARG:NH1	2.15	0.56
1:B:210:ARG:HG3	1:B:211:ARG:N	2.21	0.55
1:B:60:ARG:HH21	1:B:70:ARG:CZ	2.19	0.55
1:B:257:LEU:HD13	1:B:308:ALA:CB	2.33	0.55
1:B:309:ALA:HB1	1:B:314:ALA:HA	1.88	0.55
1:B:269:VAL:O	1:B:271:LEU:HD22	2.06	0.55
1:B:182:GLY:H	1:B:211:ARG:CZ	2.19	0.55
1:B:56:ASP:O	1:B:60:ARG:HG3	2.07	0.55
1:B:251:LEU:O	1:B:254:VAL:HG22	2.07	0.55
1:A:251:LEU:O	1:A:254:VAL:HG22	2.07	0.55
1:B:242:SER:O	2:B:401:SAH:HA	2.07	0.54
1:A:40:THR:HG22	1:A:43:GLU:CD	2.28	0.54
1:B:172:HIS:NE2	1:B:197:THR:HG23	2.22	0.54
1:B:91:ALA:HB1	1:B:93:LYS:HE3	1.91	0.53
1:A:205:ALA:HB1	1:A:223:VAL:CG1	2.39	0.53
1:B:251:LEU:C	1:B:251:LEU:HD23	2.29	0.53
1:B:131:SER:HG	1:B:134:GLU:HG2	1.73	0.52
1:B:136:LEU:CD1	1:B:143:SER:HA	2.37	0.52
1:B:199:LEU:HD23	1:B:224:VAL:HG23	1.92	0.52
1:B:304:LEU:HD23	4:B:748:HOH:O	2.09	0.52
1:B:241:LEU:CD1	1:B:241:LEU:O	2.57	0.52
1:B:96:TRP:NE1	1:B:149:LEU:HD11	2.24	0.52
1:A:89:HIS:HD2	1:A:91:ALA:HB3	1.75	0.52
1:A:152:HIS:O	1:A:156:LEU:HG	2.10	0.52
1:B:122:PRO:HG3	1:B:294:PHE:O	2.10	0.52
1:B:322:SER:OG	1:B:323:TYR:CG	2.64	0.51
1:B:302:ALA:C	1:B:304:LEU:N	2.63	0.51
1:B:192:GLU:HA	1:B:220:ARG:NH2	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:21:ALA:HA	1:B:58:LEU:HD11	1.94	0.50
1:A:199:LEU:HD22	1:A:224:VAL:O	2.12	0.50
1:A:115:HIS:HD2	4:A:741:HOH:O	1.94	0.50
1:B:168:ALA:HA	1:B:194:LEU:CD2	2.40	0.50
1:A:89:HIS:CD2	1:A:91:ALA:H	2.29	0.49
1:A:205:ALA:HB1	1:A:223:VAL:HG11	1.94	0.49
1:B:162:ALA:HB2	1:B:187:LEU:HD12	1.94	0.49
1:B:271:LEU:N	1:B:271:LEU:HD22	2.28	0.49
1:A:10:LEU:HD23	1:B:93:LYS:CB	2.42	0.49
1:B:105:ARG:HH12	1:B:135:ASP:CG	2.15	0.49
1:B:168:ALA:CA	1:B:194:LEU:HD21	2.41	0.49
1:B:145:SER:O	1:B:149:LEU:HB2	2.13	0.48
1:A:299:ARG:NE	1:A:307:LEU:HD11	2.28	0.48
1:B:162:ALA:C	1:B:164:LYS:H	2.16	0.48
1:B:16:LEU:C	1:B:19:PRO:HD2	2.34	0.48
1:A:8:ILE:HG22	1:A:12:ALA:HB3	1.94	0.48
1:A:222:GLN:HA	4:A:534:HOH:O	2.13	0.48
1:B:191:HIS:O	1:B:220:ARG:NH2	2.47	0.48
1:A:199:LEU:C	1:A:199:LEU:HD13	2.34	0.48
1:B:35:ALA:HB2	1:B:86:ARG:HE	1.78	0.48
1:B:30:VAL:HG11	1:B:59:LEU:HD21	1.96	0.47
1:A:34:ILE:HG22	1:A:44:ILE:HD11	1.96	0.47
1:B:302:ALA:C	1:B:304:LEU:H	2.17	0.47
2:B:401:SAH:HB1	2:B:401:SAH:H4'	1.97	0.47
1:B:96:TRP:HA	1:B:103:VAL:HG21	1.96	0.47
1:B:191:HIS:HB2	1:B:194:LEU:HD12	1.95	0.47
1:B:321:ILE:HG13	1:B:324:VAL:O	2.15	0.47
1:A:124:TYR:HB3	1:A:125:PRO:HD3	1.96	0.47
1:B:258:ARG:HG2	1:B:262:GLU:OE1	2.15	0.47
1:B:220:ARG:HD3	4:B:529:HOH:O	2.15	0.47
1:A:276:VAL:O	1:A:278:GLY:N	2.48	0.46
1:B:245:LEU:HG	1:B:272:VAL:HG13	1.96	0.46
1:B:102:ALA:HB1	1:B:146:PHE:HB2	1.97	0.46
1:B:59:LEU:O	1:B:63:VAL:HG23	2.15	0.46
1:A:13:LEU:HD21	1:B:65:VAL:CG1	2.44	0.46
1:B:73:GLN:HA	4:B:602:HOH:O	2.15	0.46
1:A:89:HIS:CD2	1:A:91:ALA:HB3	2.50	0.46
1:A:323:TYR:CD2	1:A:323:TYR:N	2.84	0.46
1:B:11:ARG:CG	1:B:11:ARG:HH11	2.24	0.46
1:A:110:PHE:CD2	1:A:292:THR:HG21	2.51	0.45
1:B:244:VAL:O	1:B:248:TRP:HD1	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:167:TRP:HZ3	1:B:240:VAL:HG21	1.82	0.45
1:B:246:HIS:O	1:B:296:GLY:HA3	2.15	0.45
1:B:246:HIS:HD2	1:B:247:ASP:N	2.14	0.45
1:B:241:LEU:CD1	1:B:272:VAL:HA	2.47	0.45
1:B:309:ALA:CB	1:B:314:ALA:HA	2.46	0.45
1:B:246:HIS:CD2	1:B:247:ASP:H	2.31	0.45
1:B:181:GLY:HA3	1:B:211:ARG:NE	2.31	0.45
1:B:317:ALA:HB3	1:B:328:GLU:CG	2.47	0.45
1:B:260:CYS:O	1:B:264:ALA:N	2.50	0.44
1:A:58:LEU:HD23	1:A:58:LEU:C	2.37	0.44
1:B:246:HIS:CA	1:B:297:LYS:HE2	2.41	0.44
1:B:297:LYS:HD2	4:B:748:HOH:O	2.16	0.44
1:B:167:TRP:CZ3	1:B:240:VAL:HG21	2.52	0.44
1:A:16:LEU:C	1:A:19:PRO:HD2	2.37	0.44
1:A:161:ILE:HD13	1:A:321:ILE:HD12	1.99	0.44
1:A:27:THR:HG21	1:A:99:MET:HE2	2.00	0.44
1:A:120:GLY:HA2	1:B:54:SER:OG	2.18	0.44
1:B:273:ILE:CD1	1:B:273:ILE:N	2.78	0.44
1:B:318:ALA:O	1:B:319:HIS:ND1	2.47	0.44
1:B:168:ALA:N	1:B:194:LEU:HD11	2.33	0.44
1:B:98:ASP:OD2	1:B:100:ASN:HB2	2.18	0.44
1:A:34:ILE:HD11	1:A:78:LEU:CD2	2.48	0.44
1:B:186:ALA:HA	1:B:189:THR:OG1	2.18	0.44
1:B:179:GLY:N	1:B:200:ASP:OD2	2.48	0.44
1:B:244:VAL:H	1:B:245:LEU:HD22	1.82	0.44
1:A:34:ILE:HG22	1:A:44:ILE:CD1	2.48	0.43
1:B:206:SER:O	1:B:209:HIS:HB3	2.18	0.43
1:A:161:ILE:CD1	1:A:326:ILE:HD11	2.40	0.43
1:A:136:LEU:CD2	1:A:142:LEU:HB3	2.48	0.43
1:B:232:LEU:HG	1:B:233:PRO:HD2	2.00	0.43
1:B:251:LEU:C	1:B:251:LEU:CD2	2.86	0.43
1:B:246:HIS:CB	1:B:297:LYS:HG2	2.49	0.43
1:B:184:LEU:HD23	1:B:198:VAL:HG23	2.01	0.43
1:B:188:LEU:HD22	1:B:194:LEU:O	2.18	0.43
1:A:172:HIS:CE1	1:A:197:THR:HG23	2.52	0.43
1:A:222:GLN:HB2	4:A:725:HOH:O	2.18	0.43
1:B:243:ALA:HB3	2:B:401:SAH:HG1	2.00	0.43
1:B:243:ALA:HB3	2:B:401:SAH:N	2.33	0.43
1:B:269:VAL:O	1:B:269:VAL:HG23	2.19	0.43
1:B:112:GLU:HB2	1:B:123:ALA:HB1	2.01	0.43
1:B:89:HIS:HD2	1:B:91:ALA:HB3	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:246:HIS:CD2	1:B:247:ASP:N	2.87	0.43
1:A:57:ARG:HD2	1:B:287:ASP:OD2	2.19	0.42
1:B:228:PHE:HE2	1:B:229:PHE:CE2	2.37	0.42
1:B:259:ARG:HG3	1:B:259:ARG:HH11	1.84	0.42
1:A:188:LEU:CD2	1:A:221:ALA:HB2	2.49	0.42
1:B:207:ALA:O	1:B:210:ARG:HG2	2.20	0.42
1:A:131:SER:OG	1:A:134:GLU:HG2	2.19	0.42
1:A:11:ARG:NH1	1:B:96:TRP:CD2	2.88	0.42
1:A:139:ASP:OD1	1:A:141:VAL:N	2.51	0.42
1:B:259:ARG:HG3	1:B:259:ARG:NH1	2.34	0.42
1:B:241:LEU:HD13	1:B:245:LEU:CD1	2.32	0.42
1:A:209:HIS:CE1	1:A:213:LEU:HD11	2.55	0.42
1:B:316:ARG:HE	1:B:330:THR:CG2	2.33	0.42
1:B:113:LEU:O	1:B:117:ILE:HG12	2.20	0.41
1:B:214:ASP:C	1:B:216:GLY:H	2.22	0.41
1:B:133:TRP:CE3	1:B:133:TRP:HA	2.55	0.41
1:B:105:ARG:NH1	1:B:124:TYR:OH	2.54	0.41
1:B:187:LEU:HA	1:B:187:LEU:HD12	1.87	0.41
1:B:170:LEU:HD22	1:B:237:GLY:O	2.20	0.41
1:A:6:ALA:HB1	1:B:81:PHE:CE2	2.56	0.41
1:B:306:GLU:O	1:B:310:GLN:N	2.50	0.41
1:B:40:THR:HG22	1:B:42:ALA:N	2.35	0.41
1:A:157:ASP:C	1:A:158:TYR:HD1	2.24	0.41
1:A:276:VAL:C	1:A:278:GLY:N	2.73	0.41
1:A:11:ARG:NH1	1:A:11:ARG:CG	2.79	0.41
1:A:172:HIS:HA	1:A:195:SER:O	2.21	0.41
1:B:122:PRO:HG3	1:B:294:PHE:C	2.41	0.41
1:B:173:VAL:HG23	1:B:238:GLY:O	2.21	0.41
1:B:199:LEU:C	1:B:199:LEU:HD13	2.41	0.41
1:A:249:ASP:OD1	1:A:251:LEU:HB3	2.20	0.41
1:B:124:TYR:HB3	1:B:125:PRO:HD3	2.03	0.41
1:B:251:LEU:O	1:B:252:SER:C	2.58	0.41
1:B:227:SER:O	1:B:230:ASP:OD1	2.39	0.41
1:B:159:THR:C	1:B:161:ILE:N	2.74	0.41
1:B:184:LEU:HD23	1:B:212:PHE:CZ	2.56	0.41
1:B:40:THR:O	1:B:44:ILE:HG13	2.21	0.41
1:A:180:SER:O	1:A:208:ALA:HA	2.21	0.40
1:A:323:TYR:H	1:A:323:TYR:HD2	1.68	0.40
1:B:273:ILE:HG22	1:B:274:GLU:N	2.36	0.40
1:B:89:HIS:CD2	1:B:91:ALA:HB3	2.57	0.40
1:A:55:LEU:O	1:A:59:LEU:HG	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127:ARG:HG3	4:A:523:HOH:O	2.21	0.40
1:A:22:VAL:HG12	1:A:97:LEU:HD23	2.03	0.40
1:A:112:GLU:HB2	1:A:123:ALA:HB1	2.03	0.40
1:B:236:ALA:HA	4:B:676:HOH:O	2.21	0.40
1:B:302:ALA:O	1:B:304:LEU:N	2.53	0.40
1:B:89:HIS:CD2	1:B:91:ALA:H	2.40	0.40

There are no symmetry-related clashes.

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	321/332 (97%)	313 (98%)	6 (2%)	2 (1%)	28	23
1	B	295/332 (89%)	262 (89%)	33 (11%)	0	100	100
All	All	616/664 (93%)	575 (93%)	39 (6%)	2 (0%)	44	43

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	322	SER
1	A	277	ALA

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	231/237 (98%)	224 (97%)	7 (3%)	46	48
1	B	210/237 (89%)	193 (92%)	17 (8%)	14	9
All	All	441/474 (93%)	417 (95%)	24 (5%)	26	22

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	153	HIS
1	A	157	ASP
1	A	193	ASP
1	A	214	ASP
1	A	218	SER
1	A	228	PHE
1	A	294	PHE
1	B	11	ARG
1	B	94	ARG
1	B	136	LEU
1	B	147	ASP
1	B	164	LYS
1	B	187	LEU
1	B	211	ARG
1	B	220	ARG
1	B	228	PHE
1	B	230	ASP
1	B	241	LEU
1	B	245	LEU
1	B	246	HIS
1	B	259	ARG
1	B	304	LEU
1	B	322	SER
1	B	323	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	HIS
1	A	89	HIS
1	A	209	HIS
1	B	51	HIS
1	B	73	GLN
1	B	89	HIS

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Mol	Chain	Res	Type
1	B	115	HIS
1	B	121	GLN
1	B	246	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](#)

3 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SAH	A	401	-	20,28,28	0.74	0	20,40,40	2.23	3 (15%)
2	SAH	B	401	-	20,28,28	0.75	0	20,40,40	2.22	4 (20%)
3	GOL	B	402	-	5,5,5	0.27	0	5,5,5	0.23	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
2	SAH	A	401	-	-	0/7/31/31	0/3/3/3
2	SAH	B	401	-	-	0/7/31/31	0/3/3/3
3	GOL	B	402	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	SAH	N3-C2-N1	-8.40	121.54	128.86
2	B	401	SAH	N3-C2-N1	-8.25	121.67	128.86
2	A	401	SAH	C4'-O4'-C1'	-3.91	105.61	109.77
2	B	401	SAH	C4'-O4'-C1'	-3.84	105.68	109.77
2	A	401	SAH	C4-C5-N7	-2.33	107.16	109.41
2	B	401	SAH	C5'-C4'-C3'	-2.16	109.54	115.05
2	B	401	SAH	C4-C5-N7	-2.01	107.47	109.41

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	SAH	4	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	325/332 (97%)	0.23	11 (3%) 46 51	25, 37, 53, 72	0
1	B	303/332 (91%)	1.11	70 (23%) 1 1	27, 53, 83, 89	0
All	All	628/664 (94%)	0.66	81 (12%) 4 5	25, 42, 75, 89	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	323	TYR	10.0
1	B	266	SER	7.6
1	B	321	ILE	7.5
1	B	322	SER	7.2
1	B	234	ALA	6.8
1	B	313	LEU	6.6
1	B	282	ALA	6.1
1	B	73	GLN	5.2
1	B	330	THR	5.2
1	B	161	ILE	4.6
1	B	190	ALA	4.1
1	B	216	GLY	3.9
1	A	323	TYR	3.9
1	B	318	ALA	3.9
1	A	210	ARG	3.8
1	B	148	THR	3.8
1	B	251	LEU	3.7
1	B	316	ARG	3.6
1	B	329	MET	3.6
1	B	265	GLY	3.6
1	B	183	LEU	3.5
1	B	315	VAL	3.5
1	B	320	PRO	3.3
1	B	327	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	7	HIS	3.2
1	B	72	GLY	3.1
1	B	164	LYS	3.1
1	B	215	THR	3.1
1	B	309	ALA	3.0
1	B	267	GLY	3.0
1	B	217	LEU	3.0
1	B	212	PHE	2.9
1	B	96	TRP	2.9
1	A	8	ILE	2.8
1	B	86	ARG	2.8
1	B	305	GLY	2.8
1	A	332	LEU	2.8
1	B	258	ARG	2.8
1	B	171	GLY	2.7
1	B	188	LEU	2.7
1	B	262	GLU	2.7
1	B	194	LEU	2.7
1	B	269	VAL	2.6
1	B	129	GLY	2.6
1	B	264	ALA	2.6
1	B	75	VAL	2.6
1	B	244	VAL	2.6
1	A	193	ASP	2.5
1	A	278	GLY	2.5
1	B	325	SER	2.5
1	B	273	ILE	2.5
1	B	326	ILE	2.5
1	B	270	VAL	2.4
1	B	81	PHE	2.4
1	B	240	VAL	2.4
1	B	241	LEU	2.3
1	B	304	LEU	2.3
1	B	191	HIS	2.3
1	B	163	ALA	2.3
1	B	159	THR	2.3
1	B	74	GLY	2.3
1	A	88	ASP	2.3
1	A	277	ALA	2.2
1	A	130	THR	2.2
1	B	324	VAL	2.2
1	B	213	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	239	TYR	2.1
1	B	210	ARG	2.1
1	B	283	GLY	2.1
1	A	288	LEU	2.0
1	B	307	LEU	2.0
1	B	90	ALA	2.0
1	B	254	VAL	2.0
1	B	319	HIS	2.0
1	B	306	GLU	2.0
1	B	170	LEU	2.0
1	B	229	PHE	2.0
1	B	235	GLY	2.0
1	B	80	GLU	2.0
1	B	193	ASP	2.0
1	B	214	ASP	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(\AA^2)	Q<0.9
2	SAH	A	401	26/26	0.95	0.13	-0.11	22,28,34,38	0
2	SAH	B	401	26/26	0.90	0.16	-0.14	34,37,59,61	0
3	GOL	B	402	6/6	0.46	0.18	-	69,72,73,73	0

6.5 Other polymers [i](#)

There are no such residues in this entry.