



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 04:41 PM GMT

PDB ID : 3FV3  
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 full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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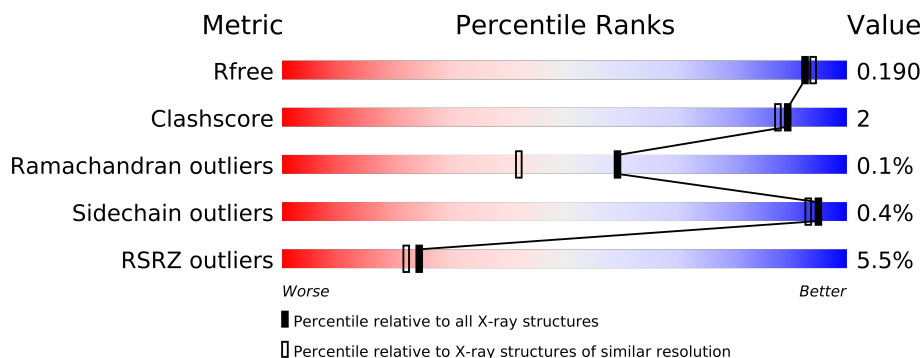
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

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))
$R_{free}$	66092	1269 (1.86-1.86)
Clashscore	79885	1470 (1.86-1.86)
Ramachandran outliers	78287	1451 (1.86-1.86)
Sidechain outliers	78261	1451 (1.86-1.86)
RSRZ outliers	66119	1269 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	339	
1	B	339	
1	C	339	
1	D	339	
1	E	339	
1	F	339	
1	G	339	
1	H	339	
2	I	6	
2	J	6	
2	K	6	
2	L	6	
2	M	6	
2	N	6	

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Mol	Chain	Length	Quality of chain
2	O	6	
2	P	6	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	SO4	D	341	-	X
3	SO4	E	340	-	X
3	SO4	F	341	-	X
4	GOL	B	343	-	X
4	GOL	D	343	-	X
4	GOL	F	342	-	X
4	GOL	G	345	-	X
4	GOL	G	346	-	X
4	GOL	H	346	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 23564 atoms, of which 0 are hydrogen and 0 are deuterium.

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

- Molecule 2 is a protein called pepstatin A.

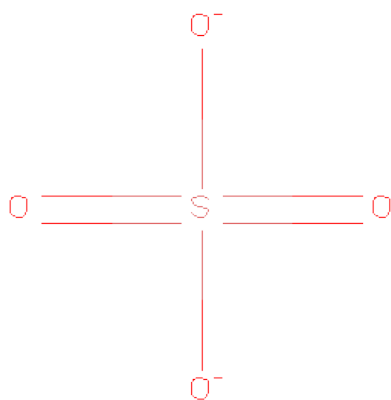
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	I	6	Total	C	N	O	0	0	0
			48	34	5	9			
2	J	6	Total	C	N	O	0	0	0
			48	34	5	9			
2	K	6	Total	C	N	O	0	0	0
			48	34	5	9			
2	L	6	Total	C	N	O	0	0	0
			48	34	5	9			
2	M	6	Total	C	N	O	0	0	0
			48	34	5	9			
2	N	6	Total	C	N	O	0	0	0
			48	34	5	9			

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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 (three-letter code: SO4) (formula: O<sub>4</sub>S).



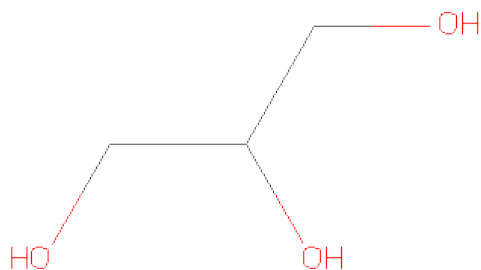
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		
3	G	1	Total	O	S	0	0
			5	4	1		
3	G	1	Total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

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



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

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

- Molecule 5 is water.

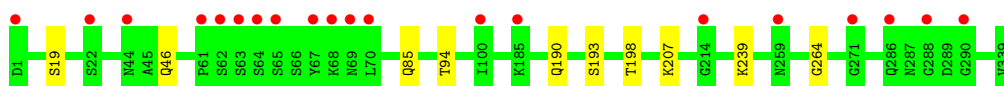
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	228	Total O 228 228	0	0
5	I	2	Total O 2 2	0	0
5	B	283	Total O 283 283	0	0
5	J	3	Total O 3 3	0	0
5	C	151	Total O 151 151	0	0
5	K	1	Total O 1 1	0	0
5	D	239	Total O 239 239	0	0
5	L	6	Total O 6 6	0	0
5	E	326	Total O 326 326	0	0
5	M	5	Total O 5 5	0	0
5	F	309	Total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

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.

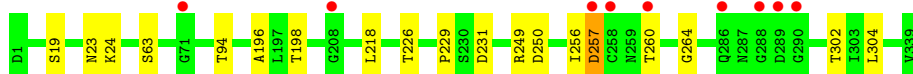
- Molecule 1: Sapp1p-secreted aspartic protease 1

Chain A: 



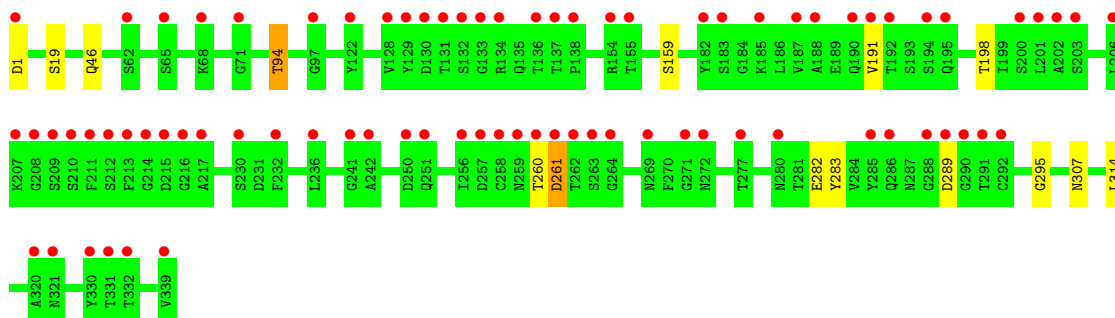
- Molecule 1: Sapp1p-secreted aspartic protease 1

Chain B: 



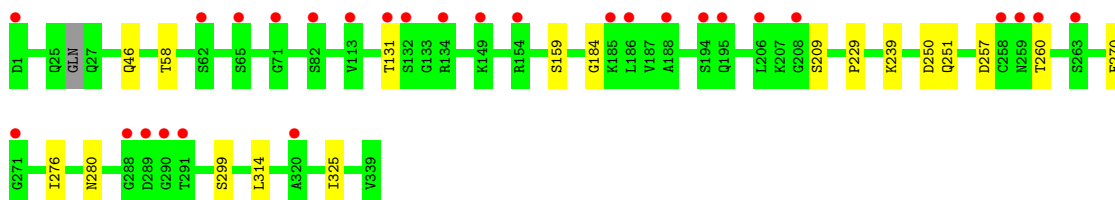
- Molecule 1: Sapp1p-secreted aspartic protease 1

Chain C: 



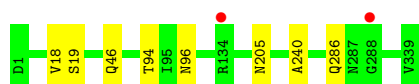
- Molecule 1: Sapp1p-secreted aspartic protease 1

Chain D: 



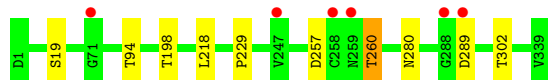
- Molecule 1: Sapp1p-secreted aspartic protease 1

Chain E: 



- Molecule 1: Sapp1p-secreted aspartic protease 1

Chain F:



- Molecule 1: Sapp1p-secreted aspartic protease 1

Chain G:



- Molecule 1: Sapp1p-secreted aspartic protease 1

Chain H:



- Molecule 2: pepstatin A

Chain I:



- Molecule 2: pepstatin A

Chain J:



- Molecule 2: pepstatin A

Chain K:



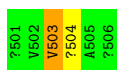
- Molecule 2: pepstatin A

Chain L:



- Molecule 2: pepstatin A

Chain M:



- Molecule 2: pepstatin A

Chain N:



- Molecule 2: pepstatin A

Chain O:



- Molecule 2: pepstatin A

Chain P:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	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) 99.1 (46.03-1.85)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.36 (at 1.84Å)	Xtriage
Refinement program	REFMAC 5.3.0037	Depositor
R, $R_{free}$	0.166 , 0.190 0.168 , 0.190	Depositor DCC
$R_{free}$ test set	1368 reflections (0.51%)	DCC
Wilson B-factor (Å <sup>2</sup> )	24.4	Xtriage
Anisotropy	0.260	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 31.8	EDS
Estimated twinning fraction	0.024 for h,-k,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 271028 reflections (0.000%)	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	23564	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	15.0	wwPDB-VP

Xtriage'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.*

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SO4, IVA, STA

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	0.50	0/2667	0.57	0/3630
1	B	0.56	0/2687	0.61	0/3657
1	C	0.46	0/2692	0.57	0/3664
1	D	0.59	0/2644	0.61	0/3597
1	E	0.62	0/2727	0.63	0/3711
1	F	0.61	0/2689	0.63	0/3660
1	G	0.59	0/2706	0.60	0/3683
1	H	0.58	0/2726	0.61	0/3709
2	I	0.50	0/17	1.02	0/21
2	J	0.58	0/17	1.04	0/21
2	K	0.41	0/17	0.78	0/21
2	L	0.64	0/17	1.07	0/21
2	M	0.70	0/17	2.06	2/21 (9.5%)
2	N	0.59	0/17	1.02	0/21
2	O	0.62	0/17	1.01	0/21
2	P	0.53	0/17	0.94	0/21
All	All	0.57	0/21674	0.61	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

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	503	VAL	CA-CB-CG1	6.32	120.38	110.90
2	M	503	VAL	CA-CB-CG2	5.07	118.51	110.90

There are no chirality outliers.

All (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,Peptide
2	N	504	STA	Mainchain,Peptide
2	O	504	STA	Mainchain,Peptide
2	P	504	STA	Mainchain,Peptide

## 5.2 Close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2585	0	2526	7	0
1	B	2596	0	2540	19	0
1	C	2598	0	2546	14	0
1	D	2569	0	2503	12	0
1	E	2621	0	2579	13	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	2	0
5	B	283	0	0	5	0
5	C	151	0	0	1	0
5	D	239	0	0	3	0
5	E	326	0	0	6	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	92	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including

hydrogens) of the entry. The overall clashscore for this entry is 2.

All (92) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:286[A]:GLN:OE1	5:E:2136:HOH:O	1.66	1.11
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:A:85[A]:GLN:OE1	5:A:1738:HOH:O	1.95	0.83
1:G:19:SER:HB2	1:G:94[A]:THR:CG2	2.07	0.83
1:F:19:SER:HB2	1:F:94[A]:THR:HG23	1.60	0.83
1:H:256:ILE:HG12	1:H:260[B]:THR:HG21	1.60	0.82
1:D:250:ASP:HB3	1:D:251:GLN:OE1	1.81	0.81
1:E:286[B]:GLN:NE2	5:E:1374:HOH:O	1.91	0.81
1:C:260:THR:CG2	1:C:261:ASP:HB3	2.11	0.81
1:F:257:ASP:O	1:F:260:THR:HG22	1.83	0.79
1:F:19:SER:HB2	1:F:94[A]:THR:CG2	2.13	0.79
1:C:260:THR:HA	1:C:261:ASP:HB2	1.64	0.78
1:C:260:THR:HG22	1:C:261:ASP:HB3	1.66	0.77
1:E:46[A]:GLN:HG2	5:E:2137:HOH:O	1.84	0.77
1:B:249[A]:ARG:HD2	5:B:1710:HOH:O	1.32	0.76
1:C:260:THR:HA	1:C:261:ASP:CB	2.16	0.75
1:G:19:SER:HB2	1:G:94[A]:THR:HG23	1.73	0.71
1:A:19:SER:HB2	1:A:94[A]:THR:HG23	1.72	0.70
1:D:46[B]:GLN:HG2	5:D:2144:HOH:O	1.92	0.68
1:B:19:SER:HB2	1:B:94[A]:THR:CG2	2.24	0.68
1:D:131:THR:HG22	5:D:815:HOH:O	1.94	0.66
1:A:19:SER:HB2	1:A:94[A]:THR:CG2	2.25	0.66
1:E:19:SER:HB2	1:E:94[A]:THR:HG22	1.80	0.64
1:G:52:ASP:OD1	5:G:2135:HOH:O	2.15	0.63
1:C:19:SER:HB2	1:C:94[A]:THR:CG2	2.28	0.63
1:B:264:GLY:HA3	4:B:341:GOL:H11	1.82	0.62
1:B:249[B]:ARG:HD3	1:B:250:ASP:OD1	2.01	0.61
1:F:280:ASN:HB3	4:F:342:GOL:H31	1.83	0.61
1:G:231:ASP:HB3	5:G:1678:HOH:O	2.01	0.61
1:B:257:ASP:HB2	1:B:260[B]:THR:HG23	1.82	0.61
1:E:18[B]:VAL:HA	4:E:344:GOL:H32	1.83	0.60
1:F:198[A]:THR:HG22	1:F:218:LEU:HA	1.83	0.60
1:E:205[B]:ASN:ND2	5:E:1276:HOH:O	2.35	0.59
1:H:198:THR:HG21	5:H:1310:HOH:O	2.03	0.58
1:B:19:SER:HB2	1:B:94[A]:THR:HG22	1.87	0.57
1:A:207:LYS:HB2	1:A:239:LYS:HE3	1.86	0.57
1:C:283:TYR:HA	1:C:295:GLY:HA3	1.87	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:94[A]:THR:HG23	4:E:344:GOL:H11	1.87	0.57
1:B:256:ILE:HG12	1:B:260[B]:THR:HG21	1.87	0.57
1:B:198:THR:HG21	5:B:691:HOH:O	2.05	0.57
1:D:276:ILE:HD11	1:D:325[A]:ILE:HG22	1.88	0.55
1:B:198:THR:HG22	1:B:218:LEU:HA	1.89	0.55
1:D:58:THR:OG1	3:D:341:SO4:O1	2.19	0.55
1:H:207:LYS:HD3	4:H:342:GOL:H31	1.89	0.54
1:H:198:THR:HG22	1:H:218:LEU:HA	1.90	0.54
1:F:198[A]:THR:HG22	1:F:218:LEU:CA	2.38	0.53
1:H:87[B]:THR:HG22	5:H:658:HOH:O	2.09	0.53
1:E:46[B]:GLN:HG2	5:E:2137:HOH:O	2.08	0.53
1:H:283:TYR:HB2	4:H:343:GOL:H31	1.92	0.52
1:C:260:THR:CA	1:C:261:ASP:CB	2.86	0.52
1:D:229:PRO:HA	1:D:299:SER:O	2.10	0.52
1:B:23:ASN:HD21	1:B:63:SER:HB3	1.77	0.50
1:E:19:SER:HB2	1:E:94[A]:THR:HG23	1.89	0.50
1:D:209:SER:HB2	1:D:239:LYS:HE2	1.93	0.49
1:B:231:ASP:HB2	5:B:1670:HOH:O	2.12	0.49
1:H:194[B]:SER:OG	1:H:195[B]:GLN:NE2	2.46	0.48
1:C:260:THR:HG23	1:C:261:ASP:HB3	1.93	0.48
1:A:264:GLY:CA	4:A:343:GOL:H32	2.44	0.48
1:E:286[B]:GLN:CD	5:E:1374:HOH:O	2.43	0.48
1:B:249[A]:ARG:CD	5:B:1710:HOH:O	1.85	0.47
1:B:24:LYS:HD3	1:B:94[A]:THR:HG21	1.97	0.47
1:G:19:SER:HB2	1:G:94[A]:THR:HG22	1.93	0.47
1:G:262:THR:HG22	4:G:345:GOL:O1	2.15	0.47
1:A:193:SER:HB3	1:A:198[A]:THR:HG22	1.96	0.46
4:H:342:GOL:H32	5:H:899:HOH:O	2.16	0.46
1:B:94[A]:THR:HG23	4:B:343:GOL:H11	1.97	0.46
1:C:282:GLU:HB3	1:C:307:ASN:HB2	1.97	0.46
1:D:184:GLY:H	4:D:344:GOL:H2	1.81	0.46
1:D:270:PHE:N	5:D:928:HOH:O	2.48	0.45
1:D:280:ASN:HB3	4:D:343:GOL:H2	1.98	0.45
1:H:87[B]:THR:HG23	5:H:499:HOH:O	2.17	0.44
1:C:159:SER:HA	1:C:314:LEU:O	2.18	0.44
1:G:282:GLU:HB3	1:G:307:ASN:HB2	1.99	0.43
1:G:226:THR:HG23	1:G:304:LEU:HD23	2.00	0.43
1:B:249[B]:ARG:NE	5:B:1710:HOH:O	2.43	0.43
1:B:229:PRO:HD3	1:B:302:THR:OG1	2.18	0.42
1:D:257:ASP:O	1:D:260:THR:HG22	2.19	0.42
1:E:96:ASN:H	4:E:344:GOL:C3	2.32	0.42
1:C:94[A]:THR:HG22	5:C:1219:HOH:O	2.20	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:191:VAL:HA	1:C:198:THR:O	2.20	0.41
1:B:226:THR:HG23	1:B:304:LEU:HD23	2.03	0.41
1:G:283:TYR:HA	1:G:295:GLY:HA3	2.03	0.41
1:B:196:ALA:HB1	1:B:218:LEU:HD13	2.02	0.41
1:B:198:THR:HG22	1:B:218:LEU:CA	2.50	0.40
1:E:240:ALA:HB1	4:E:345:GOL:H12	2.02	0.40
1:C:260:THR:HA	1:C:261:ASP:HB3	2.02	0.40
1:A:190:GLN:NE2	5:A:1865:HOH:O	2.44	0.40
1:D:159:SER:HA	1:D:314:LEU:O	2.22	0.40
1:F:229:PRO:HD3	1:F:302:THR:OG1	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	351/339 (104%)	342 (97%)	9 (3%)	0	100	100
1	B	353/339 (104%)	343 (97%)	10 (3%)	0	100	100
1	C	354/339 (104%)	346 (98%)	6 (2%)	2 (1%)	33	16
1	D	346/339 (102%)	339 (98%)	7 (2%)	0	100	100
1	E	359/339 (106%)	352 (98%)	7 (2%)	0	100	100
1	F	354/339 (104%)	348 (98%)	5 (1%)	1 (0%)	50	32
1	G	357/339 (105%)	352 (99%)	5 (1%)	0	100	100
1	H	359/339 (106%)	350 (98%)	9 (2%)	0	100	100
2	I	3/6 (50%)	3 (100%)	0	0	100	100
2	J	3/6 (50%)	3 (100%)	0	0	100	100
2	K	3/6 (50%)	3 (100%)	0	0	100	100
2	L	3/6 (50%)	2 (67%)	1 (33%)	0	100	100
2	M	3/6 (50%)	3 (100%)	0	0	100	100
2	N	3/6 (50%)	2 (67%)	1 (33%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	O	3/6 (50%)	2 (67%)	1 (33%)	0	100	100
2	P	3/6 (50%)	3 (100%)	0	0	100	100
All	All	2857/2760 (104%)	2793 (98%)	61 (2%)	3 (0%)	59	42

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	261	ASP
1	C	289	ASP
1	F	289	ASP

### 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	294/281 (105%)	293 (100%)	1 (0%)	96	94
1	B	297/281 (106%)	296 (100%)	1 (0%)	96	94
1	C	298/281 (106%)	293 (98%)	5 (2%)	73	59
1	D	292/281 (104%)	292 (100%)	0	100	100
1	E	303/281 (108%)	303 (100%)	0	100	100
1	F	297/281 (106%)	296 (100%)	1 (0%)	96	94
1	G	300/281 (107%)	300 (100%)	0	100	100
1	H	302/281 (108%)	301 (100%)	1 (0%)	96	94
2	I	2/2 (100%)	2 (100%)	0	100	100
2	J	2/2 (100%)	2 (100%)	0	100	100
2	K	2/2 (100%)	2 (100%)	0	100	100
2	L	2/2 (100%)	2 (100%)	0	100	100
2	M	2/2 (100%)	1 (50%)	1 (50%)	0	0
2	N	2/2 (100%)	2 (100%)	0	100	100
2	O	2/2 (100%)	2 (100%)	0	100	100
2	P	2/2 (100%)	2 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	2399/2264 (106%)	2389 (100%)	10 (0%)	95 93

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	GLN
1	B	257	ASP
1	C	1	ASP
1	C	46[A]	GLN
1	C	46[B]	GLN
1	C	94[A]	THR
1	C	94[B]	THR
2	M	503	VAL
1	F	260	THR
1	H	212	SER

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

Mol	Chain	Res	Type
1	B	115	GLN
1	C	287	ASN
1	F	115	GLN
1	F	287	ASN
1	H	46	GLN
1	H	115	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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	504	2	10,10,11	4.92	1 (10%)	10,12,14	0.84	0
2	STA	I	506	2	11,11,11	0.65	0	14,14,14	1.01	1 (7%)
2	STA	J	504	2	10,10,11	5.42	2 (20%)	10,12,14	1.01	0
2	STA	J	506	2	11,11,11	0.58	0	14,14,14	1.34	1 (7%)
2	STA	K	504	2	10,10,11	5.41	2 (20%)	10,12,14	0.90	1 (10%)
2	STA	K	506	2	11,11,11	0.52	0	14,14,14	1.01	1 (7%)
2	STA	L	504	2	10,10,11	5.11	3 (30%)	10,12,14	1.04	0
2	STA	L	506	2	11,11,11	0.57	0	14,14,14	1.28	1 (7%)
2	STA	M	504	2	10,10,11	6.03	2 (20%)	10,12,14	0.96	0
2	STA	M	506	2	11,11,11	0.73	0	14,14,14	0.77	0
2	STA	N	504	2	10,10,11	4.96	3 (30%)	10,12,14	1.01	0
2	STA	N	506	2	11,11,11	0.66	0	14,14,14	1.09	1 (7%)
2	STA	O	504	2	10,10,11	5.61	1 (10%)	10,12,14	0.83	0
2	STA	O	506	2	11,11,11	0.73	0	14,14,14	1.02	1 (7%)
2	STA	P	504	2	10,10,11	5.14	1 (10%)	10,12,14	0.77	0
2	STA	P	506	2	11,11,11	0.72	0	14,14,14	1.17	3 (21%)

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	504	2	-	0/10/11/12	0/0/0/0
2	STA	I	506	2	-	0/12/12/12	0/0/0/0
2	STA	J	504	2	-	0/10/11/12	0/0/0/0
2	STA	J	506	2	-	0/12/12/12	0/0/0/0
2	STA	K	504	2	-	0/10/11/12	0/0/0/0
2	STA	K	506	2	-	0/12/12/12	0/0/0/0
2	STA	L	504	2	-	0/10/11/12	0/0/0/0
2	STA	L	506	2	-	0/12/12/12	0/0/0/0
2	STA	M	504	2	-	0/10/11/12	0/0/0/0
2	STA	M	506	2	-	0/12/12/12	0/0/0/0
2	STA	N	504	2	-	0/10/11/12	0/0/0/0
2	STA	N	506	2	-	0/12/12/12	0/0/0/0
2	STA	O	504	2	-	0/10/11/12	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	STA	O	506	2	-	0/12/12/12	0/0/0/0
2	STA	P	504	2	-	0/10/11/12	0/0/0/0
2	STA	P	506	2	-	0/12/12/12	0/0/0/0

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	M	504	STA	O-C	18.89	1.24	1.11
2	O	504	STA	O-C	17.42	1.23	1.11
2	J	504	STA	O-C	16.81	1.23	1.11
2	K	504	STA	O-C	16.77	1.22	1.11
2	P	504	STA	O-C	16.07	1.22	1.11
2	I	504	STA	O-C	15.37	1.22	1.11
2	L	504	STA	O-C	15.37	1.22	1.11
2	N	504	STA	O-C	15.06	1.21	1.11
2	L	504	STA	CM-CH	3.99	1.56	1.53
2	N	504	STA	CB-CA	2.83	1.56	1.53
2	J	504	STA	CB-CA	2.71	1.55	1.53
2	K	504	STA	CM-CH	2.68	1.55	1.53
2	L	504	STA	CB-CA	2.57	1.55	1.53
2	N	504	STA	CM-CH	2.41	1.55	1.53
2	M	504	STA	CB-CA	2.05	1.55	1.53

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	506	STA	CH-CM-C	-3.76	106.24	113.93
2	J	506	STA	CH-CM-C	-3.54	106.69	113.93
2	P	506	STA	CM-CH-CA	2.36	116.56	112.65
2	K	506	STA	CH-CM-C	-2.34	109.14	113.93
2	O	506	STA	CH-CM-C	-2.20	109.44	113.93
2	I	506	STA	CH-CM-C	-2.19	109.44	113.93
2	P	506	STA	OH-CH-CA	-2.13	105.91	109.69
2	K	504	STA	OH-CH-CA	-2.08	106.00	109.69
2	N	506	STA	CM-CH-CA	-2.05	109.25	112.65
2	P	506	STA	OH-CH-CM	-2.02	104.88	109.68

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

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	A	340	-	4,4,4	0.20	0	6,6,6	0.17	0
4	GOL	A	341	-	5,5,5	0.27	0	5,5,5	0.35	0
4	GOL	A	342	-	5,5,5	0.31	0	5,5,5	0.45	0
4	GOL	A	343	-	5,5,5	0.42	0	5,5,5	0.16	0
3	SO4	B	340	-	4,4,4	0.17	0	6,6,6	0.21	0
4	GOL	B	341	-	5,5,5	0.35	0	5,5,5	0.36	0
4	GOL	B	342	-	5,5,5	0.30	0	5,5,5	0.37	0
4	GOL	B	343	-	5,5,5	0.50	0	5,5,5	0.53	0
4	GOL	B	344	-	5,5,5	0.32	0	5,5,5	0.59	0
4	GOL	C	340	-	5,5,5	0.31	0	5,5,5	0.38	0
4	GOL	C	341	-	5,5,5	0.39	0	5,5,5	0.30	0
4	GOL	C	342	-	5,5,5	0.33	0	5,5,5	0.34	0
3	SO4	D	340	-	4,4,4	0.17	0	6,6,6	0.10	0
3	SO4	D	341	-	4,4,4	0.22	0	6,6,6	0.37	0
4	GOL	D	342	-	5,5,5	0.27	0	5,5,5	0.24	0
4	GOL	D	343	-	5,5,5	0.39	0	5,5,5	0.36	0
4	GOL	D	344	-	5,5,5	0.35	0	5,5,5	0.24	0
3	SO4	E	340	-	4,4,4	0.29	0	6,6,6	0.26	0
3	SO4	E	341	-	4,4,4	0.14	0	6,6,6	0.12	0
4	GOL	E	342	-	5,5,5	0.49	0	5,5,5	0.47	0
4	GOL	E	343	-	5,5,5	0.30	0	5,5,5	0.32	0
4	GOL	E	344	-	5,5,5	0.53	0	5,5,5	0.68	0
4	GOL	E	345	-	5,5,5	0.25	0	5,5,5	0.66	0
3	SO4	F	340	-	4,4,4	0.18	0	6,6,6	0.29	0
3	SO4	F	341	-	4,4,4	0.32	0	6,6,6	0.38	0
4	GOL	F	342	-	5,5,5	0.33	0	5,5,5	0.29	0
3	SO4	G	340	-	4,4,4	0.11	0	6,6,6	0.33	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	SO4	G	341	-	4,4,4	0.13	0	6,6,6	0.12	0
3	SO4	G	342	-	4,4,4	0.12	0	6,6,6	0.17	0
3	SO4	G	343	-	4,4,4	0.23	0	6,6,6	0.13	0
3	SO4	G	344	-	4,4,4	0.20	0	6,6,6	0.16	0
4	GOL	G	345	-	5,5,5	0.27	0	5,5,5	0.66	0
4	GOL	G	346	-	5,5,5	0.35	0	5,5,5	0.15	0
3	SO4	H	340	-	4,4,4	0.22	0	6,6,6	0.18	0
3	SO4	H	341	-	4,4,4	0.34	0	6,6,6	0.18	0
4	GOL	H	342	-	5,5,5	0.36	0	5,5,5	0.39	0
4	GOL	H	343	-	5,5,5	0.26	0	5,5,5	0.49	0
4	GOL	H	344	-	5,5,5	0.49	0	5,5,5	0.53	0
4	GOL	H	345	-	5,5,5	0.32	0	5,5,5	0.49	0
4	GOL	H	346	-	5,5,5	0.33	0	5,5,5	0.29	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
3	SO4	A	340	-	-	0/0/0/0	0/0/0/0
4	GOL	A	341	-	-	0/4/4/4	0/0/0/0
4	GOL	A	342	-	-	0/4/4/4	0/0/0/0
4	GOL	A	343	-	-	0/4/4/4	0/0/0/0
3	SO4	B	340	-	-	0/0/0/0	0/0/0/0
4	GOL	B	341	-	-	0/4/4/4	0/0/0/0
4	GOL	B	342	-	-	0/4/4/4	0/0/0/0
4	GOL	B	343	-	-	0/4/4/4	0/0/0/0
4	GOL	B	344	-	-	0/4/4/4	0/0/0/0
4	GOL	C	340	-	-	0/4/4/4	0/0/0/0
4	GOL	C	341	-	-	0/4/4/4	0/0/0/0
4	GOL	C	342	-	-	0/4/4/4	0/0/0/0
3	SO4	D	340	-	-	0/0/0/0	0/0/0/0
3	SO4	D	341	-	-	0/0/0/0	0/0/0/0
4	GOL	D	342	-	-	0/4/4/4	0/0/0/0
4	GOL	D	343	-	-	0/4/4/4	0/0/0/0
4	GOL	D	344	-	-	0/4/4/4	0/0/0/0
3	SO4	E	340	-	-	0/0/0/0	0/0/0/0
3	SO4	E	341	-	-	0/0/0/0	0/0/0/0
4	GOL	E	342	-	-	0/4/4/4	0/0/0/0
4	GOL	E	343	-	-	0/4/4/4	0/0/0/0
4	GOL	E	344	-	-	0/4/4/4	0/0/0/0
4	GOL	E	345	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SO4	F	340	-	-	0/0/0/0	0/0/0/0
3	SO4	F	341	-	-	0/0/0/0	0/0/0/0
4	GOL	F	342	-	-	0/4/4/4	0/0/0/0
3	SO4	G	340	-	-	0/0/0/0	0/0/0/0
3	SO4	G	341	-	-	0/0/0/0	0/0/0/0
3	SO4	G	342	-	-	0/0/0/0	0/0/0/0
3	SO4	G	343	-	-	0/0/0/0	0/0/0/0
3	SO4	G	344	-	-	0/0/0/0	0/0/0/0
4	GOL	G	345	-	-	0/4/4/4	0/0/0/0
4	GOL	G	346	-	-	0/4/4/4	0/0/0/0
3	SO4	H	340	-	-	0/0/0/0	0/0/0/0
3	SO4	H	341	-	-	0/0/0/0	0/0/0/0
4	GOL	H	342	-	-	0/4/4/4	0/0/0/0
4	GOL	H	343	-	-	0/4/4/4	0/0/0/0
4	GOL	H	344	-	-	0/4/4/4	0/0/0/0
4	GOL	H	345	-	-	0/4/4/4	0/0/0/0
4	GOL	H	346	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å <sup>2</sup> )	Q<0.9
1	A	339/339 (100%)	0.43	20 (5%)	22	20	8, 13, 20, 27	2 (0%)
1	B	339/339 (100%)	0.10	9 (2%)	52	48	8, 13, 21, 29	2 (0%)
1	C	339/339 (100%)	1.08	79 (23%)	1	1	8, 13, 20, 31	0
1	D	338/339 (99%)	0.62	28 (8%)	11	11	7, 12, 21, 30	4 (1%)
1	E	339/339 (100%)	0.23	2 (0%)	86	87	8, 12, 21, 29	1 (0%)
1	F	339/339 (100%)	0.24	6 (1%)	65	64	8, 12, 21, 34	4 (1%)
1	G	339/339 (100%)	0.27	5 (1%)	70	69	9, 12, 20, 28	3 (0%)
1	H	339/339 (100%)	0.23	4 (1%)	75	73	7, 12, 19, 30	2 (0%)
2	I	6/6 (100%)	0.36	0	100	100	11, 14, 20, 25	0
2	J	6/6 (100%)	0.31	0	100	100	7, 11, 21, 22	0
2	K	6/6 (100%)	0.43	0	100	100	14, 16, 20, 33	0
2	L	6/6 (100%)	0.31	0	100	100	7, 8, 16, 26	0
2	M	6/6 (100%)	0.53	0	100	100	7, 9, 15, 21	0
2	N	6/6 (100%)	0.40	0	100	100	6, 8, 16, 18	0
2	O	6/6 (100%)	0.59	0	100	100	7, 10, 17, 23	0
2	P	6/6 (100%)	0.48	0	100	100	8, 8, 18, 22	0
All	All	2759/2760 (99%)	0.40	153 (5%)	24	22	6, 12, 21, 34	18 (0%)

All (153) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	134	ARG	5.3
1	A	288	GLY	5.2
1	H	288	GLY	5.2
1	D	288	GLY	5.1
1	B	288	GLY	4.8

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Mol	Chain	Res	Type	RSRZ
1	C	208	GLY	4.8
1	G	288	GLY	4.5
1	C	136	THR	4.3
1	C	261	ASP	4.0
1	C	242	ALA	4.0
1	C	258	CYS	3.9
1	C	206	LEU	3.8
1	C	210	SER	3.8
1	A	271	GLY	3.8
1	C	288	GLY	3.8
1	C	262	THR	3.7
1	C	202	ALA	3.7
1	C	129	TYR	3.6
1	A	62	SER	3.6
1	C	213	PHE	3.6
1	C	330	TYR	3.5
1	E	288	GLY	3.5
1	C	201	LEU	3.5
1	C	209	SER	3.4
1	C	212	SER	3.4
1	D	271	GLY	3.4
1	D	134	ARG	3.4
1	C	132	SER	3.4
1	C	191	VAL	3.4
1	C	263	SER	3.3
1	A	70	LEU	3.3
1	A	65	SER	3.3
1	C	232	PHE	3.3
1	C	290	GLY	3.3
1	F	288	GLY	3.3
1	A	63	SER	3.3
1	C	214	GLY	3.2
1	A	259	ASN	3.2
1	A	64	SER	3.2
1	C	251[A]	GLN	3.2
1	C	215	ASP	3.2
1	H	271	GLY	3.1
1	C	211	PHE	3.1
1	A	68	LYS	3.1
1	D	154[A]	ARG	3.1
1	D	290	GLY	3.1
1	C	236	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	133	GLY	3.1
1	C	1	ASP	3.1
1	C	216	GLY	3.1
1	C	289	ASP	3.1
1	C	128	VAL	3.0
1	C	286	GLN	3.0
1	D	131	THR	3.0
1	C	259	ASN	3.0
1	A	67	TYR	3.0
1	F	258	CYS	3.0
1	D	194[A]	SER	3.0
1	C	272	ASN	2.9
1	C	200	SER	2.9
1	C	185	LYS	2.9
1	C	154[A]	ARG	2.8
1	B	208	GLY	2.8
1	B	290	GLY	2.8
1	G	71[A]	GLY	2.8
1	G	290	GLY	2.8
1	C	250[A]	ASP	2.8
1	C	155	THR	2.8
1	D	259	ASN	2.8
1	C	130	ASP	2.8
1	C	269	ASN	2.8
1	C	194[A]	SER	2.8
1	D	208	GLY	2.7
1	A	1	ASP	2.7
1	B	289	ASP	2.7
1	C	188	ALA	2.7
1	D	1	ASP	2.7
1	C	138	PRO	2.7
1	C	256	ILE	2.7
1	B	260[A]	THR	2.6
1	D	149	LYS	2.6
1	C	182	TYR	2.6
1	D	320	ALA	2.5
1	C	195	GLN	2.5
1	H	185	LYS	2.5
1	A	61	PRO	2.5
1	D	186	LEU	2.5
1	A	44[A]	ASN	2.5
1	D	65	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	22	SER	2.5
1	B	258	CYS	2.5
1	E	134	ARG	2.5
1	D	263	SER	2.5
1	C	122	TYR	2.4
1	C	183	SER	2.4
1	C	291	THR	2.4
1	C	285	TYR	2.4
1	F	71[A]	GLY	2.4
1	C	192	THR	2.4
1	C	260	THR	2.4
1	C	320	ALA	2.4
1	A	100	ILE	2.4
1	A	290	GLY	2.4
1	C	97	GLY	2.4
1	C	271	GLY	2.3
1	D	132	SER	2.3
1	C	241	GLY	2.3
1	D	291	THR	2.3
1	C	137	THR	2.3
1	D	260	THR	2.3
1	H	259	ASN	2.3
1	C	65	SER	2.3
1	C	280	ASN	2.3
1	F	259	ASN	2.3
1	C	331	THR	2.3
1	D	289	ASP	2.3
1	C	187	VAL	2.3
1	B	71	GLY	2.3
1	C	257	ASP	2.2
1	D	195	GLN	2.2
1	B	257	ASP	2.2
1	C	230	SER	2.2
1	C	190	GLN	2.2
1	D	188	ALA	2.2
1	F	247	VAL	2.2
1	G	46[A]	GLN	2.2
1	D	71	GLY	2.2
1	B	286	GLN	2.2
1	D	206	LEU	2.2
1	C	217	ALA	2.2
1	C	321	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	F	289	ASP	2.2
1	G	1	ASP	2.2
1	C	62	SER	2.1
1	C	131[A]	THR	2.1
1	C	68	LYS	2.1
1	A	214	GLY	2.1
1	C	264	GLY	2.1
1	A	185	LYS	2.1
1	D	113	VAL	2.1
1	D	185	LYS	2.1
1	C	203	SER	2.1
1	D	62	SER	2.0
1	D	258	CYS	2.0
1	A	286	GLN	2.0
1	C	207	LYS	2.0
1	C	339	VAL	2.0
1	C	277	THR	2.0
1	C	332[A]	THR	2.0
1	D	82[A]	SER	2.0
1	C	292	CYS	2.0
1	A	69	ASN	2.0
1	C	71	GLY	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	STA	L	506	12/12	0.18	3.08	14,21,40,40	0
2	STA	J	506	12/12	0.18	2.92	11,16,36,37	0
2	STA	P	506	12/12	0.20	2.64	11,17,35,37	0
2	STA	I	506	12/12	0.16	2.15	15,19,38,39	0
2	STA	K	506	12/12	0.18	2.07	24,29,42,44	0
2	STA	J	504	11/12	0.18	1.77	5,6,9,10	0
2	STA	K	504	11/12	0.16	1.70	12,14,16,16	0
2	STA	I	504	11/12	0.16	1.33	8,11,11,12	0
2	STA	O	504	11/12	0.19	1.21	5,6,8,8	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	STA	N	504	11/12	0.18	0.81	4,5,8,8	0
2	STA	O	506	12/12	0.17	0.80	11,17,37,38	0
2	STA	N	506	12/12	0.15	0.78	8,12,29,32	0
2	STA	L	504	11/12	0.18	0.64	6,7,10,10	0
2	STA	M	506	12/12	0.16	0.25	10,14,36,36	0
2	STA	P	504	11/12	0.16	0.04	5,7,8,9	0
2	STA	M	504	11/12	0.17	-0.07	5,7,9,9	0

## 6.3 Carbohydrates i

There are no carbohydrates in this entry.

## 6.4 Ligands i

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
4	GOL	H	346	6/6	0.25	11.71	53,54,55,57	0
4	GOL	G	346	6/6	0.23	5.87	46,47,48,48	0
4	GOL	G	345	6/6	0.19	5.45	26,34,36,40	0
3	SO4	E	340	5/5	0.14	4.95	46,47,49,49	0
3	SO4	F	341	5/5	0.22	4.22	14,17,18,19	5
4	GOL	B	343	6/6	0.19	3.77	24,27,29,29	0
3	SO4	D	341	5/5	0.15	3.42	51,52,53,53	0
4	GOL	D	343	6/6	0.27	2.22	35,43,45,49	0
4	GOL	F	342	6/6	0.20	2.11	34,42,44,45	0
3	SO4	F	340	5/5	0.13	1.54	60,61,61,62	0
3	SO4	G	344	5/5	0.20	1.40	73,74,75,75	0
3	SO4	G	341	5/5	0.19	1.37	76,76,76,76	0
4	GOL	D	342	6/6	0.18	1.25	35,37,38,40	0
4	GOL	E	342	6/6	0.16	1.21	36,38,40,40	0
4	GOL	B	342	6/6	0.16	1.04	40,42,43,44	0
4	GOL	H	342	6/6	0.17	1.00	46,47,49,49	0
4	GOL	A	343	6/6	0.18	0.92	59,59,59,59	0
4	GOL	A	341	6/6	0.15	0.85	38,40,41,41	0
4	GOL	C	340	6/6	0.30	0.83	54,57,57,58	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
4	GOL	B	344	6/6	0.17	0.71	30,37,38,39	0
4	GOL	E	344	6/6	0.15	0.60	19,23,24,26	0
4	GOL	E	343	6/6	0.17	0.60	36,37,38,38	0
4	GOL	H	344	6/6	0.13	0.58	26,27,30,30	0
3	SO4	H	341	5/5	0.18	0.53	14,16,17,20	5
3	SO4	D	340	5/5	0.20	0.43	91,91,92,92	0
4	GOL	E	345	6/6	0.11	0.36	18,27,27,28	0
3	SO4	G	342	5/5	0.12	0.36	75,75,75,76	0
4	GOL	B	341	6/6	0.15	0.28	42,44,45,45	0
3	SO4	E	341	5/5	0.16	0.26	65,65,65,66	0
3	SO4	G	340	5/5	0.12	0.18	43,43,45,45	0
4	GOL	D	344	6/6	0.17	0.13	60,61,62,62	0
4	GOL	H	343	6/6	0.12	0.11	32,37,38,40	0
3	SO4	G	343	5/5	0.10	0.07	60,60,60,61	0
4	GOL	C	341	6/6	0.18	0.01	46,46,46,46	0
4	GOL	A	342	6/6	0.14	-0.47	40,42,42,43	0
4	GOL	H	345	6/6	0.12	-0.48	39,40,40,41	0
3	SO4	A	340	5/5	0.13	-0.84	93,93,94,94	0
4	GOL	C	342	6/6	0.18	-1.00	56,56,56,56	0
3	SO4	H	340	5/5	0.13	-1.04	67,67,67,68	0
3	SO4	B	340	5/5	0.06	-1.17	49,51,51,51	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.