



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 02:32 am GMT

PDB ID : 4K36
Title : His6 tagged anSMEcpe with bound AdoMet
Authors : Goldman, P.J.; Drennan, C.L.
Deposited on : 2013-04-10
Resolution : 1.62 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

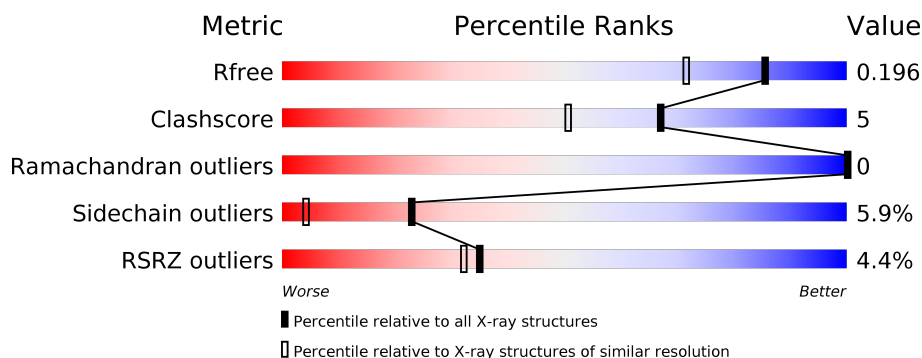
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.62 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	3539 (1.64-1.60)
Clashscore	112137	3855 (1.64-1.60)
Ramachandran outliers	110173	3764 (1.64-1.60)
Sidechain outliers	110143	3763 (1.64-1.60)
RSRZ outliers	101464	3562 (1.64-1.60)

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	392	<div> <div>7%</div> <div> <div></div> <div>83%</div> <div>12%</div> <div>• •</div> </div> </div>
1	B	392	<div> <div>2%</div> <div> <div></div> <div>83%</div> <div>9%</div> <div>• 7%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6793 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 Anaerobic sulfatase-maturing enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	381	Total	C	N	O	S	0	6	0
			3124	2008	523	568	25			
1	B	364	Total	C	N	O	S	0	2	0
			2972	1915	494	537	26			

There are 44 discrepancies between the modelled and reference sequences:

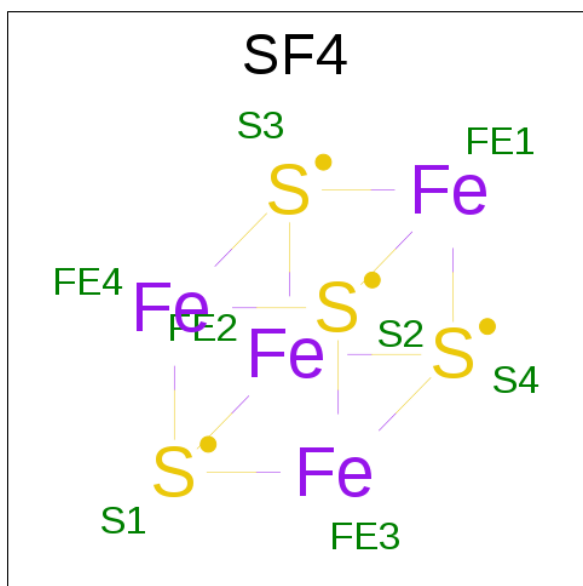
Chain	Residue	Modelled	Actual	Comment	Reference
A	371	SER	-	EXPRESSION TAG	UNP Q0TTH1
A	372	ASP	-	EXPRESSION TAG	UNP Q0TTH1
A	373	PRO	-	EXPRESSION TAG	UNP Q0TTH1
A	374	ASN	-	EXPRESSION TAG	UNP Q0TTH1
A	375	SER	-	EXPRESSION TAG	UNP Q0TTH1
A	376	SER	-	EXPRESSION TAG	UNP Q0TTH1
A	377	SER	-	EXPRESSION TAG	UNP Q0TTH1
A	378	VAL	-	EXPRESSION TAG	UNP Q0TTH1
A	379	ASP	-	EXPRESSION TAG	UNP Q0TTH1
A	380	LYS	-	EXPRESSION TAG	UNP Q0TTH1
A	381	LEU	-	EXPRESSION TAG	UNP Q0TTH1
A	382	ALA	-	EXPRESSION TAG	UNP Q0TTH1
A	383	ALA	-	EXPRESSION TAG	UNP Q0TTH1
A	384	ALA	-	EXPRESSION TAG	UNP Q0TTH1
A	385	LEU	-	EXPRESSION TAG	UNP Q0TTH1
A	386	GLU	-	EXPRESSION TAG	UNP Q0TTH1
A	387	HIS	-	EXPRESSION TAG	UNP Q0TTH1
A	388	HIS	-	EXPRESSION TAG	UNP Q0TTH1
A	389	HIS	-	EXPRESSION TAG	UNP Q0TTH1
A	390	HIS	-	EXPRESSION TAG	UNP Q0TTH1
A	391	HIS	-	EXPRESSION TAG	UNP Q0TTH1
A	392	HIS	-	EXPRESSION TAG	UNP Q0TTH1
B	371	SER	-	EXPRESSION TAG	UNP Q0TTH1
B	372	ASP	-	EXPRESSION TAG	UNP Q0TTH1
B	373	PRO	-	EXPRESSION TAG	UNP Q0TTH1

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	374	ASN	-	EXPRESSION TAG	UNP Q0TTH1
B	375	SER	-	EXPRESSION TAG	UNP Q0TTH1
B	376	SER	-	EXPRESSION TAG	UNP Q0TTH1
B	377	SER	-	EXPRESSION TAG	UNP Q0TTH1
B	378	VAL	-	EXPRESSION TAG	UNP Q0TTH1
B	379	ASP	-	EXPRESSION TAG	UNP Q0TTH1
B	380	LYS	-	EXPRESSION TAG	UNP Q0TTH1
B	381	LEU	-	EXPRESSION TAG	UNP Q0TTH1
B	382	ALA	-	EXPRESSION TAG	UNP Q0TTH1
B	383	ALA	-	EXPRESSION TAG	UNP Q0TTH1
B	384	ALA	-	EXPRESSION TAG	UNP Q0TTH1
B	385	LEU	-	EXPRESSION TAG	UNP Q0TTH1
B	386	GLU	-	EXPRESSION TAG	UNP Q0TTH1
B	387	HIS	-	EXPRESSION TAG	UNP Q0TTH1
B	388	HIS	-	EXPRESSION TAG	UNP Q0TTH1
B	389	HIS	-	EXPRESSION TAG	UNP Q0TTH1
B	390	HIS	-	EXPRESSION TAG	UNP Q0TTH1
B	391	HIS	-	EXPRESSION TAG	UNP Q0TTH1
B	392	HIS	-	EXPRESSION TAG	UNP Q0TTH1

- Molecule 2 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe₄S₄).



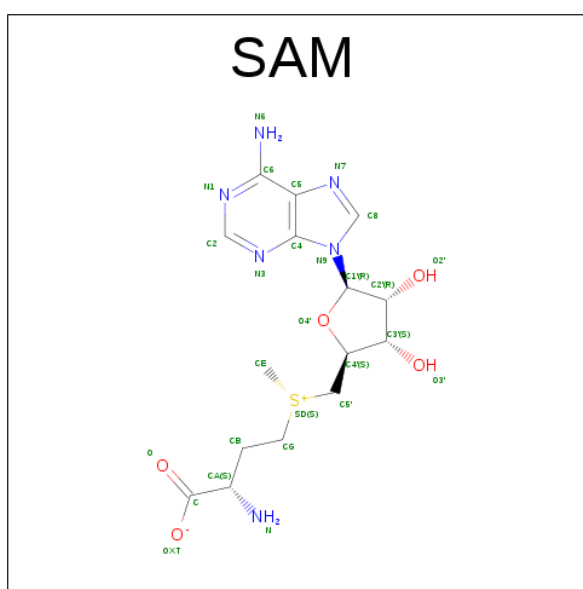
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			8	4	4		
2	A	1	Total	Fe	S	0	0
			8	4	4		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			8	4	4		
2	B	1	Total	Fe	S	0	0
			8	4	4		
2	B	1	Total	Fe	S	0	0
			8	4	4		
2	B	1	Total	Fe	S	0	0
			8	4	4		

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



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

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	Cl	0	0
			2	2		
4	A	1	Total	Cl	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	307	Total 307	O 307	0	0
5	B	285	Total 285	O 285	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	46.56Å 92.62Å 92.35Å 90.00° 94.00° 90.00°	Depositor
Resolution (Å)	37.00 – 1.62 37.00 – 1.62	Depositor EDS
% Data completeness (in resolution range)	98.4 (37.00-1.62) 92.4 (37.00-1.62)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.17 (at 1.62Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.182 , 0.195 0.182 , 0.196	Depositor DCC
R_{free} test set	4894 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	23.6	Xtriage
Anisotropy	0.481	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 51.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6793	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.48% 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: SF4, SAM, CL

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	0.55	0/3212	0.69	2/4315 (0.0%)
1	B	0.56	0/3047	0.67	0/4093
All	All	0.55	0/6259	0.68	2/8408 (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	A	50	ARG	NE-CZ-NH2	-5.30	117.65	120.30
1	A	2	PRO	N-CA-C	-5.12	98.80	112.10

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	3124	0	3029	32	0
1	B	2972	0	2897	23	0
2	A	24	0	0	0	0
2	B	24	0	0	0	0
3	A	27	0	22	0	0
3	B	27	0	22	0	0
4	A	1	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	2	0	0	0	0
5	A	307	0	0	10	0
5	B	285	0	0	5	0
All	All	6793	0	5970	55	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 5.

All (55) 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:228:ASP:O	1:B:233:ASN:ND2	2.22	0.72
1:B:233:ASN:ND2	1:B:233:ASN:H	1.93	0.67
1:A:7[B]:LEU:HD21	1:A:64:GLN:HG3	1.78	0.65
1:B:233:ASN:HD22	1:B:233:ASN:H	1.43	0.64
1:A:324:ARG:NH1	5:A:865:HOH:O	2.31	0.63
1:A:74:GLU:HG2	1:A:78:LYS:HE3	1.80	0.61
1:A:74:GLU:HG3	5:A:873:HOH:O	2.00	0.61
1:B:123:MET:HE2	1:B:146:VAL:HG12	1.82	0.61
1:B:202:LYS:HE2	5:B:877:HOH:O	2.00	0.60
1:A:74:GLU:CG	1:A:78:LYS:HE3	2.34	0.58
1:A:7[B]:LEU:CD2	1:A:64:GLN:HG3	2.34	0.57
1:A:193:ASN:ND2	5:A:690:HOH:O	2.38	0.57
1:A:74:GLU:HG2	1:A:78:LYS:CE	2.35	0.56
1:A:390:HIS:HB3	5:A:814:HOH:O	2.04	0.56
1:A:78:LYS:NZ	5:A:841:HOH:O	2.37	0.56
1:A:85:LYS:NZ	5:A:881:HOH:O	2.38	0.55
1:A:123:MET:HE2	1:A:146:VAL:HG12	1.87	0.55
1:B:305:GLU:HG3	5:B:811:HOH:O	2.08	0.54
1:A:106:GLU:HG2	1:A:156:TYR:OH	2.09	0.53
1:A:25:HIS:CD2	1:A:28:SER:HB3	2.44	0.53
1:A:199:TYR:CD2	1:A:338:GLU:HG2	2.45	0.50
1:A:74:GLU:CG	1:A:78:LYS:CE	2.90	0.50
1:B:199:TYR:CD2	1:B:338:GLU:HG2	2.48	0.49
1:B:110:LYS:HG3	1:B:156:TYR:CE2	2.48	0.48
1:B:233:ASN:HD22	1:B:233:ASN:N	2.08	0.47
1:A:390:HIS:CB	5:A:725:HOH:O	2.62	0.47
1:B:176:LYS:NZ	5:B:863:HOH:O	2.33	0.47
1:B:251:LYS:HG3	1:B:258:ASN:OD1	2.14	0.47
1:A:390:HIS:HB3	5:A:725:HOH:O	2.15	0.46
1:A:123:MET:HE3	1:A:162:ILE:HD12	1.98	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:SER:OG	1:A:388:HIS:NE2	2.44	0.46
1:B:33:VAL:N	5:B:876:HOH:O	2.50	0.45
1:A:18:LYS:HE3	1:A:18:LYS:HB2	1.72	0.45
1:B:201:GLU:HG2	5:B:818:HOH:O	2.17	0.45
1:A:173:HIS:HB3	1:A:176:LYS:HD3	1.99	0.44
1:A:175:ASN:OD1	1:A:176:LYS:HD2	2.17	0.43
1:A:81:GLU:HG2	1:A:84:ARG:NH1	2.32	0.43
1:B:216:LYS:HB2	1:B:216:LYS:NZ	2.32	0.43
1:B:305:GLU:O	1:B:309:LEU:HD23	2.17	0.43
1:A:7[B]:LEU:HD11	1:A:64:GLN:OE1	2.18	0.43
1:B:190[B]:GLN:HG2	1:B:238:ARG:HB2	2.01	0.43
1:B:233:ASN:N	1:B:233:ASN:ND2	2.59	0.43
1:A:212[A]:LYS:HE2	5:A:799:HOH:O	2.18	0.43
1:A:38[B]:ILE:HD12	1:A:71:ALA:HA	2.01	0.42
1:A:202:LYS:NZ	5:A:758:HOH:O	2.52	0.42
1:A:130:HIS:CE1	1:A:134:ARG:HD3	2.54	0.41
1:B:38:ILE:HD12	1:B:71:ALA:HA	2.01	0.41
1:B:123:MET:HE2	1:B:146:VAL:CG1	2.49	0.41
1:A:147:GLU:O	1:A:151:GLU:HG3	2.21	0.41
1:B:123:MET:HB2	1:B:123:MET:HE2	1.93	0.41
1:B:315:GLU:CD	1:B:318:LYS:HD2	2.41	0.41
1:B:337:LYS:HG2	1:B:338:GLU:N	2.37	0.40
1:A:313:VAL:O	1:A:318:LYS:HE3	2.22	0.40
1:B:23:PHE:CE1	1:B:24:TYR:CE1	3.10	0.40
1:A:245:GLU:OE2	1:A:385:LEU:HD22	2.22	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	381/392 (97%)	372 (98%)	9 (2%)	0	100 100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	361/392 (92%)	353 (98%)	8 (2%)	0	100	100
All	All	742/784 (95%)	725 (98%)	17 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	341/358 (95%)	316 (93%)	25 (7%)	16	2
1	B	327/358 (91%)	311 (95%)	16 (5%)	29	7
All	All	668/716 (93%)	627 (94%)	41 (6%)	23	4

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	LYS
1	A	44	LEU
1	A	64	GLN
1	A	103	LEU
1	A	106	GLU
1	A	147	GLU
1	A	154	LYS
1	A	163	LEU
1	A	179	LYS
1	A	212[A]	LYS
1	A	212[B]	LYS
1	A	234	ARG
1	A	251	LYS
1	A	257	MET
1	A	286	LEU
1	A	309	LEU
1	A	316[A]	GLU
1	A	316[B]	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	319	LYS
1	A	324	ARG
1	A	325	LEU
1	A	331	ARG
1	A	338	GLU
1	A	340	SER
1	A	368	ASN
1	B	33	VAL
1	B	44	LEU
1	B	73	LEU
1	B	103	LEU
1	B	128	GLU
1	B	163	LEU
1	B	179	LYS
1	B	185	ASP
1	B	233	ASN
1	B	238	ARG
1	B	251	LYS
1	B	282	ASP
1	B	283	LYS
1	B	331	ARG
1	B	342	LEU
1	B	369	ILE

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

Mol	Chain	Res	Type
1	B	233	ASN

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

Of 11 ligands modelled in this entry, 3 are monoatomic - leaving 8 for Mogul analysis.

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	SF4	A	501	1	0,12,12	0.00	-	0,24,24	0.00	-
2	SF4	A	502	1	0,12,12	0.00	-	0,24,24	0.00	-
2	SF4	A	503	1,3	0,12,12	0.00	-	0,24,24	0.00	-
3	SAM	A	504	2	21,29,29	1.01	1 (4%)	17,42,42	2.69	3 (17%)
2	SF4	B	501	1	0,12,12	0.00	-	0,24,24	0.00	-
2	SF4	B	502	1	0,12,12	0.00	-	0,24,24	0.00	-
2	SF4	B	503	1,3	0,12,12	0.00	-	0,24,24	0.00	-
3	SAM	B	504	2	21,29,29	0.93	1 (4%)	17,42,42	1.97	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	SF4	A	501	1	-	0/0/48/48	0/6/5/5
2	SF4	A	502	1	-	0/0/48/48	0/6/5/5
2	SF4	A	503	1,3	-	0/0/48/48	0/6/5/5
3	SAM	A	504	2	-	0/8/33/33	0/3/3/3
2	SF4	B	501	1	-	0/0/48/48	0/6/5/5
2	SF4	B	502	1	-	0/0/48/48	0/6/5/5
2	SF4	B	503	1,3	-	0/0/48/48	0/6/5/5
3	SAM	B	504	2	-	0/8/33/33	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	504	SAM	C2-N3	3.01	1.37	1.32
3	B	504	SAM	C2-N3	3.19	1.37	1.32

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	504	SAM	N3-C2-N1	-9.70	120.41	128.86
3	B	504	SAM	N3-C2-N1	-6.89	122.86	128.86
3	A	504	SAM	C4'-O4'-C1'	-3.00	106.58	109.77
3	A	504	SAM	C4-C5-N7	-2.26	107.23	109.41
3	B	504	SAM	C4'-O4'-C1'	-2.11	107.53	109.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	381/392 (97%)	0.40	26 (6%) 18 17	22, 33, 51, 69	0
1	B	364/392 (92%)	0.18	7 (1%) 67 66	23, 32, 45, 60	0
All	All	745/784 (95%)	0.29	33 (4%) 35 32	22, 33, 50, 69	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	30	ASN	5.5
1	A	379	ASP	5.4
1	A	369	ILE	5.3
1	A	311	PHE	4.9
1	A	381	LEU	4.5
1	A	249	LEU	4.4
1	A	378	VAL	4.2
1	A	29	ASP	3.9
1	A	368	ASN	3.8
1	B	33	VAL	3.7
1	A	380	LYS	3.7
1	A	366	ALA	3.7
1	A	385	LEU	3.4
1	A	363	ILE	3.4
1	B	369	ILE	3.2
1	A	383	ALA	3.1
1	A	33	VAL	3.0
1	A	28	SER	2.7
1	A	234	ARG	2.7
1	A	316[A]	GLU	2.6
1	A	27	LEU	2.6
1	A	192	ILE	2.5
1	A	319	LYS	2.5
1	B	233	ASN	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	365	VAL	2.5
1	A	315	GLU	2.4
1	A	324	ARG	2.4
1	B	387	HIS	2.4
1	B	201	GLU	2.3
1	B	34	LYS	2.3
1	A	392	HIS	2.2
1	B	339	ASP	2.1
1	A	7[A]	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
2	SF4	B	501	8/8	1.00	0.09	0.52	25,26,26,26	0
2	SF4	B	502	8/8	0.99	0.09	-0.46	26,26,27,28	0
3	SAM	A	504	27/27	0.95	0.12	-0.49	23,27,31,31	0
3	SAM	B	504	27/27	0.95	0.11	-0.63	22,29,31,33	7
2	SF4	A	502	8/8	0.99	0.07	-0.88	32,33,34,35	0
2	SF4	B	503	8/8	0.98	0.08	-1.24	24,26,27,28	0
2	SF4	A	503	8/8	0.99	0.09	-1.71	23,24,26,27	0
2	SF4	A	501	8/8	0.99	0.05	-1.72	29,30,30,31	0
4	CL	B	506	1/1	0.97	0.05	-1.99	37,37,37,37	0
4	CL	B	505	1/1	0.98	0.04	-2.75	29,29,29,29	0
4	CL	A	505	1/1	0.99	0.06	-5.16	28,28,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.