



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:52 PM GMT

PDB ID : 1HNN  
Title : CRYSTAL STRUCTURE OF HUMAN PNMT COMPLEXED WITH SK&F  
29661 AND ADOHCY(SAH)  
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Deposited on : 2000-12-07  
Resolution : 2.40 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
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) : trunk26865

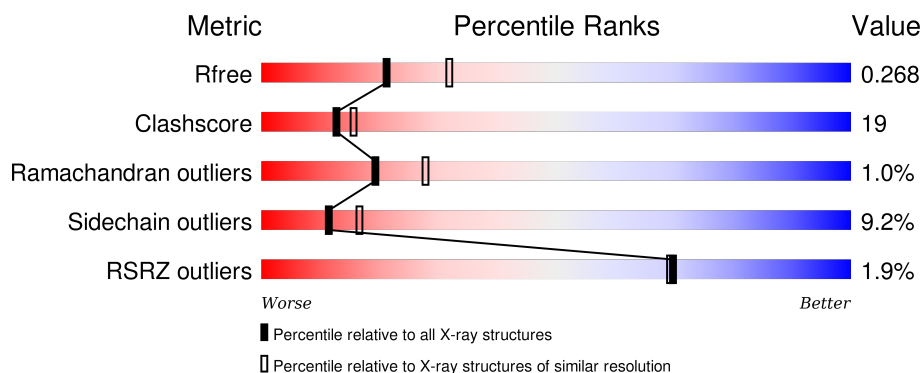
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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  $\geq 3$ , 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	282	<div> <div>2%</div> <div>56% 32% 5% 7%</div> </div>
1	B	282	<div> <div>%</div> <div>60% 30% • 5%</div> </div>

## 2 Entry composition [i](#)

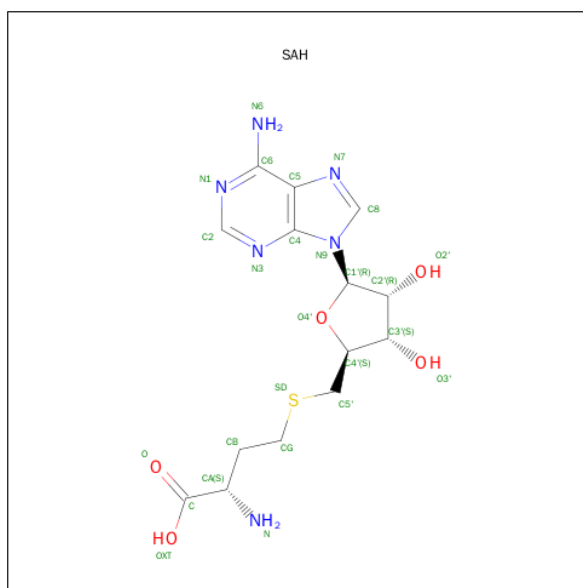
There are 4 unique types of molecules in this entry. The entry contains 4360 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PHENYLETHANOLAMINE N-METHYLTRANSFERASE.

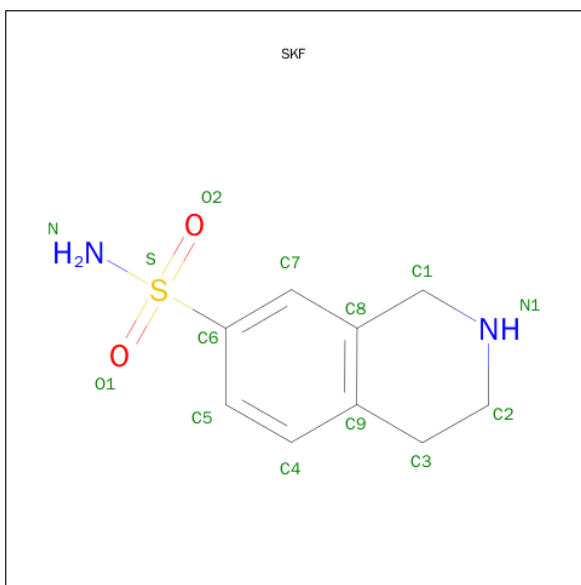
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	261	Total	C	N	O	S	0	6	0
			2055	1304	362	378	11			
1	B	267	Total	C	N	O	S	0	5	0
			2099	1331	369	389	10			

- 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
2	A	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
2	B	1	Total	C	N	O	S	0	0
			26	14	6	5	1		

- Molecule 3 is 1,2,3,4-TETRAHYDRO-ISOQUINOLINE-7-SULFONIC ACID AMIDE (three-letter code: SKF) (formula:  $C_9H_{12}N_2O_2S$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			14	9	2	2	1		
3	B	1	Total	C	N	O	S	0	0
			14	9	2	2	1		

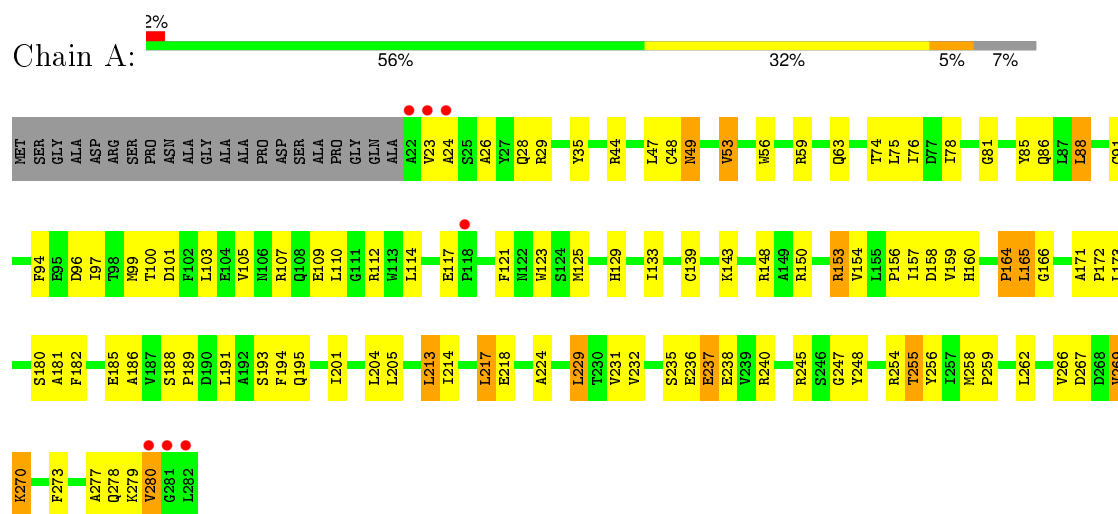
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	66	Total	O	0	0
			66	66		
4	B	60	Total	O	0	0
			60	60		

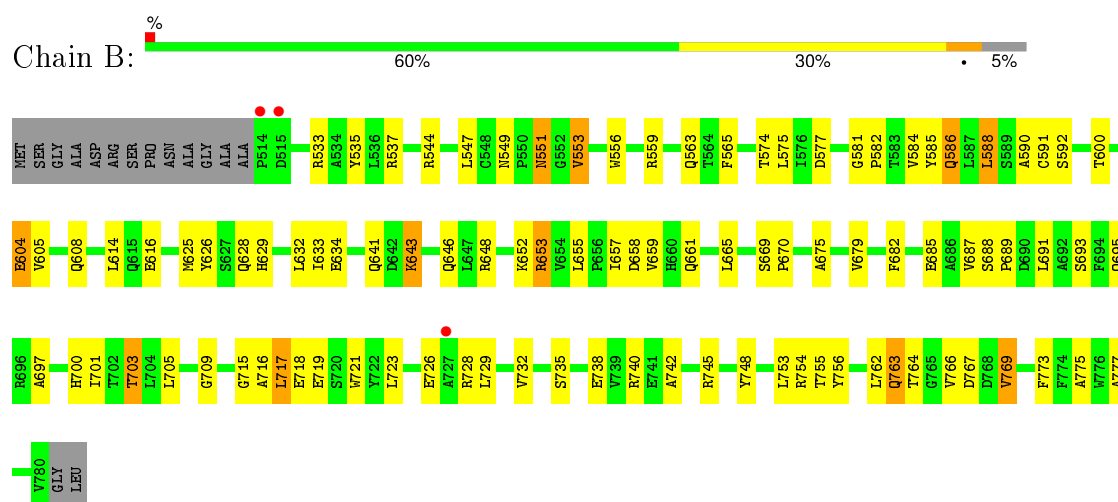
### 3 Residue-property plots

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: PHENYLETHANOLAMINE N-METHYLTRANSFERASE



#### • Molecule 1: PHENYLETHANOLAMINE N-METHYLTRANSFERASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	94.35Å 94.35Å 187.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.48 – 2.40 38.48 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.0 (38.48-2.40) 90.5 (38.48-2.30)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.63 (at 2.29Å)	Xtriage
Refinement program	CNS 0.9	Depositor
R, $R_{free}$	0.230 , 0.273 0.230 , 0.268	Depositor DCC
$R_{free}$ test set	3236 reflections (10.03%)	DCC
Wilson B-factor (Å <sup>2</sup> )	42.2	Xtriage
Anisotropy	0.357	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 30.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 34796 reflections (0.003%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4360	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: SAH, SKF

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	0.39	0/2108	0.66	0/2875
1	B	0.41	0/2152	0.66	0/2932
All	All	0.40	0/4260	0.66	0/5807

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2055	0	1985	90	0
1	B	2099	0	2050	79	0
2	A	26	0	19	5	0
2	B	26	0	19	3	0
3	A	14	0	12	2	0
3	B	14	0	12	3	0
4	A	66	0	0	2	0
4	B	60	0	0	1	0
All	All	4360	0	4097	161	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 19.

All (161) 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:B:559:ARG:HG2	1:B:563:GLN:HE21	1.25	1.01
1:A:237:GLU:HG3	1:B:559:ARG:HH12	1.26	0.99
1:B:695:GLN:NE2	1:B:742:ALA:HA	1.86	0.91
1:A:188:SER:HB2	1:A:193:SER:HB2	1.51	0.90
1:B:719:GLU:OE1	1:B:769:VAL:HG13	1.73	0.87
1:A:188:SER:O	1:A:232:VAL:HG23	1.76	0.86
1:A:75:LEU:HD23	1:A:76:ILE:N	1.96	0.80
1:A:53:VAL:HG13	1:A:256:TYR:CE2	2.19	0.78
1:A:237:GLU:HG3	1:B:559:ARG:NH1	1.99	0.78
1:A:195:GLN:HE22	1:A:245:ARG:CB	1.96	0.78
1:B:700:HIS:O	1:B:703:THR:HB	1.85	0.76
1:A:44:ARG:HD3	3:A:3001:SKF:N	2.01	0.74
1:A:217:LEU:HD13	1:A:273:PHE:HD2	1.53	0.73
1:B:735:SER:OG	1:B:738:GLU:HG3	1.90	0.71
1:A:182:PHE:HB2	2:A:2001:SAH:HB2	1.72	0.71
1:A:48[B]:CYS:O	1:A:49[B]:ASN:HB3	1.90	0.71
1:A:254:ARG:HG2	1:B:740:ARG:NH2	2.07	0.69
1:A:235:SER:OG	1:A:238:GLU:HG3	1.93	0.69
1:A:35:TYR:OH	2:A:2001:SAH:HB1	1.92	0.68
1:B:585:TYR:HA	1:B:588:LEU:HD22	1.75	0.68
1:A:171:ALA:O	1:A:173:LEU:HD23	1.94	0.67
1:A:59:ARG:O	1:A:63:GLN:HG3	1.95	0.66
1:B:553:VAL:HG22	1:B:762:LEU:HD13	1.78	0.65
1:B:695:GLN:HE22	1:B:745:ARG:HB2	1.61	0.65
1:B:659:VAL:CG1	1:B:697:ALA:HB1	2.27	0.65
1:B:653:ARG:HG3	1:B:655:LEU:HG	1.78	0.65
1:B:717:LEU:HD13	1:B:773:PHE:HD2	1.62	0.63
1:B:682:PHE:HB2	2:B:2002:SAH:HB2	1.81	0.62
1:A:217:LEU:HD13	1:A:273:PHE:CD2	2.35	0.62
1:B:688:SER:HB2	1:B:693:SER:HB3	1.80	0.61
1:B:559:ARG:HG2	1:B:563:GLN:NE2	2.08	0.61
1:B:695:GLN:HE22	1:B:742:ALA:HA	1.65	0.60
1:A:53:VAL:HG13	1:A:256:TYR:CD2	2.37	0.60
1:A:109:GLU:HA	1:A:112:ARG:NH1	2.17	0.59
1:B:659:VAL:HG22	1:B:701:ILE:HD11	1.83	0.59
1:A:48[B]:CYS:O	1:A:49[B]:ASN:CB	2.52	0.58
1:B:717:LEU:HD13	1:B:773:PHE:CD2	2.39	0.58
1:A:107:ARG:NH1	1:A:156:PRO:HB3	2.19	0.57
1:B:575:LEU:HD23	1:B:575:LEU:C	2.24	0.57
1:A:255:THR:CG2	1:B:755:THR:HB	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:ILE:HA	1:A:204:LEU:HD12	1.86	0.57
1:A:44:ARG:HD3	3:A:3001:SKF:H1N	1.67	0.57
1:A:173:LEU:HD11	1:A:204:LEU:HD23	1.87	0.57
1:A:237:GLU:CG	1:B:559:ARG:NH1	2.67	0.57
1:B:723:LEU:HD23	1:B:728:ARG:HA	1.86	0.56
1:A:53:VAL:HG22	1:A:262:LEU:HD13	1.87	0.56
1:A:157:ILE:HA	1:A:165:LEU:HB3	1.86	0.56
1:B:753:LEU:HD13	1:B:775:ALA:HB2	1.85	0.56
1:A:229:LEU:H	1:A:229:LEU:HD22	1.70	0.56
1:A:236:GLU:OE2	1:A:255:THR:HG22	2.05	0.56
1:A:267:ASP:OD1	1:A:269:VAL:HG23	2.05	0.56
1:B:716:ALA:HB1	1:B:769:VAL:HG12	1.88	0.56
1:A:255:THR:HG21	1:B:755:THR:HB	1.87	0.56
1:A:229:LEU:N	1:A:229:LEU:HD22	2.20	0.56
1:B:535:TYR:OH	2:B:2002:SAH:HB1	2.06	0.55
1:A:109:GLU:HA	1:A:112:ARG:HH11	1.71	0.55
1:A:188:SER:O	1:A:232:VAL:CG2	2.51	0.55
1:B:547:LEU:H	1:B:625[A]:MET:HE2	1.71	0.55
1:B:553:VAL:CG2	1:B:762:LEU:HD13	2.37	0.55
1:B:658:ASP:HB3	1:B:661:GLN:HG2	1.89	0.54
1:B:553:VAL:HG13	1:B:756:TYR:CE2	2.43	0.54
1:A:129[A]:HIS:CD2	1:A:133:ILE:HD11	2.42	0.54
1:B:767:ASP:OD1	3:B:3002:SKF:HC11	2.08	0.54
1:B:769:VAL:HG11	3:B:3002:SKF:HC12	1.89	0.53
1:A:91:CYS:O	1:A:150:ARG:NE	2.39	0.53
1:B:695:GLN:NE2	1:B:745:ARG:HD2	2.24	0.53
1:A:96:ASP:OD2	1:A:153:ARG:NH2	2.43	0.52
1:A:85:TYR:HA	1:A:88:LEU:HD22	1.92	0.52
1:B:748:TYR:HB3	1:B:777:ALA:HB1	1.91	0.51
1:B:767:ASP:OD2	1:B:769:VAL:HG22	2.10	0.51
1:B:763:GLN:HA	1:B:763:GLN:HE21	1.76	0.51
1:B:687:VAL:O	1:B:687:VAL:HG12	2.10	0.51
1:A:153:ARG:NH2	1:A:172:PRO:HD3	2.26	0.51
1:A:154:VAL:O	1:A:154:VAL:HG23	2.11	0.51
1:A:159:VAL:HG22	2:A:2001:SAH:N1	2.26	0.51
1:A:112:ARG:HA	1:A:117:GLU:HB2	1.93	0.50
1:B:559:ARG:NE	1:B:563:GLN:NE2	2.59	0.50
1:A:180:SER:HB3	1:A:213:LEU:HD23	1.93	0.50
1:B:721:TRP:HA	1:B:729:LEU:O	2.11	0.50
1:A:247:GLY:O	1:A:280:VAL:CG1	2.59	0.50
1:A:139:CYS:HA	4:A:1123:HOH:O	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:634:GLU:OE1	1:B:643:LYS:NZ	2.45	0.49
1:A:121:PHE:HD2	1:A:123:TRP:CE2	2.30	0.49
1:A:254:ARG:HG2	1:B:740:ARG:HH22	1.77	0.49
1:B:556:TRP:CE2	1:B:756:TYR:HB2	2.48	0.49
1:A:110:LEU:O	1:A:114:LEU:HD23	2.12	0.48
1:A:247:GLY:O	1:A:280:VAL:HG12	2.13	0.48
1:A:218:GLU:HB2	1:A:270:LYS:HB2	1.95	0.48
1:A:240:ARG:NH1	1:B:754:ARG:HD3	2.29	0.48
1:B:565:PHE:CZ	1:B:590:ALA:HB2	2.49	0.47
1:A:78:ILE:HA	1:A:100:THR:OG1	2.14	0.47
1:B:559:ARG:CG	1:B:563:GLN:HE21	2.10	0.47
1:B:547:LEU:H	1:B:625[A]:MET:CE	2.28	0.47
1:A:237:GLU:OE2	1:A:237:GLU:HA	2.14	0.47
1:B:705:LEU:HD11	1:B:709:GLY:HA3	1.96	0.47
1:B:549:ASN:OD1	1:B:551[A]:ASN:ND2	2.48	0.47
1:A:248:TYR:HB3	1:A:277:ALA:HB1	1.97	0.47
1:A:143:LYS:HG3	4:A:1006:HOH:O	2.15	0.46
1:B:732:VAL:HG13	1:B:732:VAL:O	2.15	0.46
1:A:121:PHE:CD2	1:A:123:TRP:CE2	3.03	0.46
1:A:99:MET:CE	1:A:110:LEU:HD13	2.46	0.46
1:B:577:ASP:OD1	1:B:679[B]:VAL:HG23	2.15	0.46
1:A:114:LEU:O	1:A:148:ARG:HD3	2.17	0.45
1:B:653:ARG:HD2	1:B:655:LEU:HD11	1.99	0.45
1:B:669:SER:OG	1:B:670:PRO:HD2	2.15	0.45
1:A:181:ALA:HA	1:A:214:ILE:O	2.17	0.45
1:B:652:LYS:O	1:B:653:ARG:HB3	2.16	0.45
1:B:600:THR:HA	1:B:655:LEU:O	2.17	0.45
1:B:688:SER:HA	1:B:689:PRO:HD3	1.78	0.45
1:B:764:THR:OG1	1:B:766:VAL:HG22	2.17	0.45
1:A:23:VAL:HG23	1:A:24:ALA:N	2.32	0.45
1:A:266:VAL:O	1:A:267:ASP:HB3	2.17	0.44
1:B:691:LEU:HD22	1:B:738:GLU:OE1	2.17	0.44
1:A:81:GLY:HA2	2:A:2001:SAH:O3'	2.17	0.44
1:B:581:GLY:HA2	2:B:2002:SAH:O3'	2.18	0.44
1:A:280:VAL:HG22	1:A:280:VAL:O	2.17	0.44
1:A:158:ASP:OD1	1:A:160:HIS:HB2	2.17	0.44
1:B:657:ILE:HG12	1:B:658:ASP:N	2.32	0.44
1:A:100:THR:O	1:A:101:ASP:HB2	2.18	0.44
1:B:634:GLU:OE2	1:B:643:LYS:NZ	2.50	0.44
1:B:715:GLY:HA3	1:B:773:PHE:CE2	2.53	0.44
1:B:547:LEU:HD12	1:B:625[B]:MET:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:182:PHE:O	2:A:2001:SAH:HG1	2.18	0.43
1:A:110:LEU:C	1:A:114:LEU:HD23	2.39	0.43
1:A:191:LEU:O	1:A:194:PHE:HB3	2.18	0.43
1:B:629:HIS:CD2	1:B:633:ILE:HD11	2.53	0.43
1:A:24:ALA:O	1:A:28:GLN:HG2	2.18	0.43
1:A:186:ALA:HA	1:A:231:VAL:CG1	2.48	0.43
1:A:56:TRP:CE2	1:A:256:TYR:HB2	2.54	0.43
1:A:53:VAL:HG22	1:A:262:LEU:CD1	2.48	0.43
1:B:604:GLU:O	1:B:608:GLN:HG3	2.19	0.42
1:A:237:GLU:OE2	1:A:237:GLU:CA	2.67	0.42
1:A:157:ILE:HG12	1:A:158:ASP:N	2.34	0.42
1:A:189:PRO:C	1:A:232:VAL:HG22	2.40	0.42
1:B:604:GLU:HG3	1:B:605:VAL:N	2.34	0.42
1:A:75:LEU:HD12	1:A:94:PHE:CD1	2.54	0.42
1:A:267:ASP:OD1	1:A:269:VAL:CG2	2.66	0.42
1:A:110:LEU:O	1:A:114:LEU:CD2	2.67	0.42
1:A:188:SER:HB2	1:A:193:SER:CB	2.36	0.42
1:A:47:LEU:HD12	1:A:125[B]:MET:C	2.40	0.42
1:B:753:LEU:CD1	1:B:775:ALA:HB2	2.49	0.42
1:A:91:CYS:HA	1:A:97:ILE:HD11	2.01	0.42
1:A:47:LEU:HD12	1:A:125[A]:MET:HB3	2.01	0.42
1:A:103:LEU:HB3	1:A:105:VAL:HG22	2.02	0.42
1:B:586:GLN:HE21	1:B:586:GLN:HB2	1.69	0.42
1:B:551[A]:ASN:HD22	1:B:551[A]:ASN:C	2.23	0.42
1:A:53:VAL:CG2	1:A:262:LEU:HD13	2.50	0.41
1:B:687:VAL:CG1	1:B:687:VAL:O	2.68	0.41
1:B:559:ARG:NE	1:B:563:GLN:HE22	2.18	0.41
1:A:258:MET:HA	1:A:259:PRO:HD3	1.88	0.41
1:B:537:ARG:HG3	1:B:537:ARG:HH11	1.85	0.41
1:B:559:ARG:HD2	4:B:1126:HOH:O	2.21	0.41
1:A:23:VAL:O	1:A:26:ALA:N	2.54	0.41
1:A:205:LEU:O	1:A:279:LYS:HD3	2.20	0.41
1:B:616:GLU:HG3	1:B:648:ARG:NH2	2.36	0.41
1:B:688:SER:HB2	1:B:693:SER:CB	2.49	0.41
1:B:574:THR:CG2	1:B:675:ALA:HA	2.50	0.41
1:B:544:ARG:HD3	3:B:3002:SKF:N	2.36	0.41
1:A:248:TYR:HA	1:A:278:GLN:O	2.21	0.41
1:A:164:PRO:O	1:A:166:GLY:N	2.55	0.40
1:B:547:LEU:CD1	1:B:626:TYR:HA	2.52	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	265/282 (94%)	244 (92%)	16 (6%)	5 (2%)	10	12
1	B	270/282 (96%)	252 (93%)	17 (6%)	1 (0%)	39	56
All	All	535/564 (95%)	496 (93%)	33 (6%)	6 (1%)	19	25

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	165	LEU
1	A	164	PRO
1	A	49[A]	ASN
1	A	49[B]	ASN
1	A	224	ALA
1	B	592	SER

### 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	213/226 (94%)	198 (93%)	15 (7%)	19	29
1	B	219/226 (97%)	194 (89%)	25 (11%)	7	9
All	All	432/452 (96%)	392 (91%)	40 (9%)	11	16

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	ARG
1	A	53	VAL
1	A	74	THR
1	A	86	GLN
1	A	88	LEU
1	A	153	ARG
1	A	185	GLU
1	A	213	LEU
1	A	217	LEU
1	A	229	LEU
1	A	237	GLU
1	A	255	THR
1	A	269	VAL
1	A	270	LYS
1	A	280	VAL
1	B	533	ARG
1	B	551[A]	ASN
1	B	551[B]	ASN
1	B	553	VAL
1	B	582	PRO
1	B	584	VAL
1	B	586	GLN
1	B	588	LEU
1	B	591	CYS
1	B	604	GLU
1	B	614	LEU
1	B	628	GLN
1	B	632	LEU
1	B	641	GLN
1	B	643	LYS
1	B	646	GLN
1	B	653	ARG
1	B	665	LEU
1	B	685	GLU
1	B	703	THR
1	B	717	LEU
1	B	718	GLU
1	B	726	GLU
1	B	763	GLN
1	B	769	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	195	GLN
1	B	563	GLN
1	B	629	HIS
1	B	646	GLN
1	B	695	GLN
1	B	763	GLN

### 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	$\# Z  > 2$
2	SAH	A	2001	-	20,28,28	1.19	1 (5%)	19,40,40	2.09	4 (21%)
3	SKF	A	3001	-	15,15,15	0.65	0	21,22,22	2.05	3 (14%)
2	SAH	B	2002	-	20,28,28	1.20	2 (10%)	19,40,40	2.00	4 (21%)
3	SKF	B	3002	-	15,15,15	0.64	0	21,22,22	2.01	3 (14%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means

no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SAH	A	2001	-	-	0/7/31/31	0/3/3/3
3	SKF	A	3001	-	-	0/6/13/13	0/2/2/2
2	SAH	B	2002	-	-	0/7/31/31	0/3/3/3
3	SKF	B	3002	-	-	0/6/13/13	0/2/2/2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2002	SAH	C8-N7	-2.53	1.29	1.34
2	A	2001	SAH	C3'-C4'	2.97	1.61	1.53
2	B	2002	SAH	C3'-C4'	2.97	1.61	1.53

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	3001	SKF	O2-S-O1	-7.02	108.93	118.80
3	B	3002	SKF	O2-S-O1	-6.95	109.03	118.80
2	A	2001	SAH	C4'-O4'-C1'	-6.22	102.89	109.72
2	B	2002	SAH	C4'-O4'-C1'	-5.97	103.16	109.72
2	B	2002	SAH	C5'-SD-CG	-2.64	94.48	102.41
2	A	2001	SAH	O4'-C4'-C5'	-2.39	102.35	108.85
2	B	2002	SAH	O4'-C4'-C5'	-2.27	102.68	108.85
2	A	2001	SAH	O4'-C1'-N9	-2.26	103.37	108.10
2	B	2002	SAH	C2'-C1'-N9	2.54	118.18	114.29
3	B	3002	SKF	O1-S-C6	2.89	110.96	107.39
3	A	3001	SKF	O1-S-C6	2.90	110.96	107.39
2	A	2001	SAH	C2'-C1'-N9	3.43	119.53	114.29
3	B	3002	SKF	O2-S-C6	4.03	112.36	107.39
3	A	3001	SKF	O2-S-C6	4.24	112.62	107.39

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2001	SAH	5	0
3	A	3001	SKF	2	0
2	B	2002	SAH	3	0
3	B	3002	SKF	3	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	261/282 (92%)	-0.21	7 (2%) 58 57	31, 50, 70, 84	0
1	B	267/282 (94%)	-0.31	3 (1%) 82 82	29, 45, 58, 86	0
All	All	528/564 (93%)	-0.26	10 (1%) 70 69	29, 47, 66, 86	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	282	LEU	6.3
1	A	281	GLY	4.8
1	B	514	PRO	4.7
1	B	515	ASP	4.4
1	A	22	ALA	3.0
1	A	24	ALA	3.0
1	A	118	PRO	2.9
1	A	23	VAL	2.8
1	B	727	ALA	2.5
1	A	280	VAL	2.3

### 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
3	SKF	A	3001	14/14	0.98	0.19	1.72	38,42,42,43	0
2	SAH	B	2002	26/26	0.92	0.21	1.44	33,42,46,47	0
2	SAH	A	2001	26/26	0.90	0.20	1.38	48,57,62,62	0
3	SKF	B	3002	14/14	0.97	0.16	0.84	46,49,50,50	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.