W O R L D W I D E

wwPDB

PROTEIN DATA BANK

wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 1, 2017 – 09:58 PM EDT

PDB ID	:	3SUF
Title	:	Crystal structure of NS3/4A protease variant D168A in complex with MK-5172
Authors	:	Schiffer, C.A.; Romano, K.P.
Deposited on	:	unknown
Resolution	:	2.19 Å(reported)

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

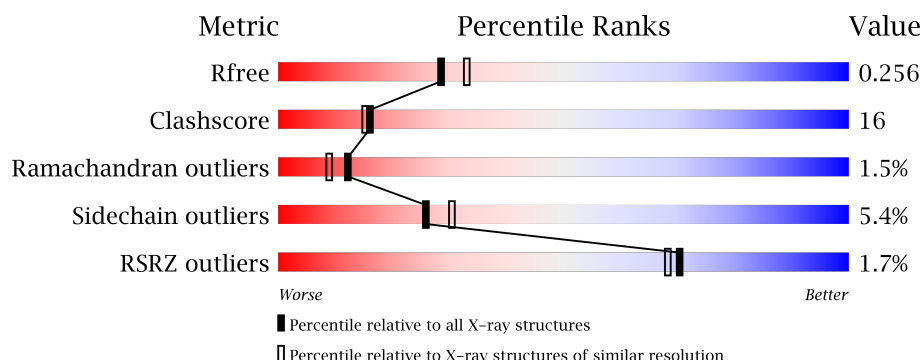|		
MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030345
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030345

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.19 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	Percentile Ranks	Value
Rfree		0.256
Clashscore		16
Ramachandran outliers		1.5%
Sidechain outliers		5.4%
RSRZ outliers		1.7%

Worse   *Better*

∎ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (2.20-2.20)

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	203	2% 67% 25% • 6%
1	B	203	67% 23% • 6%
1	C	203	.% 52% 38% • 6%
1	D	203	2% 69% 21% 10%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SUE	A	1201	X	-	-	-
2	SUE	B	1201	X	-	-	-
2	SUE	C	1201	X	-	-	-
2	SUE	D	1201	X	-	-	-
3	ZN	B	1203	-	-	X	-
3	ZN	C	1202	-	-	X	-

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5672 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 NS3 protease, NS4A protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	190	Total	C	N	O	S	0	0	0
			1335	831	233	264	7			
1	B	190	Total	C	N	O	S	0	0	0
			1341	834	236	264	7			
1	C	190	Total	C	N	O	S	0	0	0
			1341	835	239	260	7			
1	D	183	Total	C	N	O	S	0	0	0
			1286	804	228	248	6			

There are 108 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	980	GLY	-	EXPRESSION TAG	UNP A8DG50
A	981	SER	-	EXPRESSION TAG	UNP A8DG50
A	982	HIS	-	EXPRESSION TAG	UNP A8DG50
A	983	MET	-	EXPRESSION TAG	UNP A8DG50
A	984	ALA	-	EXPRESSION TAG	UNP A8DG50
A	985	SER	-	EXPRESSION TAG	UNP A8DG50
A	986	MET	CYS	ENGINEERED MUTATION	UNP A8DG50
A	987	LYS	LEU	ENGINEERED MUTATION	UNP A8DG50
A	988	LYS	SER	ENGINEERED MUTATION	UNP A8DG50
A	989	LYS	THR	ENGINEERED MUTATION	UNP A8DG50
A	991	SER	CYS	SEE REMARK 999	UNP A8DG50
A	998	ILE	VAL	SEE REMARK 999	UNP A8DG50
A	999	ASN	ILE	SEE REMARK 999	UNP A8DG50
A	1001	SER	ALA	ENGINEERED MUTATION	UNP A8DG50
A	1002	GLY	PRO	ENGINEERED MUTATION	UNP A8DG50
A	1003	ASP	ILE	ENGINEERED MUTATION	UNP A8DG50
A	1013	GLU	LEU	ENGINEERED MUTATION	UNP A8DG50
A	1014	GLU	LEU	ENGINEERED MUTATION	UNP A8DG50
A	1017	GLN	ILE	ENGINEERED MUTATION	UNP A8DG50
A	1018	GLU	ILE	ENGINEERED MUTATION	UNP A8DG50
A	1021	GLN	LEU	ENGINEERED MUTATION	UNP A8DG50

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1040	THR	ALA	ENGINEERED MUTATION	UNP A8DG50
A	1047	SER	CYS	ENGINEERED MUTATION	UNP A8DG50
A	1052	LEU	CYS	ENGINEERED MUTATION	UNP A8DG50
A	1072	THR	ILE	ENGINEERED MUTATION	UNP A8DG50
A	1086	GLN	PRO	ENGINEERED MUTATION	UNP A8DG50
A	1168	ALA	ASP	ENGINEERED MUTATION	UNP A8DG50
B	980	GLY	-	EXPRESSION TAG	UNP A8DG50
B	981	SER	-	EXPRESSION TAG	UNP A8DG50
B	982	HIS	-	EXPRESSION TAG	UNP A8DG50
B	983	MET	-	EXPRESSION TAG	UNP A8DG50
B	984	ALA	-	EXPRESSION TAG	UNP A8DG50
B	985	SER	-	EXPRESSION TAG	UNP A8DG50
B	986	MET	CYS	ENGINEERED MUTATION	UNP A8DG50
B	987	LYS	LEU	ENGINEERED MUTATION	UNP A8DG50
B	988	LYS	SER	ENGINEERED MUTATION	UNP A8DG50
B	989	LYS	THR	ENGINEERED MUTATION	UNP A8DG50
B	991	SER	CYS	SEE REMARK 999	UNP A8DG50
B	998	ILE	VAL	SEE REMARK 999	UNP A8DG50
B	999	ASN	ILE	SEE REMARK 999	UNP A8DG50
B	1001	SER	ALA	ENGINEERED MUTATION	UNP A8DG50
B	1002	GLY	PRO	ENGINEERED MUTATION	UNP A8DG50
B	1003	ASP	ILE	ENGINEERED MUTATION	UNP A8DG50
B	1013	GLU	LEU	ENGINEERED MUTATION	UNP A8DG50
B	1014	GLU	LEU	ENGINEERED MUTATION	UNP A8DG50
B	1017	GLN	ILE	ENGINEERED MUTATION	UNP A8DG50
B	1018	GLU	ILE	ENGINEERED MUTATION	UNP A8DG50
B	1021	GLN	LEU	ENGINEERED MUTATION	UNP A8DG50
B	1040	THR	ALA	ENGINEERED MUTATION	UNP A8DG50
B	1047	SER	CYS	ENGINEERED MUTATION	UNP A8DG50
B	1052	LEU	CYS	ENGINEERED MUTATION	UNP A8DG50
B	1072	THR	ILE	ENGINEERED MUTATION	UNP A8DG50
B	1086	GLN	PRO	ENGINEERED MUTATION	UNP A8DG50
B	1168	ALA	ASP	ENGINEERED MUTATION	UNP A8DG50
C	980	GLY	-	EXPRESSION TAG	UNP A8DG50
C	981	SER	-	EXPRESSION TAG	UNP A8DG50
C	982	HIS	-	EXPRESSION TAG	UNP A8DG50
C	983	MET	-	EXPRESSION TAG	UNP A8DG50
C	984	ALA	-	EXPRESSION TAG	UNP A8DG50
C	985	SER	-	EXPRESSION TAG	UNP A8DG50
C	986	MET	CYS	ENGINEERED MUTATION	UNP A8DG50
C	987	LYS	LEU	ENGINEERED MUTATION	UNP A8DG50
C	988	LYS	SER	ENGINEERED MUTATION	UNP A8DG50

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Chain	Residue	Modelled	Actual	Comment	Reference
C	989	LYS	THR	ENGINEERED MUTATION	UNP A8DG50
C	991	SER	CYS	SEE REMARK 999	UNP A8DG50
C	998	ILE	VAL	SEE REMARK 999	UNP A8DG50
C	999	ASN	ILE	SEE REMARK 999	UNP A8DG50
C	1001	SER	ALA	ENGINEERED MUTATION	UNP A8DG50
C	1002	GLY	PRO	ENGINEERED MUTATION	UNP A8DG50
C	1003	ASP	ILE	ENGINEERED MUTATION	UNP A8DG50
C	1013	GLU	LEU	ENGINEERED MUTATION	UNP A8DG50
C	1014	GLU	LEU	ENGINEERED MUTATION	UNP A8DG50
C	1017	GLN	ILE	ENGINEERED MUTATION	UNP A8DG50
C	1018	GLU	ILE	ENGINEERED MUTATION	UNP A8DG50
C	1021	GLN	LEU	ENGINEERED MUTATION	UNP A8DG50
C	1040	THR	ALA	ENGINEERED MUTATION	UNP A8DG50
C	1047	SER	CYS	ENGINEERED MUTATION	UNP A8DG50
C	1052	LEU	CYS	ENGINEERED MUTATION	UNP A8DG50
C	1072	THR	ILE	ENGINEERED MUTATION	UNP A8DG50
C	1086	GLN	PRO	ENGINEERED MUTATION	UNP A8DG50
C	1168	ALA	ASP	ENGINEERED MUTATION	UNP A8DG50
D	980	GLY	-	EXPRESSION TAG	UNP A8DG50
D	981	SER	-	EXPRESSION TAG	UNP A8DG50
D	982	HIS	-	EXPRESSION TAG	UNP A8DG50
D	983	MET	-	EXPRESSION TAG	UNP A8DG50
D	984	ALA	-	EXPRESSION TAG	UNP A8DG50
D	985	SER	-	EXPRESSION TAG	UNP A8DG50
D	986	MET	CYS	ENGINEERED MUTATION	UNP A8DG50
D	987	LYS	LEU	ENGINEERED MUTATION	UNP A8DG50
D	988	LYS	SER	ENGINEERED MUTATION	UNP A8DG50
D	989	LYS	THR	ENGINEERED MUTATION	UNP A8DG50
D	991	SER	CYS	SEE REMARK 999	UNP A8DG50
D	998	ILE	VAL	SEE REMARK 999	UNP A8DG50
D	999	ASN	ILE	SEE REMARK 999	UNP A8DG50
D	1001	SER	ALA	ENGINEERED MUTATION	UNP A8DG50
D	1002	GLY	PRO	ENGINEERED MUTATION	UNP A8DG50
D	1003	ASP	ILE	ENGINEERED MUTATION	UNP A8DG50
D	1013	GLU	LEU	ENGINEERED MUTATION	UNP A8DG50
D	1014	GLU	LEU	ENGINEERED MUTATION	UNP A8DG50
D	1017	GLN	ILE	ENGINEERED MUTATION	UNP A8DG50
D	1018	GLU	ILE	ENGINEERED MUTATION	UNP A8DG50
D	1021	GLN	LEU	ENGINEERED MUTATION	UNP A8DG50
D	1040	THR	ALA	ENGINEERED MUTATION	UNP A8DG50
D	1047	SER	CYS	ENGINEERED MUTATION	UNP A8DG50
D	1052	LEU	CYS	ENGINEERED MUTATION	UNP A8DG50

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Chain	Residue	Modelled	Actual	Comment	Reference
D	1072	THR	ILE	ENGINEERED MUTATION	UNP A8DG50
D	1086	GLN	PRO	ENGINEERED MUTATION	UNP A8DG50
D	1168	ALA	ASP	ENGINEERED MUTATION	UNP A8DG50

- Molecule 2 is (1aR,5S,8S,10R,22aR)-5-tert-butyl-N-{(1R,2S)-1-[(cyclopropylsulfonyl)carba moyl]-2-ethenylcyclopropyl}-14-methoxy-3,6-dioxo-1,1a,3,4,5,6,9,10,18,19,20,21,22,22a-tetra decahydro-8H-7,10-methanocyclopropa[18,19][1,10,3,6]dioxadiazacyclononadecino[11,12-b]q uinoxaline-8-carboxamide (three-letter code: SUE) (formula: $C_{38}H_{50}N_6O_9S$).
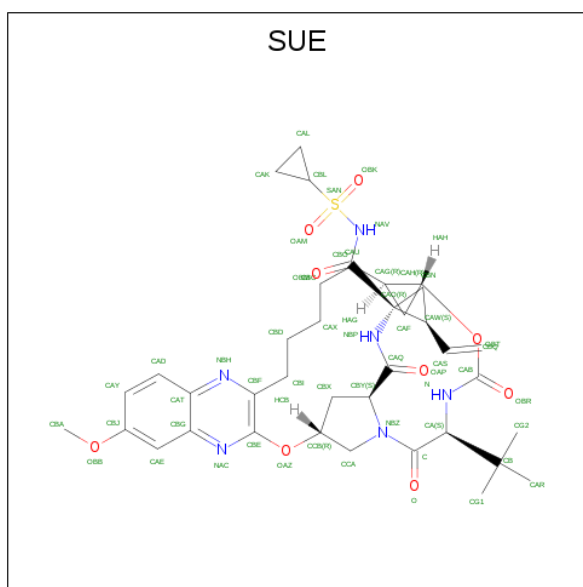
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			54	38	6	9	1		
2	B	1	Total	C	N	O	S	0	0
			54	38	6	9	1		
2	C	1	Total	C	N	O	S	0	0
			54	38	6	9	1		
2	D	1	Total	C	N	O	S	0	0
			54	38	6	9	1		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		
3	A	1	Total	Zn	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total 1	Zn 1	0	0
3	C	1	Total 1	Zn 1	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).
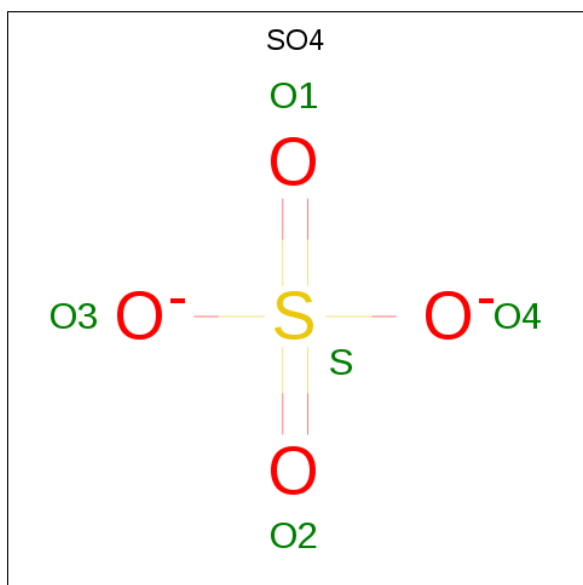
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total 5	O 4	S 1	0	0

- Molecule 5 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	38	Total 38	O 38	0	0
5	B	38	Total 40	O 40	0	2
5	C	35	Total 36	O 36	0	1
5	D	30	Total 30	O 30	0	0

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 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.
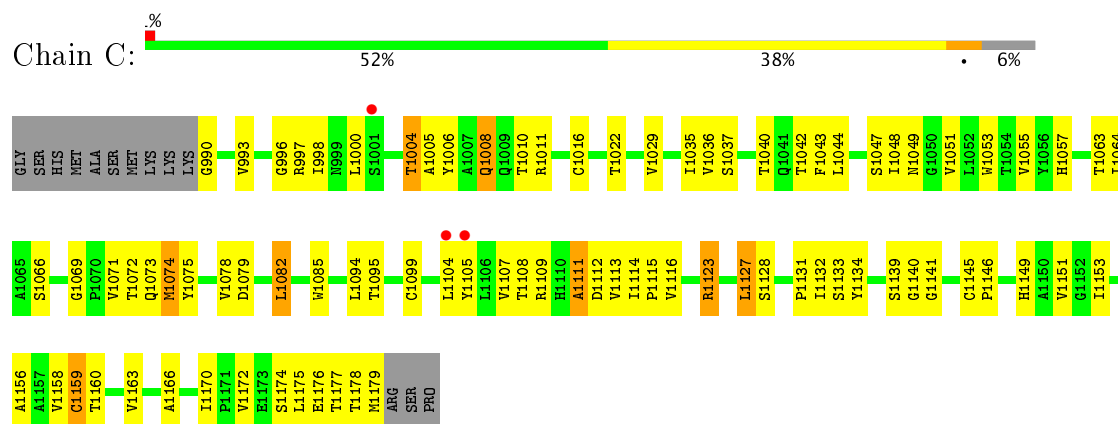
- Molecule 1: NS3 protease, NS4A protein
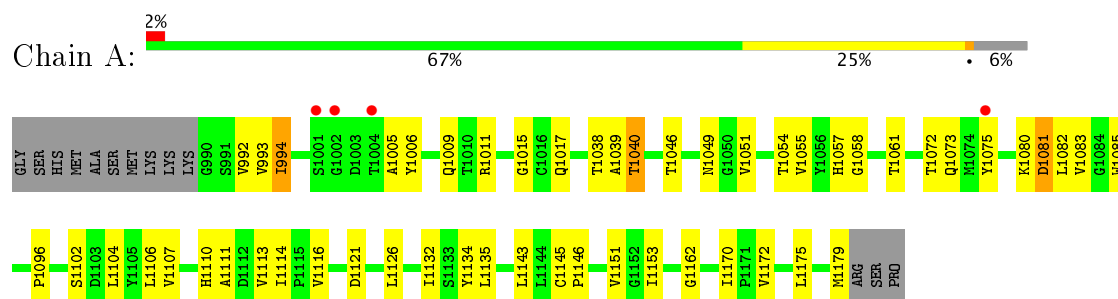
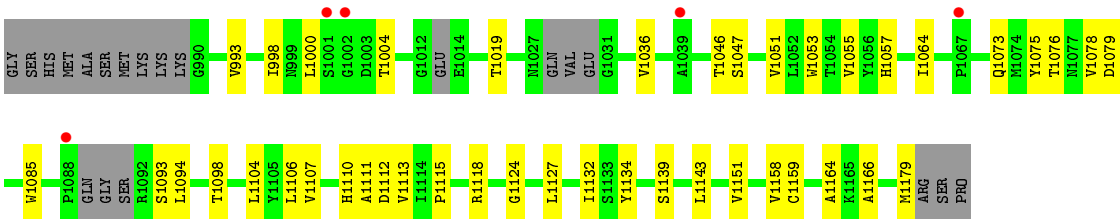- Molecule 1: NS3 protease, NS4A protein
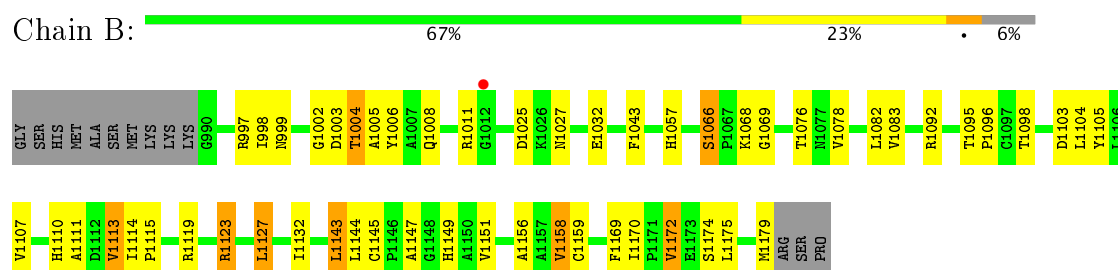
- Molecule 1: NS3 protease, NS4A protein

- Molecule 1: NS3 protease, NS4A protein

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	56.00Å 103.56Å 73.51Å 90.00° 112.04° 90.00°	Depositor
Resolution (Å)	29.22 – 2.19 29.21 – 2.19	Depositor EDS
% Data completeness (in resolution range)	94.3 (29.22-2.19) 94.5 (29.21-2.19)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.39 (at 2.20Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.199 , 0.259 0.199 , 0.256	Depositor DCC
R_{free} test set	1881 reflections (5.26%)	DCC
Wilson B-factor (Å^2)	34.1	Xtriage
Anisotropy	0.281	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 20.2	EDS
L-test for twinning[2]	$< |L| > = 0.39, < L^2 > = 0.22$	Xtriage
Estimated twinning fraction	0.427 for h,-k,-h-l	Xtriage
Reported twinning fraction	0.568 for H, K, L 0.432 for -H, -K, H+L	Depositor
Outliers	0 of 37653 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5672	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.45% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: SUE, ZN, SO4

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	0.54	0/1359	0.67	0/1861
1	B	0.59	1/1365 (0.1%)	0.69	2/1868 (0.1%)
1	C	0.54	1/1365 (0.1%)	0.68	0/1867
1	D	0.50	0/1307	0.62	0/1785
All	All	0.54	2/5396 (0.0%)	0.67	2/7381 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1066	SER	CB-OG	-6.20	1.34	1.42
1	C	1085	TRP	CD2-CE2	5.07	1.47	1.41

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1123	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	B	1066	SER	CB-CA-C	-5.13	100.35	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	1335	0	1285	40	0
1	B	1341	0	1296	44	0
1	C	1341	0	1304	71	0
1	D	1286	0	1249	29	0
2	A	54	0	50	2	0
2	B	54	0	50	5	0
2	C	54	0	50	3	0
2	D	54	0	49	4	0
3	A	1	0	0	0	0
3	B	1	0	0	2	0
3	C	1	0	0	2	0
3	D	1	0	0	0	0
4	B	5	0	0	0	0
5	A	38	0	0	0	0
5	B	40	0	0	3	0
5	C	36	0	0	3	0
5	D	30	0	0	0	0
All	All	5672	0	5333	176	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 16.

The worst 5 of 176 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:C:1095:THR:HG23	1:C:1149:HIS:CD2	1.99	0.98
1:C:1123:ARG:HG3	1:C:1123:ARG:HH11	1.32	0.94
1:C:1095:THR:HG23	1:C:1149:HIS:HD2	1.30	0.94
1:B:1145:CYS:HG	3:B:1203:ZN:ZN	0.73	0.92
1:C:1099:CYS:SG	3:C:1202:ZN:ZN	1.59	0.91

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	188/203 (93%)	170 (90%)	14 (7%)	4 (2%)	8	5
1	B	188/203 (93%)	179 (95%)	6 (3%)	3 (2%)	11	8
1	C	188/203 (93%)	169 (90%)	15 (8%)	4 (2%)	8	5
1	D	175/203 (86%)	161 (92%)	14 (8%)	0	100	100
All	All	739/812 (91%)	679 (92%)	49 (7%)	11 (2%)	12	9

5 of 11 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	1111	ALA
1	A	1121	ASP
1	B	1003	ASP
1	C	1177	THR
1	A	1017	GLN

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	139/164 (85%)	134 (96%)	5 (4%)	40	50
1	B	140/164 (85%)	131 (94%)	9 (6%)	20	23
1	C	139/164 (85%)	128 (92%)	11 (8%)	14	14
1	D	133/164 (81%)	128 (96%)	5 (4%)	38	47
All	All	551/656 (84%)	521 (95%)	30 (5%)	26	30

5 of 30 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	1179	MET
1	C	1047	SER
1	D	1064	ILE

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Mol	Chain	Res	Type
1	C	1040	THR
1	C	1074	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1089	GLN
1	B	1149	HIS
1	C	1073	GLN
1	C	1149	HIS
1	D	999	ASN

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)

Of 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 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	SUE	A	1201	-	57,60,60	2.82	22 (38%)	74,92,92	1.97	14 (18%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	#|Z| > 2	Counts	RMSZ	#|Z| > 2
2	SUE	B	1201	-	57,60,60	2.60	17 (29%)	74,92,92	1.81	19 (25%)
4	SO4	B	1202	-	4,4,4	0.46	0	6,6,6	0.40	0
2	SUE	C	1201	-	57,60,60	2.52	24 (42%)	74,92,92	2.12	21 (28%)
2	SUE	D	1201	-	57,60,60	2.72	21 (36%)	74,92,92	2.05	23 (31%)

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	SUE	A	1201	-	1/1/15/19	0/58/91/91	0/2/7/7
2	SUE	B	1201	-	1/1/15/19	0/58/91/91	0/2/7/7
4	SO4	B	1202	-	-	0/0/0/0	0/0/0/0
2	SUE	C	1201	-	1/1/15/19	0/58/91/91	0/2/7/7
2	SUE	D	1201	-	1/1/15/19	0/58/91/91	0/2/7/7

The worst 5 of 84 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1201	SUE	CA-C	-10.91	1.36	1.53
2	D	1201	SUE	SAN-NAV	-8.59	1.41	1.62
2	B	1201	SUE	CA-C	-8.49	1.40	1.53
2	C	1201	SUE	CA-C	-8.11	1.40	1.53
2	D	1201	SUE	CA-C	-7.92	1.41	1.53

The worst 5 of 77 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	1201	SUE	OBM-CBO-CAO	-5.88	113.77	120.77
2	A	1201	SUE	OBR-CAB-N	-5.67	115.11	124.87
2	C	1201	SUE	CAW-CAO-NBP	-5.31	106.86	117.61
2	D	1201	SUE	CAW-CAO-NBP	-4.84	107.82	117.61
2	A	1201	SUE	CAW-CAO-NBP	-4.53	108.44	117.61

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	1201	SUE	NBZ
2	A	1201	SUE	NBZ
2	C	1201	SUE	NBZ

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Mol	Chain	Res	Type	Atom
2	D	1201	SUE	NBZ

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1201	SUE	2	0
2	B	1201	SUE	5	0
2	C	1201	SUE	3	0
2	D	1201	SUE	4	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (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, 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	190/203 (93%)	0.30	4 (2%)	64	61	19, 32, 50, 60	0
1	B	190/203 (93%)	0.25	1 (0%)	90	90	21, 32, 49, 61	0
1	C	190/203 (93%)	0.34	3 (1%)	72	70	22, 34, 47, 54	0
1	D	183/203 (90%)	0.44	5 (2%)	55	52	24, 39, 61, 66	0
All	All	753/812 (92%)	0.33	13 (1%)	70	68	19, 34, 52, 66	0

The worst 5 of 13 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1001	SER	6.1
1	A	1004	THR	4.7
1	D	1001	SER	4.0
1	D	1002	GLY	4.0
1	A	1002	GLY	3.1

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(Å²)	Q<0.9
2	SUE	A	1201	54/54	0.94	0.13	-0.17	25,29,32,32	0
2	SUE	B	1201	54/54	0.94	0.13	-0.45	22,29,33,34	0
3	ZN	A	1202	1/1	0.98	0.12	-0.52	36,36,36,36	0
2	SUE	C	1201	54/54	0.95	0.12	-0.56	20,27,36,39	0
3	ZN	B	1203	1/1	0.98	0.10	-0.69	36,36,36,36	0
4	SO4	B	1202	5/5	0.93	0.11	-0.77	40,43,44,44	0
2	SUE	D	1201	54/54	0.94	0.12	-0.95	22,28,30,31	0
3	ZN	D	1202	1/1	0.99	0.07	-2.94	33,33,33,33	0
3	ZN	C	1202	1/1	0.99	0.05	-3.93	47,47,47,47	0

6.5 Other polymers (i)

There are no such residues in this entry.