Full wwPDB X-ray Structure Validation Report (i)

Feb 19, 2016 – 08:01 PM GMT

PDB ID	:	4X3Q
Title	:	Crystal structure of S-adenosylmethionine-dependent methyltransferase SibL in complex with SAH
Authors	:	liu, J.S.; Chen, S.C.; Yang, C.S.; Huang, C.H.; Chen, Y.
Deposited on	:	2014-12-01
Resolution	:	2.59 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

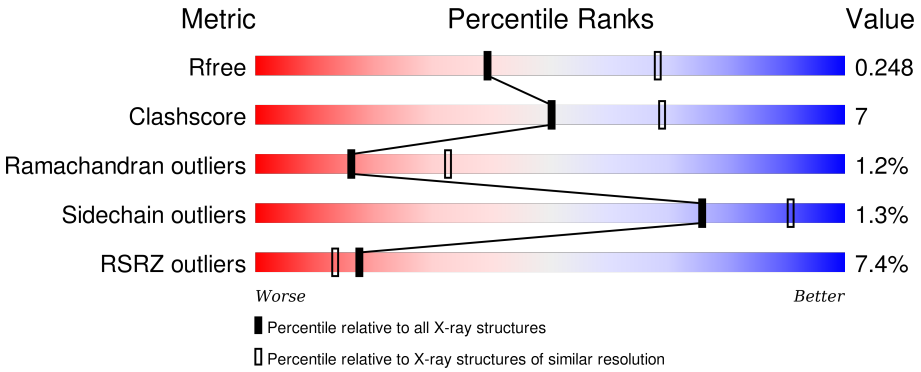
MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026982
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20026982

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 2.59 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	2636 (2.60-2.56)
Clashscore	102246	3003 (2.60-2.56)
Ramachandran outliers	100387	2956 (2.60-2.56)
Sidechain outliers	100360	2956 (2.60-2.56)
RSRZ outliers	91569	2642 (2.60-2.56)

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 <=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	355	6% 82% 13%
1	B	355	6% 80% 15%
1	C	355	7% 81% 14%
1	D	355	10% 80% 15%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10783 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 SibL.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	S	Se	0	0	0
			2639	1673	480	482	3	1			
1	B	341	Total	C	N	O	S	Se	0	0	0
			2639	1673	480	482	3	1			
1	C	341	Total	C	N	O	S	Se	0	0	0
			2639	1673	480	482	3	1			
1	D	341	Total	C	N	O	S	Se	0	0	0
			2639	1673	480	482	3	1			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	345	ALA	-	expression tag	UNP C0LTM6
A	346	ALA	-	expression tag	UNP C0LTM6
A	347	ALA	-	expression tag	UNP C0LTM6
A	348	LEU	-	expression tag	UNP C0LTM6
A	349	GLU	-	expression tag	UNP C0LTM6
A	350	HIS	-	expression tag	UNP C0LTM6
A	351	HIS	-	expression tag	UNP C0LTM6
A	352	HIS	-	expression tag	UNP C0LTM6
A	353	HIS	-	expression tag	UNP C0LTM6
A	354	HIS	-	expression tag	UNP C0LTM6
A	355	HIS	-	expression tag	UNP C0LTM6
B	345	ALA	-	expression tag	UNP C0LTM6
B	346	ALA	-	expression tag	UNP C0LTM6
B	347	ALA	-	expression tag	UNP C0LTM6
B	348	LEU	-	expression tag	UNP C0LTM6
B	349	GLU	-	expression tag	UNP C0LTM6
B	350	HIS	-	expression tag	UNP C0LTM6
B	351	HIS	-	expression tag	UNP C0LTM6
B	352	HIS	-	expression tag	UNP C0LTM6
B	353	HIS	-	expression tag	UNP C0LTM6
B	354	HIS	-	expression tag	UNP C0LTM6

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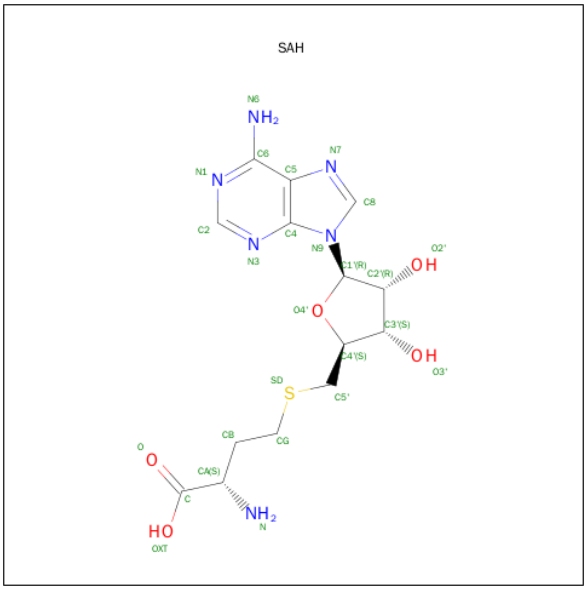
Chain	Residue	Modelled	Actual	Comment	Reference
B	355	HIS	-	expression tag	UNP C0LTM6
C	345	ALA	-	expression tag	UNP C0LTM6
C	346	ALA	-	expression tag	UNP C0LTM6
C	347	ALA	-	expression tag	UNP C0LTM6
C	348	LEU	-	expression tag	UNP C0LTM6
C	349	GLU	-	expression tag	UNP C0LTM6
C	350	HIS	-	expression tag	UNP C0LTM6
C	351	HIS	-	expression tag	UNP C0LTM6
C	352	HIS	-	expression tag	UNP C0LTM6
C	353	HIS	-	expression tag	UNP C0LTM6
C	354	HIS	-	expression tag	UNP C0LTM6
C	355	HIS	-	expression tag	UNP C0LTM6
D	345	ALA	-	expression tag	UNP C0LTM6
D	346	ALA	-	expression tag	UNP C0LTM6
D	347	ALA	-	expression tag	UNP C0LTM6
D	348	LEU	-	expression tag	UNP C0LTM6
D	349	GLU	-	expression tag	UNP C0LTM6
D	350	HIS	-	expression tag	UNP C0LTM6
D	351	HIS	-	expression tag	UNP C0LTM6
D	352	HIS	-	expression tag	UNP C0LTM6
D	353	HIS	-	expression tag	UNP C0LTM6
D	354	HIS	-	expression tag	UNP C0LTM6
D	355	HIS	-	expression tag	UNP C0LTM6

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

- Molecule 3 is water.

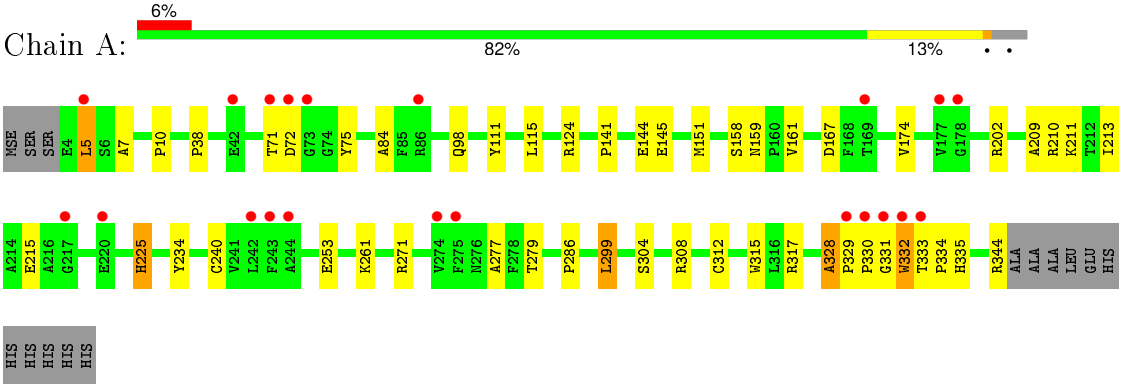
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	41	Total	O	0	0
			41	41		
3	B	29	Total	O	0	0
			29	29		
3	C	28	Total	O	0	0
			28	28		
3	D	25	Total	O	0	0
			25	25		

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.

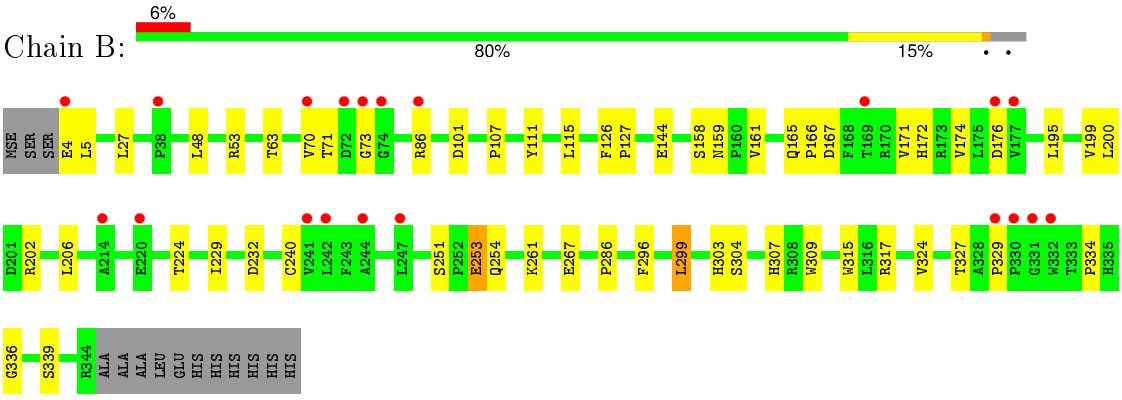
- Molecule 1: SibL

Chain A:



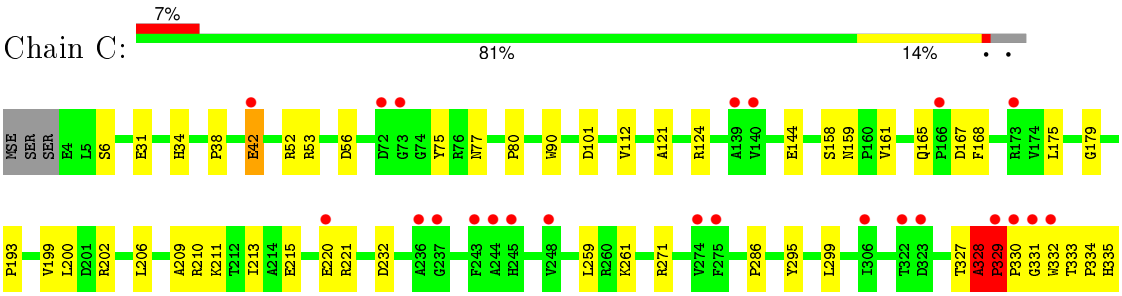
- Molecule 1: SibL

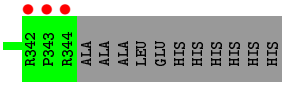
Chain B:



- Molecule 1: SibL

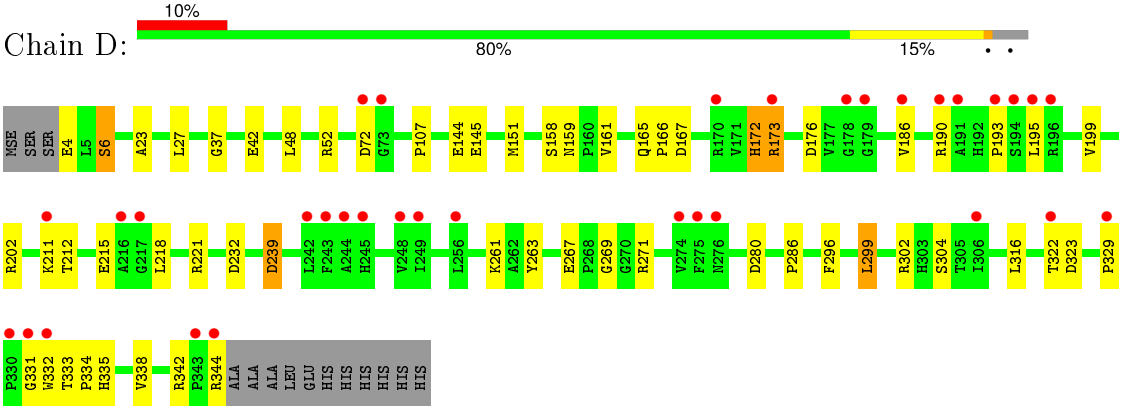
Chain C:





- Molecule 1: SibL

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	91.79Å 112.11Å 143.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.15 – 2.59 28.15 – 2.59	Depositor EDS
% Data completeness (in resolution range)	98.8 (28.15-2.59) 95.4 (28.15-2.59)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.74 (at 2.57Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.184 , 0.248 0.185 , 0.248	Depositor DCC
R_{free} test set	3788 reflections (4.29%)	DCC
Wilson B-factor (Å^2)	48.1	Xtriage
Anisotropy	0.280	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 51.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 46306 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	10783	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 3.44% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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

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	0.49	0/2706	0.69	1/3693 (0.0%)
1	B	0.46	0/2706	0.65	0/3693
1	C	0.46	0/2706	0.65	1/3693 (0.0%)
1	D	0.46	0/2706	0.68	0/3693
All	All	0.47	0/10824	0.67	2/14772 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	C	0	2
All	All	0	3

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	329	PRO	C-N-CD	-7.47	104.16	120.60
1	A	5	LEU	CA-CB-CG	6.18	129.52	115.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	332	TRP	Peptide
1	C	328	ALA	Peptide

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Mol	Chain	Res	Type	Group
1	C	329	PRO	Peptide

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	2639	0	2587	35	1
1	B	2639	0	2587	36	0
1	C	2639	0	2587	40	0
1	D	2639	0	2587	49	1
2	A	26	0	19	2	0
2	B	26	0	19	1	0
2	C	26	0	19	1	0
2	D	26	0	19	2	0
3	A	41	0	0	3	0
3	B	29	0	0	2	0
3	C	28	0	0	3	0
3	D	25	0	0	3	0
All	All	10783	0	10424	148	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 7.

All (148) 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:D:271:ARG:NH1	1:D:323:ASP:OD2	1.95	0.98
1:C:101:ASP:OD1	3:C:509:HOH:O	1.95	0.84
1:D:144:GLU:OE2	1:D:202:ARG:NH2	2.09	0.84
1:A:167:ASP:OD1	1:A:271:ARG:NH1	2.17	0.78
1:A:144:GLU:OE2	1:A:202:ARG:NH2	2.14	0.78
1:C:286:PRO:HG2	1:C:334:PRO:HB3	1.65	0.76
1:A:253:GLU:N	1:A:253:GLU:OE1	2.19	0.75
1:D:48:LEU:O	3:D:509:HOH:O	2.04	0.74
1:A:329:PRO:O	3:A:522:HOH:O	2.05	0.74
1:D:173:ARG:CG	1:D:239:ASP:HB2	2.19	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:144:GLU:OE2	1:C:202:ARG:NH2	2.23	0.71
1:B:71:THR:O	1:B:73:GLY:N	2.22	0.71
1:A:210:ARG:NH2	3:A:535:HOH:O	2.22	0.71
1:A:328:ALA:HB3	1:A:329:PRO:HD3	1.73	0.71
1:D:173:ARG:HG2	1:D:239:ASP:HB2	1.73	0.69
1:C:42:GLU:CD	1:C:52:ARG:HE	1.95	0.69
1:C:330:PRO:O	3:C:524:HOH:O	2.10	0.69
1:B:286:PRO:HG2	1:B:334:PRO:HB3	1.74	0.68
1:D:173:ARG:NH2	1:D:267:GLU:OE2	2.26	0.68
1:C:211:LYS:O	1:C:215:GLU:HG3	1.93	0.68
1:D:286:PRO:HG2	1:D:334:PRO:HB3	1.77	0.67
1:C:167:ASP:OD1	1:C:271:ARG:HD3	1.95	0.67
1:D:190:ARG:NH2	3:D:525:HOH:O	2.28	0.66
1:D:42:GLU:OE1	1:D:52:ARG:NH1	2.23	0.66
1:C:202:ARG:NH1	2:C:401:SAH:O2'	2.28	0.66
1:B:267:GLU:OE1	3:B:501:HOH:O	2.13	0.65
1:A:211:LYS:O	1:A:215:GLU:HG3	1.96	0.65
1:D:322:THR:O	1:D:323:ASP:HB3	1.95	0.65
1:D:167:ASP:OD2	1:D:271:ARG:NH2	2.31	0.63
1:C:333:THR:HG1	1:C:335:HIS:CE1	2.15	0.63
1:B:161:VAL:HG23	1:B:329:PRO:HG2	1.81	0.63
1:D:193:PRO:HA	1:D:221:ARG:HH12	1.63	0.63
1:D:151:MSE:HE2	2:D:401:SAH:HG1	1.80	0.62
1:C:329:PRO:HB2	1:C:330:PRO:HA	1.81	0.61
1:B:253:GLU:HG2	1:B:254:GLN:N	2.17	0.60
1:D:195:LEU:O	1:D:221:ARG:NH2	2.30	0.59
1:D:186:VAL:HG21	1:D:212:THR:HG22	1.86	0.58
1:D:172:HIS:ND1	1:D:172:HIS:N	2.53	0.57
1:C:179:GLY:HA3	1:C:199:VAL:HG13	1.87	0.56
1:B:144:GLU:OE2	1:B:202:ARG:NH2	2.22	0.56
1:A:174:VAL:HG22	1:A:240:CYS:HB3	1.87	0.56
1:C:329:PRO:HB2	1:C:330:PRO:CA	2.35	0.56
1:C:158:SER:O	1:C:161:VAL:HG12	2.06	0.56
1:D:173:ARG:HG3	1:D:239:ASP:HB2	1.86	0.55
1:C:80:PRO:HA	1:D:4:GLU:N	2.22	0.55
1:D:211:LYS:O	1:D:215:GLU:HG3	2.07	0.55
1:A:124:ARG:HD3	1:C:124:ARG:NH1	2.22	0.55
1:B:174:VAL:HG22	1:B:240:CYS:HB3	1.89	0.55
1:B:171:VAL:C	1:B:172:HIS:ND1	2.61	0.54
1:A:333:THR:CG2	1:A:334:PRO:HD2	2.38	0.53
1:A:333:THR:HG21	1:A:335:HIS:CE1	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:158:SER:O	1:D:161:VAL:HG12	2.09	0.52
1:A:141:PRO:CG	1:D:145:GLU:HB3	2.39	0.52
1:D:173:ARG:HB2	1:D:239:ASP:H	1.74	0.52
1:D:202:ARG:HD3	2:D:401:SAH:O2'	2.09	0.52
1:B:202:ARG:HD3	2:B:401:SAH:O2'	2.10	0.52
1:A:124:ARG:HD3	1:C:124:ARG:CZ	2.41	0.51
1:C:328:ALA:HB1	1:C:329:PRO:HA	1.93	0.51
1:C:332:TRP:HZ3	1:D:6:SER:HG	1.56	0.50
1:C:193:PRO:HA	1:C:221:ARG:NH2	2.26	0.50
1:D:280:ASP:OD1	3:D:520:HOH:O	2.19	0.50
1:D:269:GLY:HA3	1:D:344:ARG:HE	1.76	0.50
1:B:327:THR:HG22	1:B:329:PRO:HD3	1.94	0.50
1:D:107:PRO:HG2	1:D:299:LEU:HD11	1.94	0.49
1:B:71:THR:C	1:B:73:GLY:H	2.16	0.49
1:C:121:ALA:HB1	1:C:124:ARG:HH21	1.76	0.49
1:C:193:PRO:HA	1:C:221:ARG:HH22	1.77	0.49
1:A:71:THR:O	1:A:72:ASP:HB3	2.13	0.48
1:A:151:MSE:HE2	2:A:401:SAH:HG1	1.96	0.48
1:B:174:VAL:HG23	1:B:195:LEU:HD21	1.95	0.48
1:B:176:ASP:HB3	1:B:199:VAL:HA	1.96	0.48
1:B:317:ARG:NH2	1:B:324:VAL:O	2.47	0.47
1:A:304:SER:O	1:B:53:ARG:NH2	2.40	0.47
1:C:52:ARG:NH1	1:C:56:ASP:OD2	2.47	0.47
1:C:295:TYR:HA	3:C:514:HOH:O	2.14	0.47
1:C:333:THR:OG1	1:C:334:PRO:HD2	2.15	0.47
1:A:225:HIS:ND1	1:A:225:HIS:C	2.68	0.47
1:B:309:TRP:CH2	1:B:336:GLY:HA3	2.50	0.47
1:A:202:ARG:HD3	2:A:401:SAH:O2'	2.15	0.47
1:B:317:ARG:CZ	1:B:324:VAL:HB	2.44	0.47
1:D:176:ASP:HB3	1:D:199:VAL:HA	1.97	0.47
1:D:173:ARG:H	1:D:173:ARG:HG2	1.51	0.46
1:B:158:SER:O	1:B:161:VAL:HG12	2.16	0.46
1:D:72:ASP:OD1	1:D:72:ASP:O	2.33	0.46
1:B:101:ASP:OD1	3:B:525:HOH:O	2.21	0.46
1:A:333:THR:HG23	1:A:334:PRO:HD2	1.96	0.46
1:A:279:THR:O	1:A:308:ARG:HA	2.15	0.46
1:A:317:ARG:NE	3:A:501:HOH:O	1.88	0.46
1:D:167:ASP:OD1	1:D:167:ASP:N	2.49	0.46
1:C:175:LEU:HD11	1:C:200:LEU:HD11	1.97	0.46
1:C:232:ASP:O	1:C:261:LYS:NZ	2.25	0.46
1:C:220:GLU:OE1	1:C:220:GLU:HA	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:277:ALA:HB1	1:A:334:PRO:HG2	1.98	0.45
1:D:329:PRO:HB3	1:D:333:THR:O	2.17	0.45
1:C:206:LEU:O	1:C:210:ARG:HG3	2.16	0.45
1:D:232:ASP:O	1:D:261:LYS:NZ	2.34	0.45
1:B:307:HIS:CG	1:B:315:TRP:HH2	2.35	0.45
1:D:263:TYR:CZ	1:D:342:ARG:HG3	2.53	0.44
1:A:145:GLU:N	1:A:145:GLU:OE1	2.43	0.44
1:B:286:PRO:CG	1:B:334:PRO:HB3	2.45	0.44
1:A:141:PRO:HG2	1:D:145:GLU:HB3	2.00	0.44
1:C:31:GLU:OE2	1:C:90:TRP:NE1	2.47	0.44
1:B:309:TRP:HH2	1:B:336:GLY:HA3	1.83	0.44
1:C:6:SER:HB3	1:D:332:TRP:CB	2.48	0.44
1:A:286:PRO:CG	1:A:334:PRO:HB3	2.47	0.43
1:D:333:THR:HB	1:D:335:HIS:ND1	2.34	0.43
1:A:158:SER:O	1:A:161:VAL:HG12	2.18	0.43
1:B:111:TYR:CZ	1:B:115:LEU:HD11	2.54	0.43
1:A:312:CYS:HA	1:A:315:TRP:CE3	2.54	0.43
1:A:111:TYR:CZ	1:A:115:LEU:HD11	2.53	0.43
1:D:186:VAL:HG13	1:D:218:LEU:HD12	2.01	0.43
1:B:206:LEU:HD22	1:B:224:THR:HB	2.01	0.43
1:A:299:LEU:HA	1:A:299:LEU:HD23	1.77	0.43
1:D:27:LEU:O	1:D:48:LEU:HD22	2.18	0.43
1:D:296:PHE:O	1:D:304:SER:HB2	2.18	0.43
1:B:126:PHE:HA	1:B:127:PRO:HD3	1.91	0.43
1:C:53:ARG:NH1	1:D:304:SER:O	2.52	0.43
1:A:84:ALA:HB2	1:B:5:LEU:HD11	2.01	0.42
1:D:173:ARG:HG3	1:D:239:ASP:H	1.83	0.42
1:A:38:PRO:HA	1:A:75:TYR:O	2.19	0.42
1:C:38:PRO:HA	1:C:75:TYR:O	2.19	0.42
1:B:296:PHE:O	1:B:304:SER:HB2	2.19	0.42
1:C:333:THR:HG23	1:C:335:HIS:HB2	2.01	0.42
1:C:209:ALA:O	1:C:213:ILE:HG13	2.20	0.42
1:B:251:SER:HB3	1:B:303:HIS:CG	2.55	0.42
1:B:107:PRO:HG2	1:B:299:LEU:HD11	2.02	0.42
1:B:27:LEU:O	1:B:48:LEU:HD22	2.20	0.42
1:A:333:THR:HG22	1:A:334:PRO:HD2	2.02	0.41
1:B:200:LEU:HD22	1:B:229:ILE:HG22	2.02	0.41
1:D:316:LEU:HD22	1:D:338:VAL:HG12	2.02	0.41
1:B:165:GLN:HA	1:B:166:PRO:HD3	1.84	0.41
1:D:202:ARG:HB2	1:D:202:ARG:CZ	2.50	0.41
1:A:209:ALA:O	1:A:213:ILE:HG13	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:307:HIS:ND1	1:B:315:TRP:HH2	2.18	0.41
1:C:112:VAL:HG22	1:D:23:ALA:HA	2.02	0.41
1:A:7:ALA:O	1:A:10:PRO:HD2	2.21	0.41
1:B:63:THR:HG21	1:B:70:VAL:HG12	2.01	0.41
1:C:328:ALA:HB1	1:C:329:PRO:CA	2.51	0.41
1:D:165:GLN:HA	1:D:166:PRO:HD3	1.91	0.41
1:C:259:LEU:HA	1:C:259:LEU:HD23	1.89	0.41
1:D:167:ASP:OD2	1:D:271:ARG:NE	2.51	0.41
1:B:324:VAL:HA	1:B:339:SER:O	2.20	0.41
1:C:6:SER:HB3	1:D:332:TRP:HB3	2.03	0.40
1:D:158:SER:HB2	1:D:335:HIS:CG	2.57	0.40
1:C:165:GLN:HB3	1:C:168:PHE:CD2	2.57	0.40
1:C:34:HIS:HB2	1:C:77:ASN:ND2	2.37	0.40
1:B:232:ASP:O	1:B:261:LYS:NZ	2.35	0.40
1:A:234:TYR:CE1	1:A:261:LYS:HE2	2.56	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:344:ARG:NH2	1:D:37:GLY:O[4_455]	2.12	0.08

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	339/355 (96%)	327 (96%)	6 (2%)	6 (2%)	11	20
1	B	339/355 (96%)	333 (98%)	4 (1%)	2 (1%)	30	54
1	C	339/355 (96%)	329 (97%)	5 (2%)	5 (2%)	13	25
1	D	339/355 (96%)	331 (98%)	5 (2%)	3 (1%)	21	41

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1356/1420 (96%)	1320 (97%)	20 (2%)	16 (1%)	16 33

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	328	ALA
1	A	330	PRO
1	C	328	ALA
1	C	331	GLY
1	C	329	PRO
1	A	299	LEU
1	A	331	GLY
1	B	299	LEU
1	C	299	LEU
1	D	299	LEU
1	A	332	TRP
1	B	159	ASN
1	A	159	ASN
1	D	159	ASN
1	C	159	ASN
1	D	331	GLY

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	270/279 (97%)	267 (99%)	3 (1%)	80 93
1	B	270/279 (97%)	266 (98%)	4 (2%)	72 89
1	C	270/279 (97%)	268 (99%)	2 (1%)	88 96
1	D	270/279 (97%)	265 (98%)	5 (2%)	65 85
All	All	1080/1116 (97%)	1066 (99%)	14 (1%)	76 91

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LEU
1	A	98	GLN
1	A	225	HIS
1	B	4	GLU
1	B	86	ARG
1	B	167	ASP
1	B	253	GLU
1	C	42	GLU
1	C	327	THR
1	D	6	SER
1	D	172	HIS
1	D	173	ARG
1	D	239	ASP
1	D	302	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	GLN

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

There are no carbohydrates in this entry.

5.6 Ligand geometry

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	401	-	22,28,28	0.81	0	18,40,40	1.16	2 (11%)
2	SAH	B	401	-	22,28,28	0.87	0	18,40,40	0.98	1 (5%)
2	SAH	C	401	-	22,28,28	0.86	0	18,40,40	1.04	1 (5%)
2	SAH	D	401	-	22,28,28	0.86	0	18,40,40	1.16	2 (11%)

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
2	SAH	C	401	-	-	0/7/31/31	0/3/3/3
2	SAH	D	401	-	-	0/7/31/31	0/3/3/3

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	401	SAH	C4'-O4'-C1'	-3.05	106.41	109.64
2	D	401	SAH	C4'-O4'-C1'	-2.75	106.73	109.64
2	B	401	SAH	C4'-O4'-C1'	-2.61	106.88	109.64
2	A	401	SAH	C4'-O4'-C1'	-2.42	107.08	109.64
2	A	401	SAH	O4'-C1'-N9	2.39	112.62	108.11
2	D	401	SAH	O4'-C1'-N9	2.57	112.97	108.11

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	SAH	2	0
2	B	401	SAH	1	0
2	C	401	SAH	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	401	SAH	2	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, 95^{th} 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	340/355 (95%)	0.24	21 (6%) 24 20	34, 48, 71, 100	0
1	B	340/355 (95%)	0.24	20 (5%) 26 21	34, 49, 72, 88	0
1	C	340/355 (95%)	0.36	26 (7%) 17 13	33, 51, 74, 98	0
1	D	340/355 (95%)	0.44	34 (10%) 9 7	34, 52, 84, 94	0
All	All	1360/1420 (95%)	0.32	101 (7%) 17 13	33, 50, 77, 100	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	332	TRP	8.4
1	A	332	TRP	7.7
1	A	331	GLY	6.5
1	B	330	PRO	6.4
1	D	332	TRP	5.8
1	C	73	GLY	5.6
1	D	73	GLY	5.5
1	B	73	GLY	5.4
1	B	331	GLY	4.8
1	C	236	ALA	4.5
1	D	248	VAL	4.4
1	A	330	PRO	4.3
1	C	344	ARG	4.2
1	B	72	ASP	4.1
1	D	193	PRO	4.1
1	D	244	ALA	4.1
1	C	72	ASP	4.1
1	D	331	GLY	4.1
1	A	73	GLY	4.1
1	D	217	GLY	3.9
1	C	330	PRO	3.9

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Mol	Chain	Res	Type	RSRZ
1	D	170	ARG	3.8
1	A	329	PRO	3.8
1	B	242	LEU	3.7
1	D	330	PRO	3.7
1	D	72	ASP	3.7
1	D	211	LYS	3.7
1	B	244	ALA	3.5
1	C	140	VAL	3.4
1	C	244	ALA	3.4
1	B	332	TRP	3.3
1	D	343	PRO	3.3
1	A	242	LEU	3.2
1	D	195	LEU	3.2
1	D	344	ARG	3.2
1	A	333	THR	3.2
1	A	42	GLU	3.1
1	A	244	ALA	3.1
1	A	243	PHE	3.0
1	D	216	ALA	3.0
1	D	329	PRO	2.9
1	B	329	PRO	2.9
1	A	71	THR	2.9
1	B	247	LEU	2.9
1	B	4	GLU	2.8
1	C	275	PHE	2.8
1	C	331	GLY	2.8
1	A	86	ARG	2.7
1	C	237	GLY	2.7
1	A	178	GLY	2.7
1	B	220	GLU	2.7
1	A	177	VAL	2.7
1	C	343	PRO	2.6
1	C	243	PHE	2.6
1	D	275	PHE	2.6
1	D	306	ILE	2.6
1	C	274	VAL	2.6
1	C	166	PRO	2.6
1	C	248	VAL	2.5
1	D	179	GLY	2.5
1	D	243	PHE	2.5
1	D	245	HIS	2.5
1	B	177	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	190	ARG	2.5
1	A	217	GLY	2.5
1	A	72	ASP	2.4
1	D	242	LEU	2.4
1	B	70	VAL	2.4
1	B	241	VAL	2.4
1	D	196	ARG	2.4
1	C	306	ILE	2.4
1	A	169	THR	2.4
1	C	173	ARG	2.4
1	C	329	PRO	2.4
1	D	276	ASN	2.3
1	D	186	VAL	2.3
1	C	220	GLU	2.3
1	B	176	ASP	2.3
1	C	245	HIS	2.3
1	B	74	GLY	2.3
1	D	178	GLY	2.3
1	A	220	GLU	2.2
1	D	194	SER	2.2
1	A	5	LEU	2.2
1	D	256	LEU	2.2
1	B	38	PRO	2.2
1	B	86	ARG	2.2
1	D	274	VAL	2.2
1	D	322	THR	2.1
1	C	139	ALA	2.1
1	A	274	VAL	2.1
1	C	342	ARG	2.1
1	D	173	ARG	2.1
1	A	275	PHE	2.1
1	D	191	ALA	2.1
1	C	322	THR	2.1
1	C	323	ASP	2.0
1	B	214	ALA	2.0
1	B	169	THR	2.0
1	D	249	ILE	2.0
1	C	42	GLU	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, 95^{th} 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
2	SAH	A	401	26/26	0.93	0.23	0.61	33,41,47,49	0
2	SAH	C	401	26/26	0.89	0.18	-0.16	43,48,58,61	0
2	SAH	B	401	26/26	0.95	0.17	-0.19	39,44,47,51	0
2	SAH	D	401	26/26	0.92	0.19	-0.21	46,54,60,69	0

6.5 Other polymers (i)

There are no such residues in this entry.