



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 04:59 PM UTC

PDB ID : 3FV3 / pdb_00003fv3
Title : Secreted aspartic protease 1 from *Candida parapsilosis* in complex with pepstatin A
Authors : Dostal, J.; Brynda, J.; Hruskova-Heidingsfeldova, O.; Sieglova, I.; Pichova, I.; Rezacova, P.
Deposited on : 2009-01-15
Resolution : 1.85 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

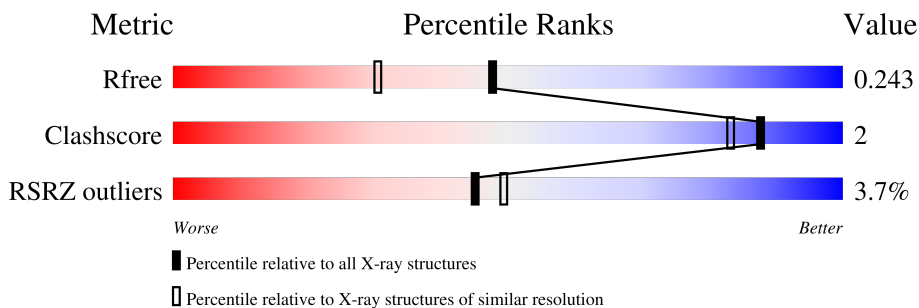
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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.



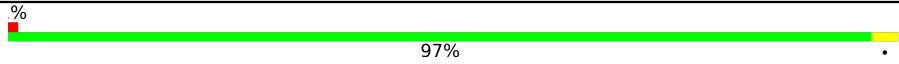








Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	339	 7% 98% 5%
1	B	339	 % 95% 5%
1	C	339	 13% 96% 5%
1	D	339	 4% 94% 5%
1	E	339	 % 98% 5%
1	F	339	 % 97% 5%
1	G	339	 % 97% 5%

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Mol	Chain	Length	Quality of chain
1	H	339	 97%
2	I	6	 67% 33%
2	J	6	 67% 33%
2	K	6	 67% 33%
2	L	6	 67% 33%
2	M	6	 67% 33%
2	N	6	 67% 33%
2	O	6	 67% 33%
2	P	6	 67% 33%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 23564 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 Sapp1p-secreted aspartic protease 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	339	Total 2585	C 1614	N 422	O 545	S 4	5	14	0
1	B	339	Total 2596	C 1620	N 423	O 549	S 4	5	16	0
1	C	339	Total 2598	C 1623	N 422	O 549	S 4	5	17	0
1	D	338	Total 2569	C 1603	N 420	O 542	S 4	6	12	0
1	E	339	Total 2621	C 1639	N 424	O 554	S 4	5	22	0
1	F	339	Total 2598	C 1623	N 421	O 550	S 4	5	17	0
1	G	339	Total 2609	C 1629	N 424	O 552	S 4	5	20	0
1	H	339	Total 2620	C 1637	N 426	O 553	S 4	5	22	0

- Molecule 2 is a protein called pepstatin A.

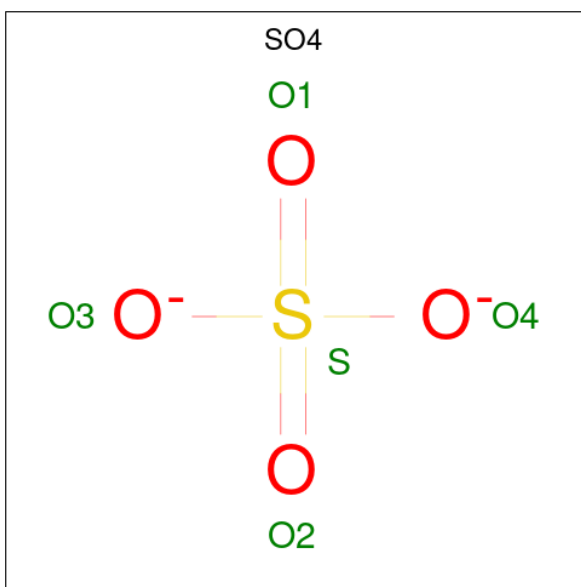
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	I	6	Total 48	C 34	N 5	O 9	0	0	0
2	J	6	Total 48	C 34	N 5	O 9	0	0	0
2	K	6	Total 48	C 34	N 5	O 9	0	0	0
2	L	6	Total 48	C 34	N 5	O 9	0	0	0
2	M	6	Total 48	C 34	N 5	O 9	0	0	0
2	N	6	Total 48	C 34	N 5	O 9	0	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	O	6	Total	C	N	O	0	0	0
			48	34	5	9			
2	P	6	Total	C	N	O	0	0	0
			48	34	5	9			

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



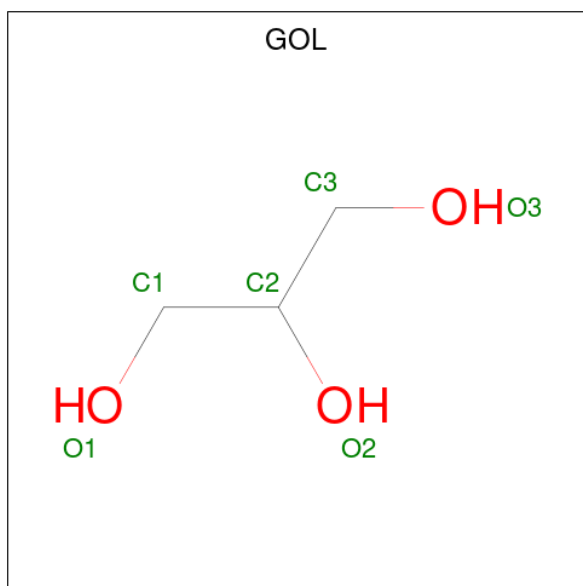
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		
3	G	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	G	1	Total	O	S	0	0
			5	4	1		
3	G	1	Total	O	S	0	0
			5	4	1		
3	G	1	Total	O	S	0	0
			5	4	1		
3	H	1	Total	O	S	0	0
			5	4	1		
3	H	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total 6	C 3	O 3	0	0
4	B	1	Total 6	C 3	O 3	0	0
4	C	1	Total 6	C 3	O 3	0	0
4	C	1	Total 6	C 3	O 3	0	0
4	C	1	Total 6	C 3	O 3	0	0
4	D	1	Total 6	C 3	O 3	0	0
4	D	1	Total 6	C 3	O 3	0	0
4	D	1	Total 6	C 3	O 3	0	0
4	E	1	Total 6	C 3	O 3	0	0
4	E	1	Total 6	C 3	O 3	0	0
4	E	1	Total 6	C 3	O 3	0	0
4	E	1	Total 6	C 3	O 3	0	0
4	E	1	Total 6	C 3	O 3	0	0
4	F	1	Total 6	C 3	O 3	0	0
4	G	1	Total 6	C 3	O 3	0	0
4	G	1	Total 6	C 3	O 3	0	0
4	H	1	Total 6	C 3	O 3	0	0
4	H	1	Total 6	C 3	O 3	0	0
4	H	1	Total 6	C 3	O 3	0	0
4	H	1	Total 6	C 3	O 3	0	0

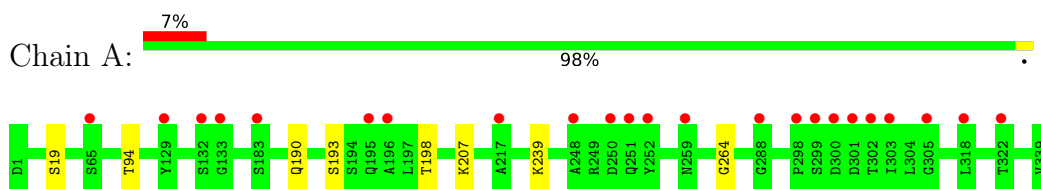
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	228	Total O 228 228	0	0
5	I	2	Total O 2 2	0	0
5	B	283	Total O 283 283	0	0
5	J	3	Total O 3 3	0	0
5	C	151	Total O 151 151	0	0
5	K	1	Total O 1 1	0	0
5	D	239	Total O 239 239	0	0
5	L	6	Total O 6 6	0	0
5	E	326	Total O 326 326	0	0
5	M	5	Total O 5 5	0	0
5	F	309	Total O 309 309	0	0
5	N	3	Total O 3 3	0	0
5	G	295	Total O 295 295	0	0
5	O	4	Total O 4 4	0	0
5	H	301	Total O 301 301	0	0
5	P	3	Total O 3 3	0	0

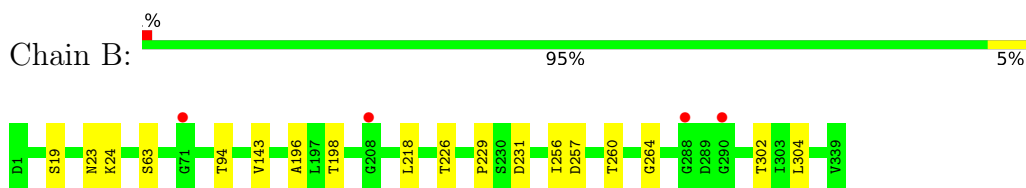
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

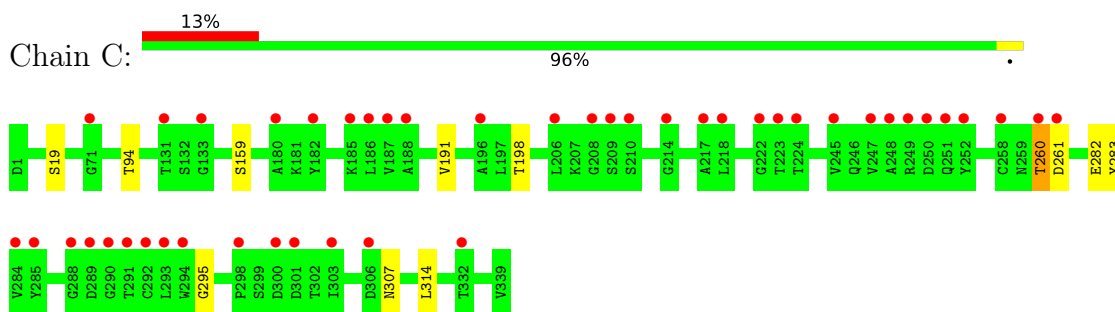
- Molecule 1: Sapp1p-secreted aspartic protease 1



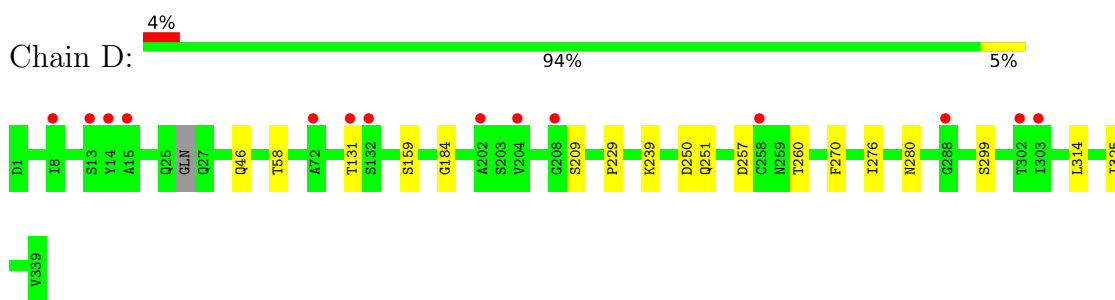
- Molecule 1: Sapp1p-secreted aspartic protease 1



- Molecule 1: Sapp1p-secreted aspartic protease 1



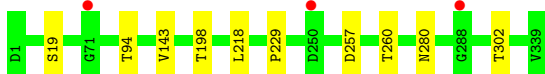
- Molecule 1: Sapp1p-secreted aspartic protease 1



- Molecule 1: Sapp1p-secreted aspartic protease 1



- Molecule 1: Sapp1p-secreted aspartic protease 1



- Molecule 1: Sapp1p-secreted aspartic protease 1



- Molecule 1: Sapp1p-secreted aspartic protease 1



- Molecule 2: pepstatin A



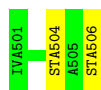
- Molecule 2: pepstatin A



- Molecule 2: pepstatin A



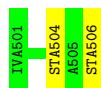
● Molecule 2: pepstatin A

Chain L:  67% 33%

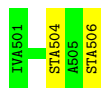
● Molecule 2: pepstatin A

Chain M:  67% 33%

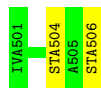
● Molecule 2: pepstatin A

Chain N:  67% 33%

● Molecule 2: pepstatin A

Chain O:  67% 33%

● Molecule 2: pepstatin A

Chain P:  67% 33%

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	86.49Å 194.25Å 97.15Å 90.00° 91.52° 90.00°	Depositor
Resolution (Å)	46.03 – 1.85 46.03 – 1.85	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.03-1.85) 99.3 (46.03-1.85)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 1.84Å)	Xtrriage
Refinement program	REFMAC 5.3.0037	Depositor
R, R_{free}	0.166 , 0.190 (Not available) , 0.243	Depositor DCC
R_{free} test set	1368 reflections (0.50%)	wwPDB-VP
Wilson B-factor (Å ²)	24.4	Xtrriage
Anisotropy	0.260	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 42.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.024 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	23564	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.91% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, STA, IVA, 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	0.59	0/2667	0.72	0/3630
1	B	0.64	1/2687 (0.0%)	0.76	0/3657
1	C	0.55	0/2692	0.74	1/3664 (0.0%)
1	D	0.69	0/2644	0.77	0/3597
1	E	0.70	0/2727	0.75	0/3711
1	F	0.70	1/2689 (0.0%)	0.76	0/3660
1	G	0.69	0/2706	0.75	0/3683
1	H	0.66	0/2726	0.73	0/3709
2	I	0.56	0/17	0.99	0/21
2	J	0.60	0/17	1.03	0/21
2	K	0.43	0/17	0.74	0/21
2	L	0.63	0/17	1.05	0/21
2	M	0.73	0/17	2.03	1/21 (4.8%)
2	N	0.61	0/17	1.07	0/21
2	O	0.55	0/17	1.00	0/21
2	P	0.58	0/17	0.94	0/21
All	All	0.65	2/21674 (0.0%)	0.75	2/29479 (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
2	I	0	2
2	J	0	2
2	K	0	2
2	L	0	2
2	M	0	2
2	N	0	2
2	O	0	2

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Mol	Chain	#Chirality outliers	#Planarity outliers
2	P	0	2
All	All	0	16

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	143	VAL	CA-CB	6.08	1.57	1.54
1	B	143	VAL	CA-CB	5.25	1.56	1.54

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	503	VAL	CA-CB-CG1	5.87	120.38	110.40
1	C	260	THR	N-CA-C	-5.57	102.05	110.24

There are no chirality outliers.

5 of 16 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	I	504	STA	Mainchain,Peptide
2	J	504	STA	Mainchain,Peptide
2	K	504	STA	Mainchain,Peptide
2	L	504	STA	Mainchain,Peptide
2	M	504	STA	Mainchain

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	2585	0	2526	6	0
1	B	2596	0	2540	15	0
1	C	2598	0	2546	14	0
1	D	2569	0	2503	12	0
1	E	2621	0	2579	10	0
1	F	2598	0	2542	8	0
1	G	2609	0	2558	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	H	2620	0	2578	9	0
2	I	48	0	60	0	0
2	J	48	0	60	0	0
2	K	48	0	60	0	0
2	L	48	0	60	0	0
2	M	48	0	60	0	0
2	N	48	0	60	0	0
2	O	48	0	60	0	0
2	P	48	0	60	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	D	10	0	0	1	0
3	E	10	0	0	0	0
3	F	10	0	0	0	0
3	G	25	0	0	0	0
3	H	10	0	0	0	0
4	A	18	0	24	1	0
4	B	24	0	32	2	0
4	C	18	0	24	0	0
4	D	18	0	24	2	0
4	E	24	0	32	4	0
4	F	6	0	8	1	0
4	G	12	0	16	1	0
4	H	30	0	40	3	0
5	A	228	0	0	1	0
5	B	283	0	0	2	0
5	C	151	0	0	1	0
5	D	239	0	0	3	0
5	E	326	0	0	3	0
5	F	309	0	0	1	0
5	G	295	0	0	2	0
5	H	301	0	0	4	0
5	I	2	0	0	0	0
5	J	3	0	0	0	0
5	K	1	0	0	0	0
5	L	6	0	0	0	0
5	M	5	0	0	0	0
5	N	3	0	0	0	0
5	O	4	0	0	0	0
5	P	3	0	0	0	0
All	All	23564	0	21052	84	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 2.

The worst 5 of 84 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:H:257:ASP:O	1:H:260[B]:THR:HG23	1.69	0.93
1:C:19:SER:HB2	1:C:94[A]:THR:HG23	1.53	0.88
1:F:198[A]:THR:HG21	5:F:632:HOH:O	1.74	0.85
1:E:19:SER:HB2	1:E:94[A]:THR:CG2	2.06	0.84
1:G:19:SER:HB2	1:G:94[A]:THR:CG2	2.07	0.83

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

16 non-standard protein/DNA/RNA residues 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	STA	I	506	2	11,11,11	0.68	0	11,14,14	0.96	1 (9%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	STA	L	506	2	11,11,11	0.61	0	11,14,14	1.29	1 (9%)
2	STA	K	504	2	10,10,11	0.67	0	9,12,14	0.99	0
2	STA	N	506	2	11,11,11	0.81	0	11,14,14	1.22	1 (9%)
2	STA	J	506	2	11,11,11	0.69	0	11,14,14	1.45	2 (18%)
2	STA	M	504	2	10,10,11	0.58	0	9,12,14	0.76	0
2	STA	N	504	2	10,10,11	0.88	0	9,12,14	0.73	0
2	STA	M	506	2	11,11,11	0.75	0	11,14,14	0.82	0
2	STA	J	504	2	10,10,11	0.70	0	9,12,14	1.00	0
2	STA	I	504	2	10,10,11	0.47	0	9,12,14	0.82	0
2	STA	O	504	2	10,10,11	0.64	0	9,12,14	0.88	0
2	STA	K	506	2	11,11,11	0.59	0	11,14,14	1.05	1 (9%)
2	STA	P	506	2	11,11,11	0.83	0	11,14,14	1.28	1 (9%)
2	STA	O	506	2	11,11,11	0.76	0	11,14,14	1.01	1 (9%)
2	STA	L	504	2	10,10,11	0.88	0	9,12,14	0.87	0
2	STA	P	504	2	10,10,11	0.52	0	9,12,14	0.73	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	STA	I	506	2	-	2/12/12/12	-
2	STA	L	506	2	-	1/12/12/12	-
2	STA	K	504	2	-	3/11/11/12	-
2	STA	N	506	2	-	0/12/12/12	-
2	STA	J	506	2	-	0/12/12/12	-
2	STA	M	504	2	-	1/11/11/12	-
2	STA	N	504	2	-	3/11/11/12	-
2	STA	M	506	2	-	0/12/12/12	-
2	STA	J	504	2	-	2/11/11/12	-
2	STA	I	504	2	-	3/11/11/12	-
2	STA	O	504	2	-	3/11/11/12	-
2	STA	K	506	2	-	0/12/12/12	-
2	STA	P	506	2	-	0/12/12/12	-
2	STA	O	506	2	-	0/12/12/12	-
2	STA	L	504	2	-	1/11/11/12	-
2	STA	P	504	2	-	2/11/11/12	-

There are no bond length outliers.

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	L	506	STA	CH-CM-C	-3.49	106.24	113.93
2	J	506	STA	CH-CM-C	-3.28	106.69	113.93
2	N	506	STA	CM-CH-CA	-2.38	109.25	112.93
2	P	506	STA	CM-CH-CA	2.35	116.56	112.93
2	K	506	STA	CH-CM-C	-2.17	109.14	113.93

There are no chirality outliers.

5 of 21 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	K	504	STA	O-C-CM-CH
2	M	504	STA	O-C-CM-CH
2	N	504	STA	N-CA-CB-CG
2	O	504	STA	O-C-CM-CH
2	I	504	STA	CA-CH-CM-C

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

40 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	$\# Z > 2$
3	SO4	G	340	-	4,4,4	0.27	0	6,6,6	0.41	0
4	GOL	B	344	-	5,5,5	0.35	0	5,5,5	0.62	0
4	GOL	H	344	-	5,5,5	0.51	0	5,5,5	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	G	343	-	4,4,4	0.25	0	6,6,6	0.15	0
3	SO4	E	341	-	4,4,4	0.26	0	6,6,6	0.13	0
4	GOL	E	345	-	5,5,5	0.33	0	5,5,5	0.55	0
3	SO4	F	341	-	4,4,4	0.22	0	6,6,6	0.50	0
4	GOL	A	343	-	5,5,5	0.45	0	5,5,5	0.30	0
3	SO4	A	340	-	4,4,4	0.25	0	6,6,6	0.17	0
4	GOL	C	340	-	5,5,5	0.34	0	5,5,5	0.39	0
4	GOL	H	342	-	5,5,5	0.40	0	5,5,5	0.45	0
4	GOL	B	342	-	5,5,5	0.33	0	5,5,5	0.36	0
4	GOL	H	343	-	5,5,5	0.28	0	5,5,5	0.64	0
4	GOL	H	346	-	5,5,5	0.37	0	5,5,5	0.26	0
4	GOL	D	343	-	5,5,5	0.43	0	5,5,5	0.24	0
3	SO4	H	340	-	4,4,4	0.25	0	6,6,6	0.17	0
4	GOL	D	344	-	5,5,5	0.39	0	5,5,5	0.20	0
4	GOL	E	343	-	5,5,5	0.34	0	5,5,5	0.39	0
3	SO4	E	340	-	4,4,4	0.22	0	6,6,6	0.42	0
3	SO4	D	341	-	4,4,4	0.20	0	6,6,6	0.48	0
4	GOL	C	342	-	5,5,5	0.36	0	5,5,5	0.39	0
3	SO4	D	340	-	4,4,4	0.24	0	6,6,6	0.14	0
3	SO4	H	341	-	4,4,4	0.23	0	6,6,6	0.27	0
4	GOL	C	341	-	5,5,5	0.41	0	5,5,5	0.35	0
4	GOL	G	345	-	5,5,5	0.31	0	5,5,5	0.81	0
4	GOL	H	345	-	5,5,5	0.38	0	5,5,5	0.41	0
4	GOL	B	343	-	5,5,5	0.54	0	5,5,5	0.73	0
3	SO4	G	342	-	4,4,4	0.25	0	6,6,6	0.18	0
4	GOL	E	342	-	5,5,5	0.53	0	5,5,5	0.49	0
3	SO4	B	340	-	4,4,4	0.22	0	6,6,6	0.24	0
4	GOL	F	342	-	5,5,5	0.36	0	5,5,5	0.36	0
4	GOL	E	344	-	5,5,5	0.55	0	5,5,5	0.91	0
3	SO4	G	341	-	4,4,4	0.26	0	6,6,6	0.14	0
4	GOL	D	342	-	5,5,5	0.29	0	5,5,5	0.21	0
4	GOL	A	341	-	5,5,5	0.28	0	5,5,5	0.40	0
4	GOL	B	341	-	5,5,5	0.38	0	5,5,5	0.25	0
3	SO4	F	340	-	4,4,4	0.24	0	6,6,6	0.30	0
3	SO4	G	344	-	4,4,4	0.27	0	6,6,6	0.16	0
4	GOL	A	342	-	5,5,5	0.38	0	5,5,5	0.33	0
4	GOL	G	346	-	5,5,5	0.39	0	5,5,5	0.24	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	B	344	-	-	2/4/4/4	-
4	GOL	H	344	-	-	2/4/4/4	-
4	GOL	E	345	-	-	1/4/4/4	-
4	GOL	A	343	-	-	2/4/4/4	-
4	GOL	C	340	-	-	4/4/4/4	-
4	GOL	H	342	-	-	4/4/4/4	-
4	GOL	B	342	-	-	2/4/4/4	-
4	GOL	H	343	-	-	4/4/4/4	-
4	GOL	H	346	-	-	4/4/4/4	-
4	GOL	D	343	-	-	1/4/4/4	-
4	GOL	D	344	-	-	0/4/4/4	-
4	GOL	E	343	-	-	0/4/4/4	-
4	GOL	C	342	-	-	0/4/4/4	-
4	GOL	C	341	-	-	1/4/4/4	-
4	GOL	G	345	-	-	2/4/4/4	-
4	GOL	H	345	-	-	4/4/4/4	-
4	GOL	B	343	-	-	4/4/4/4	-
4	GOL	E	342	-	-	4/4/4/4	-
4	GOL	F	342	-	-	2/4/4/4	-
4	GOL	E	344	-	-	2/4/4/4	-
4	GOL	D	342	-	-	2/4/4/4	-
4	GOL	A	341	-	-	0/4/4/4	-
4	GOL	B	341	-	-	0/4/4/4	-
4	GOL	A	342	-	-	0/4/4/4	-
4	GOL	G	346	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 49 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	343	GOL	O1-C1-C2-O2
4	A	343	GOL	O1-C1-C2-C3
4	B	342	GOL	C1-C2-C3-O3
4	B	343	GOL	C1-C2-C3-O3
4	B	344	GOL	C1-C2-C3-O3

There are no ring outliers.

12 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	345	GOL	1	0
4	A	343	GOL	1	0
4	H	342	GOL	2	0
4	H	343	GOL	1	0
4	D	343	GOL	1	0
4	D	344	GOL	1	0
3	D	341	SO4	1	0
4	G	345	GOL	1	0
4	B	343	GOL	1	0
4	F	342	GOL	1	0
4	E	344	GOL	3	0
4	B	341	GOL	1	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	339/339 (100%)	0.51	23 (6%) 23 25	6, 13, 20, 27	16 (4%)
1	B	339/339 (100%)	-0.05	4 (1%) 76 80	6, 12, 20, 29	18 (5%)
1	C	339/339 (100%)	0.94	45 (13%) 7 6	5, 13, 20, 31	17 (5%)
1	D	338/339 (99%)	0.55	14 (4%) 41 45	6, 12, 21, 30	16 (4%)
1	E	339/339 (100%)	-0.11	4 (1%) 76 80	6, 12, 20, 29	23 (6%)
1	F	339/339 (100%)	-0.17	3 (0%) 81 84	6, 12, 21, 34	21 (6%)
1	G	339/339 (100%)	0.06	4 (1%) 76 80	6, 12, 20, 28	23 (6%)
1	H	339/339 (100%)	0.01	4 (1%) 76 80	6, 12, 19, 30	24 (7%)
2	I	3/6 (50%)	-0.01	0 100 100	12, 12, 14, 15	0
2	J	3/6 (50%)	-0.14	0 100 100	9, 9, 11, 11	0
2	K	3/6 (50%)	0.11	0 100 100	15, 15, 16, 20	0
2	L	3/6 (50%)	-0.15	0 100 100	7, 7, 8, 15	0
2	M	3/6 (50%)	-0.01	0 100 100	9, 9, 9, 10	0
2	N	3/6 (50%)	-0.33	0 100 100	7, 7, 8, 9	0
2	O	3/6 (50%)	-0.30	0 100 100	8, 8, 10, 12	0
2	P	3/6 (50%)	-0.24	0 100 100	8, 8, 8, 10	0
All	All	2735/2760 (99%)	0.21	101 (3%) 45 49	5, 12, 20, 34	158 (5%)

The worst 5 of 101 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	288	GLY	4.5
1	C	222	GLY	4.4
1	D	288	GLY	4.1
1	C	251[A]	GLN	4.1
1	C	285	TYR	4.1

6.2 Non-standard residues in protein, DNA, RNA chains [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. 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	B-factors(Å ²)	Q<0.9
2	STA	K	506	12/12	0.81	0.13	24,29,42,44	0
2	STA	L	506	12/12	0.85	0.12	14,21,40,40	0
2	STA	P	506	12/12	0.86	0.12	11,17,35,37	0
2	STA	M	506	12/12	0.88	0.11	10,14,36,36	0
2	STA	O	506	12/12	0.89	0.11	11,17,37,38	0
2	STA	I	506	12/12	0.89	0.10	15,19,38,39	0
2	STA	J	506	12/12	0.90	0.10	11,16,36,37	0
2	STA	N	506	12/12	0.90	0.10	8,12,29,32	0
2	STA	K	504	11/12	0.93	0.07	12,14,16,16	0
2	STA	L	504	11/12	0.95	0.06	6,7,10,10	0
2	STA	M	504	11/12	0.96	0.06	5,7,9,9	0
2	STA	N	504	11/12	0.96	0.06	4,5,8,8	0
2	STA	O	504	11/12	0.96	0.06	5,6,8,8	0
2	STA	P	504	11/12	0.96	0.06	5,7,8,9	0
2	STA	I	504	11/12	0.96	0.06	8,11,11,12	0
2	STA	J	504	11/12	0.96	0.06	5,6,9,10	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
3	SO4	G	341	5/5	0.64	0.15	76,76,76,76	0
3	SO4	D	340	5/5	0.69	0.17	91,91,92,92	0
3	SO4	E	341	5/5	0.74	0.14	65,65,65,66	0
3	SO4	G	342	5/5	0.74	0.13	75,75,75,76	0
3	SO4	G	344	5/5	0.74	0.18	73,74,75,75	0
4	GOL	D	344	6/6	0.74	0.14	60,61,62,62	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	GOL	H	346	6/6	0.74	0.21	53,54,55,57	0
3	SO4	A	340	5/5	0.75	0.11	93,93,94,94	0
4	GOL	C	342	6/6	0.76	0.12	56,56,56,56	0
4	GOL	B	341	6/6	0.78	0.14	42,44,45,45	0
4	GOL	E	342	6/6	0.79	0.17	36,38,40,40	0
4	GOL	A	343	6/6	0.80	0.16	59,59,59,59	0
4	GOL	C	341	6/6	0.80	0.13	46,46,46,46	0
4	GOL	G	346	6/6	0.81	0.14	46,47,48,48	0
4	GOL	H	342	6/6	0.81	0.16	46,47,49,49	0
3	SO4	F	340	5/5	0.81	0.10	60,61,61,62	0
4	GOL	F	342	6/6	0.82	0.12	34,42,44,45	0
4	GOL	D	343	6/6	0.82	0.15	35,43,45,49	0
4	GOL	G	345	6/6	0.83	0.17	26,34,36,40	0
4	GOL	C	340	6/6	0.83	0.15	54,57,57,58	0
4	GOL	D	342	6/6	0.83	0.12	35,37,38,40	0
3	SO4	H	340	5/5	0.83	0.11	67,67,67,68	0
4	GOL	B	342	6/6	0.84	0.14	40,42,43,44	0
4	GOL	E	343	6/6	0.86	0.11	36,37,38,38	0
3	SO4	B	340	5/5	0.86	0.11	49,51,51,51	0
3	SO4	G	343	5/5	0.87	0.09	60,60,60,61	0
4	GOL	A	341	6/6	0.87	0.11	38,40,41,41	0
3	SO4	D	341	5/5	0.87	0.11	51,52,53,53	0
3	SO4	E	340	5/5	0.88	0.10	46,47,49,49	0
4	GOL	H	343	6/6	0.88	0.10	32,37,38,40	0
4	GOL	H	344	6/6	0.88	0.15	26,27,30,30	0
3	SO4	G	340	5/5	0.88	0.11	43,43,45,45	0
4	GOL	E	344	6/6	0.90	0.14	19,23,24,26	0
4	GOL	B	344	6/6	0.90	0.11	30,37,38,39	0
4	GOL	A	342	6/6	0.90	0.11	40,42,42,43	0
4	GOL	H	345	6/6	0.90	0.12	39,40,40,41	0
4	GOL	B	343	6/6	0.90	0.15	24,27,29,29	0
3	SO4	F	341	5/5	0.91	0.13	14,17,18,19	5
3	SO4	H	341	5/5	0.95	0.09	14,16,17,20	5
4	GOL	E	345	6/6	0.97	0.08	18,27,27,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.