



Full wwPDB X-ray Structure Validation Report

Sep 9, 2014 – 05:14 PM EDT

PDB ID : 4MLG
Title : Structure of RS223-Beta-xylosidase
Authors : Jordan, D.; Braker, J.; Wagschal, K.; Lee, C.; Dubrovskaya, I.; Anderson, S.;
Wawrzak, Z.
Deposited on : 2013-09-06
Resolution : 2.70 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

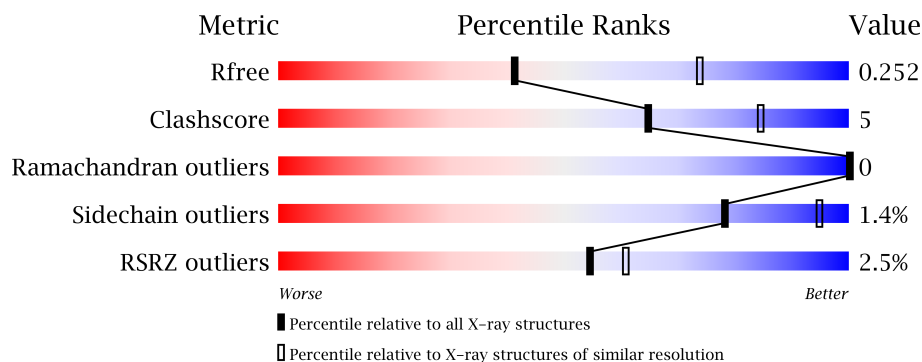
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1439
EDS : stable23489
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) : stable23489

1 Overall quality at a glance

The reported resolution of this entry is 2.70 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	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (2.70-2.70)

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. 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	324	
1	B	324	
1	C	324	
1	D	324	
1	E	324	
1	F	324	
1	G	324	
1	H	324	
1	I	324	
1	J	324	
1	K	324	
1	L	324	

The following table lists non-polymeric compounds that are outliers for geometric or electron-

density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	CA	A	401	-	X
2	CA	B	401	-	X
2	CA	C	401	-	X
2	CA	E	401	-	X
2	CA	F	401	-	X
2	CA	G	401	-	X
2	CA	H	401	-	X
2	CA	I	401	-	X
2	CA	K	401	-	X
3	SO4	A	402	-	X
3	SO4	A	403	-	X
3	SO4	C	403	-	X
3	SO4	D	402	-	X
3	SO4	D	403	-	X
3	SO4	E	402	-	X
3	SO4	F	402	-	X
3	SO4	F	403	-	X
3	SO4	G	402	-	X
3	SO4	G	403	-	X
3	SO4	H	403	-	X
3	SO4	I	402	-	X
3	SO4	I	403	-	X
3	SO4	J	403	-	X
3	SO4	K	402	-	X
3	SO4	K	403	-	X
3	SO4	L	403	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 32402 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 Beta-xylosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	323	Total	C	N	O	S	0	0	0
			2640	1699	438	488	15			
1	B	323	Total	C	N	O	S	0	0	0
			2640	1699	438	488	15			
1	C	323	Total	C	N	O	S	0	0	0
			2636	1696	437	488	15			
1	D	322	Total	C	N	O	S	0	0	0
			2635	1696	437	487	15			
1	E	323	Total	C	N	O	S	0	0	0
			2640	1699	438	488	15			
1	F	323	Total	C	N	O	S	0	0	0
			2636	1696	437	488	15			
1	G	324	Total	C	N	O	S	0	0	0
			2641	1699	438	489	15			
1	H	324	Total	C	N	O	S	0	0	0
			2645	1701	438	491	15			
1	I	321	Total	C	N	O	S	0	0	0
			2615	1684	434	483	14			
1	J	323	Total	C	N	O	S	0	0	0
			2636	1696	437	488	15			
1	K	324	Total	C	N	O	S	0	0	0
			2645	1701	438	491	15			
1	L	322	Total	C	N	O	S	0	0	0
			2631	1694	436	486	15			

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

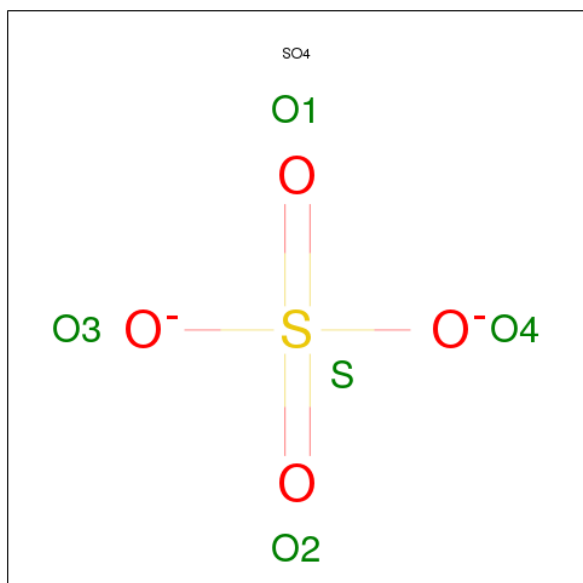
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	1	Total	Ca	0	0
			1	1		
2	J	1	Total	Ca	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total	Ca	0	0
			1	1		
2	K	1	Total	Ca	0	0
			1	1		
2	E	1	Total	Ca	0	0
			1	1		
2	H	1	Total	Ca	0	0
			1	1		
2	B	1	Total	Ca	0	0
			1	1		
2	I	1	Total	Ca	0	0
			1	1		
2	C	1	Total	Ca	0	0
			1	1		
2	A	1	Total	Ca	0	0
			1	1		
2	L	1	Total	Ca	0	0
			1	1		
2	F	1	Total	Ca	0	0
			1	1		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

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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	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	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		
3	H	1	Total	O	S	0	0
			5	4	1		
3	H	1	Total	O	S	0	0
			5	4	1		
3	I	1	Total	O	S	0	0
			5	4	1		
3	I	1	Total	O	S	0	0
			5	4	1		
3	J	1	Total	O	S	0	0
			5	4	1		
3	J	1	Total	O	S	0	0
			5	4	1		
3	K	1	Total	O	S	0	0
			5	4	1		
3	K	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 4 is water.

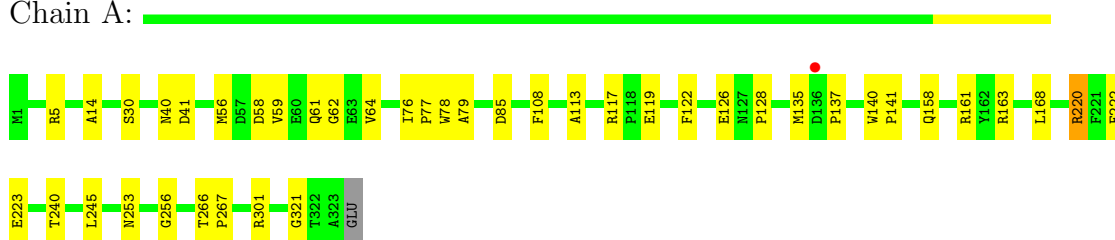
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	72	Total	O	0	0
			72	72		
4	B	64	Total	O	0	0
			64	64		
4	C	45	Total	O	0	0
			45	45		
4	D	68	Total	O	0	0
			68	68		
4	E	50	Total	O	0	0
			50	50		
4	F	58	Total	O	0	0
			58	58		
4	G	47	Total	O	0	0
			47	47		
4	H	50	Total	O	0	0
			50	50		
4	I	49	Total	O	0	0
			49	49		
4	J	50	Total	O	0	0
			50	50		
4	K	44	Total	O	0	0
			44	44		
4	L	33	Total	O	0	0
			33	33		

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.

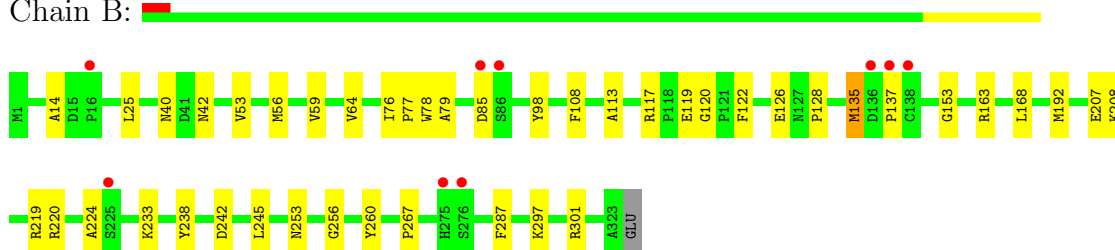
- Molecule 1: Beta-xylosidase

Chain A:



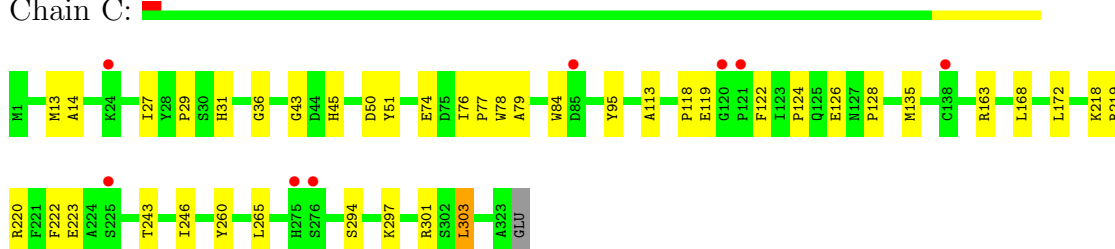
- Molecule 1: Beta-xylosidase

Chain B:



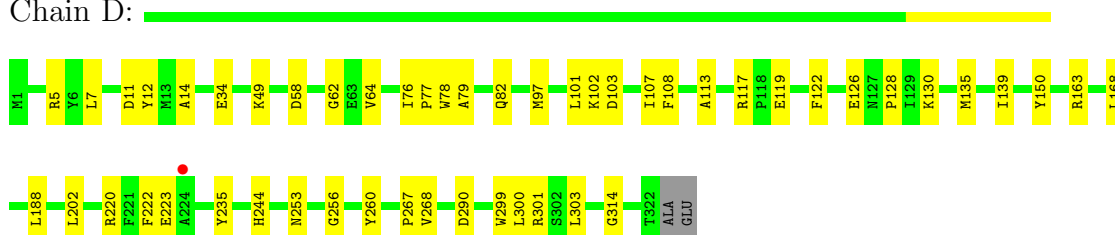
- Molecule 1: Beta-xylosidase

Chain C:



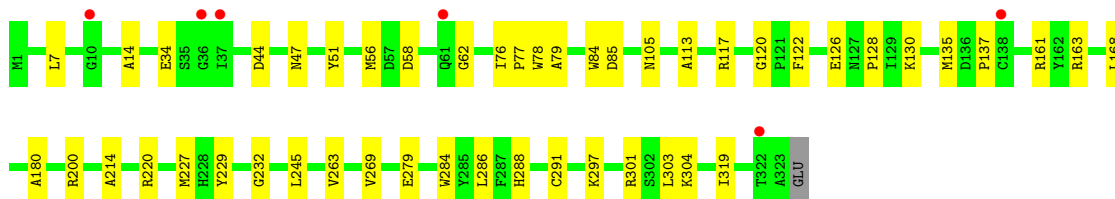
- Molecule 1: Beta-xylosidase

Chain D:



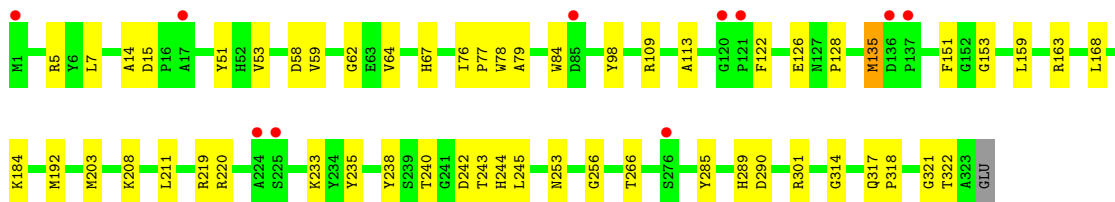
- Molecule 1: Beta-xylosidase

Chain E:



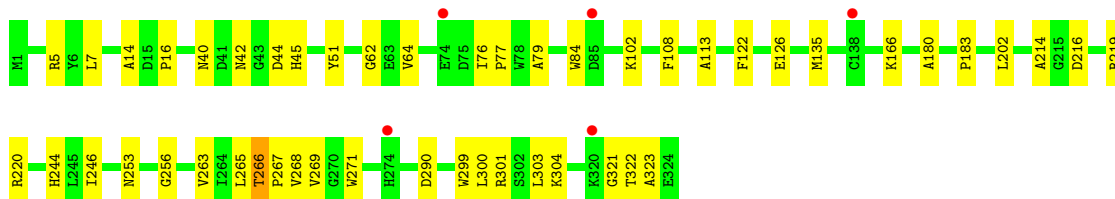
- Molecule 1: Beta-xylosidase

Chain F:



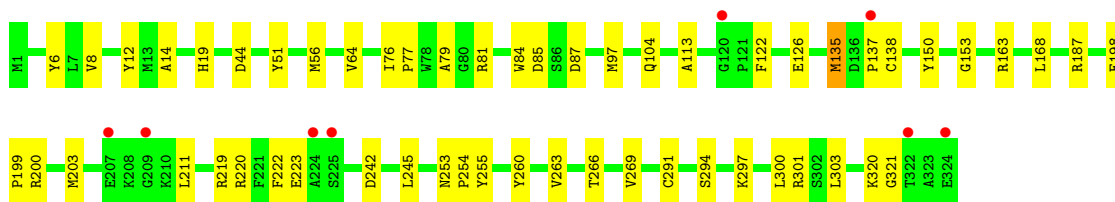
- Molecule 1: Beta-xylosidase

Chain G:



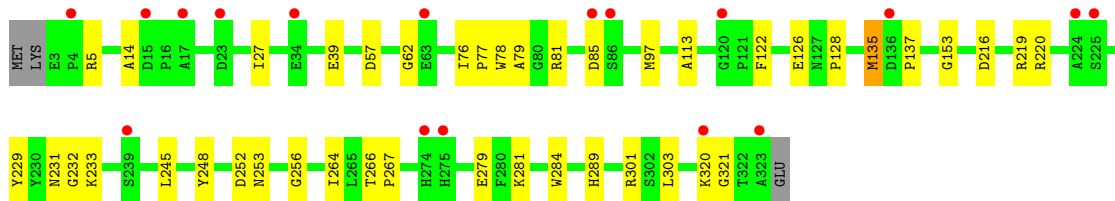
- Molecule 1: Beta-xylosidase

Chain H:

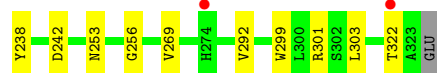


- Molecule 1: Beta-xylosidase

Chain I:



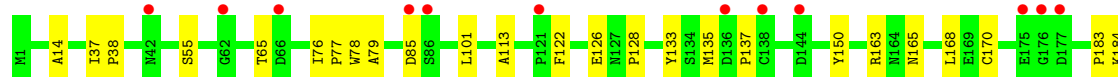
- Molecule 1: Beta-xylosidase

Chain J: 

- Molecule 1: Beta-xylosidase

Chain K: 

- Molecule 1: Beta-xylosidase

Chain L: 

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	124.53Å 148.70Å 125.65Å 90.00° 112.87° 90.00°	Depositor
Resolution (Å)	29.20 – 2.70 29.17 – 2.69	Depositor EDS
% Data completeness (in resolution range)	98.3 (29.20-2.70) 91.8 (29.17-2.69)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 2.68Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.210 , 0.257 0.205 , 0.252	Depositor DCC
R_{free} test set	5339 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	40.3	Xtriage
Anisotropy	0.594	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 37.4	EDS
Estimated twinning fraction	0.018 for l,-k,h	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 115104 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	32402	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.78% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities 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: CA, 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	0.22	0/2729	0.39	0/3706
1	B	0.22	0/2729	0.38	0/3706
1	C	0.21	0/2725	0.38	0/3702
1	D	0.22	0/2724	0.38	0/3699
1	E	0.22	0/2729	0.38	0/3706
1	F	0.23	0/2725	0.38	0/3702
1	G	0.21	0/2730	0.37	0/3709
1	H	0.22	0/2734	0.39	0/3714
1	I	0.21	0/2704	0.37	0/3675
1	J	0.22	0/2725	0.38	0/3702
1	K	0.22	0/2734	0.38	0/3714
1	L	0.21	0/2720	0.37	0/3694
All	All	0.22	0/32708	0.38	0/44429

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2640	0	2490	20	0
1	B	2640	0	2490	21	0
1	C	2636	0	2479	25	0
1	D	2635	0	2485	26	0
1	E	2640	0	2490	28	0
1	F	2636	0	2479	41	0
1	G	2641	0	2481	34	0
1	H	2645	0	2485	30	0
1	I	2615	0	2455	26	0
1	J	2636	0	2479	27	0
1	K	2645	0	2485	30	0
1	L	2631	0	2479	27	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
3	A	10	0	0	0	0
3	B	10	0	0	0	0
3	C	10	0	0	0	0
3	D	10	0	0	0	0
3	E	10	0	0	0	0
3	F	10	0	0	0	0
3	G	10	0	0	0	0
3	H	10	0	0	0	0
3	I	10	0	0	0	0
3	J	10	0	0	0	0
3	K	10	0	0	0	0
3	L	10	0	0	0	0
4	A	72	0	0	0	0
4	B	64	0	0	0	0
4	C	45	0	0	0	0
4	D	68	0	0	0	0
4	E	50	0	0	2	0
4	F	58	0	0	0	0
4	G	47	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	50	0	0	2	0
4	I	49	0	0	1	0
4	J	50	0	0	1	0
4	K	44	0	0	1	0
4	L	33	0	0	0	0
All	All	32402	0	29777	328	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 5.

All (328) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:265:LEU:HD12	1:G:266:THR:H	1.21	1.05
1:F:240:THR:CG2	1:F:243:THR:HB	2.13	0.79
1:G:265:LEU:HD12	1:G:266:THR:N	1.98	0.78
1:F:240:THR:HG22	1:F:243:THR:HB	1.69	0.74
1:L:266:THR:HG23	1:L:267:PRO:HD2	1.70	0.72
1:H:104:GLN:HG3	4:H:507:HOH:O	1.90	0.71
1:J:118:PRO:HG2	1:J:119:GLU:OE1	1.89	0.71
1:F:243:THR:HG22	1:F:245:LEU:HG	1.75	0.67
1:A:220:ARG:NH2	1:A:240:THR:O	2.28	0.67
1:G:290:ASP:HB2	1:G:304:LYS:HD2	1.75	0.66
1:G:76:ILE:HB	1:G:79:ALA:HB2	1.76	0.66
1:E:77:PRO:HG2	1:E:126:GLU:HA	1.78	0.66
1:C:76:ILE:HB	1:C:79:ALA:HB2	1.78	0.66
1:K:76:ILE:HB	1:K:79:ALA:HB2	1.78	0.66
1:J:76:ILE:HB	1:J:79:ALA:HB2	1.78	0.65
1:H:203:MET:HB2	1:H:211:LEU:HD21	1.79	0.64
1:E:269:VAL:HG23	1:E:291:CYS:H	1.62	0.64
1:I:77:PRO:HG2	1:I:126:GLU:HA	1.80	0.64
1:E:44:ASP:O	1:E:47:ASN:ND2	2.32	0.62
1:H:76:ILE:HB	1:H:79:ALA:HB2	1.80	0.62
1:F:76:ILE:HB	1:F:79:ALA:HB2	1.81	0.62
1:B:253:ASN:HD21	1:B:256:GLY:HA3	1.65	0.62
1:E:117:ARG:HG3	1:E:120:GLY:H	1.66	0.61
1:I:57:ASP:O	1:I:281:LYS:NZ	2.33	0.61
1:G:246:ILE:HB	1:G:265:LEU:HB3	1.82	0.61
1:L:266:THR:CG2	1:L:267:PRO:HD2	2.30	0.61
1:E:14:ALA:HB2	1:E:301:ARG:HB2	1.83	0.61
1:H:320:LYS:NZ	4:H:535:HOH:O	2.34	0.61
1:L:77:PRO:HG2	1:L:126:GLU:HA	1.81	0.61
1:L:76:ILE:HB	1:L:79:ALA:HB2	1.83	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:K:14:ALA:HB2	1:K:301:ARG:HB2	1.83	0.60
1:K:294:SER:HB3	1:K:297:LYS:HB2	1.84	0.60
1:I:113:ALA:HB1	1:I:122:PHE:HB3	1.82	0.60
1:K:269:VAL:HG23	1:K:322:THR:HB	1.85	0.59
1:B:14:ALA:HB2	1:B:301:ARG:HB2	1.85	0.59
1:F:240:THR:HG22	1:F:240:THR:O	2.03	0.59
1:A:76:ILE:HB	1:A:79:ALA:HB2	1.84	0.59
1:J:52:HIS:NE2	4:J:510:HOH:O	2.32	0.59
1:F:113:ALA:HB1	1:F:122:PHE:HB3	1.85	0.59
1:C:95:TYR:CZ	1:C:118:PRO:HG3	2.38	0.58
1:A:14:ALA:HB2	1:A:301:ARG:HB2	1.85	0.58
1:C:51:TYR:OH	1:C:84:TRP:O	2.20	0.58
1:J:117:ARG:HB3	1:J:119:GLU:OE2	2.03	0.58
1:E:7:LEU:HD13	1:E:56:MET:CE	2.34	0.58
1:C:77:PRO:HG2	1:C:126:GLU:HA	1.86	0.58
1:F:14:ALA:HB2	1:F:301:ARG:HB2	1.86	0.58
1:J:77:PRO:HG2	1:J:126:GLU:HA	1.86	0.58
1:H:297:LYS:HD2	1:H:300:LEU:HD12	1.86	0.58
1:D:7:LEU:HB3	1:D:64:VAL:HG21	1.86	0.57
1:B:77:PRO:HG2	1:B:126:GLU:HA	1.86	0.57
1:C:113:ALA:HB1	1:C:122:PHE:HB3	1.87	0.57
1:F:243:THR:HG21	1:F:245:LEU:HD12	1.86	0.56
1:D:103:ASP:OD1	1:D:107:ILE:N	2.33	0.56
1:G:113:ALA:HB1	1:G:122:PHE:HB3	1.87	0.56
1:L:113:ALA:HB1	1:L:122:PHE:HB3	1.87	0.56
1:E:7:LEU:HD12	1:E:303:LEU:CD1	2.35	0.56
1:L:37:ILE:HD12	1:L:38:PRO:HD2	1.88	0.56
1:F:266:THR:OG1	1:F:321:GLY:N	2.37	0.56
1:F:5:ARG:NH1	1:F:59:VAL:O	2.38	0.56
1:J:14:ALA:HB2	1:J:301:ARG:HB2	1.88	0.55
1:B:207:GLU:HG2	1:B:208:LYS:HD2	1.88	0.55
1:C:14:ALA:HB2	1:C:301:ARG:HB2	1.88	0.55
1:K:220:ARG:NH2	1:K:240:THR:O	2.38	0.55
1:L:290:ASP:HB2	1:L:304:LYS:HD2	1.88	0.55
1:L:133:TYR:OH	1:L:165:ASN:ND2	2.39	0.55
1:F:243:THR:CG2	1:F:245:LEU:HG	2.36	0.55
1:I:81:ARG:NH1	4:I:508:HOH:O	2.39	0.55
1:L:14:ALA:HB2	1:L:301:ARG:HB2	1.89	0.54
1:D:253:ASN:ND2	1:D:256:GLY:H	2.07	0.53
1:I:248:TYR:HB3	1:I:264:ILE:HD11	1.89	0.53
1:D:76:ILE:HB	1:D:79:ALA:HB2	1.90	0.53
1:E:163:ARG:HG2	1:E:168:LEU:HD11	1.89	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:5:ARG:NH2	1:D:62:GLY:O	2.36	0.53
1:G:102:LYS:NZ	4:G:513:HOH:O	2.41	0.53
1:F:203:MET:HB3	1:F:211:LEU:HD21	1.91	0.53
1:F:242:ASP:OD1	1:F:242:ASP:N	2.42	0.53
1:I:97:MET:HB3	1:I:113:ALA:HB3	1.91	0.53
1:I:289:HIS:CE1	1:I:301:ARG:HD2	2.44	0.53
1:J:253:ASN:OD1	1:J:256:GLY:N	2.40	0.52
1:K:252:ASP:OD1	1:K:253:ASN:ND2	2.37	0.52
1:E:7:LEU:HD13	1:E:56:MET:HE2	1.90	0.52
1:A:253:ASN:HD21	1:A:256:GLY:HA3	1.74	0.52
1:H:14:ALA:HB2	1:H:301:ARG:HB2	1.92	0.52
1:H:8:VAL:HG23	1:H:303:LEU:HD11	1.92	0.52
1:K:113:ALA:HB1	1:K:122:PHE:HB3	1.91	0.52
1:G:263:VAL:HG13	1:G:265:LEU:O	2.10	0.52
1:E:200:ARG:NH2	4:E:530:HOH:O	2.43	0.52
1:B:219:ARG:HG2	1:B:260:TYR:CZ	2.45	0.51
1:F:15:ASP:H	1:F:289:HIS:CD2	2.28	0.51
1:H:113:ALA:HB1	1:H:122:PHE:HB3	1.92	0.51
1:I:5:ARG:NH2	1:I:62:GLY:O	2.29	0.51
1:J:113:ALA:HB1	1:J:122:PHE:HB3	1.92	0.51
1:F:242:ASP:O	1:F:244:HIS:NE2	2.43	0.51
1:H:51:TYR:OH	1:H:84:TRP:O	2.21	0.51
1:J:51:TYR:OH	1:J:84:TRP:O	2.23	0.51
1:G:268:VAL:HG13	1:G:290:ASP:OD1	2.11	0.51
1:I:14:ALA:HB2	1:I:301:ARG:HB2	1.92	0.51
1:K:1:MET:HG2	1:K:1:MET:O	2.11	0.50
1:L:55:SER:HB3	1:L:65:THR:HB	1.93	0.50
1:D:113:ALA:HB1	1:D:122:PHE:HB3	1.94	0.50
1:G:77:PRO:HG2	1:G:126:GLU:HA	1.93	0.50
1:F:289:HIS:CE1	1:F:301:ARG:HD2	2.47	0.50
1:E:76:ILE:HB	1:E:79:ALA:HB2	1.93	0.50
1:G:267:PRO:HB2	1:G:323:ALA:O	2.12	0.50
1:L:273:THR:O	1:L:273:THR:HG23	2.11	0.50
1:B:53:VAL:HG11	1:B:119:GLU:HA	1.94	0.50
1:K:77:PRO:HG2	1:K:126:GLU:HA	1.93	0.50
1:A:77:PRO:HG2	1:A:126:GLU:HA	1.94	0.49
1:J:8:VAL:HG23	1:J:303:LEU:HD11	1.94	0.49
1:K:180:ALA:HB2	1:K:214:ALA:HA	1.95	0.49
1:D:97:MET:HB3	1:D:113:ALA:HB3	1.94	0.49
1:D:163:ARG:HG2	1:D:168:LEU:HD11	1.94	0.49
1:I:252:ASP:OD1	1:I:253:ASN:ND2	2.37	0.49
1:K:253:ASN:OD1	1:K:256:GLY:N	2.42	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:82:GLN:OE1	1:D:102:LYS:HD2	2.12	0.49
1:C:74:GLU:N	1:C:74:GLU:OE2	2.46	0.49
1:H:242:ASP:N	1:H:242:ASP:OD1	2.43	0.49
1:E:279:GLU:HB2	1:E:284:TRP:CE2	2.48	0.49
1:J:78:TRP:CG	1:J:128:PRO:HD3	2.48	0.49
1:L:183:PRO:HG2	1:L:202:LEU:HB3	1.95	0.49
1:D:14:ALA:HB2	1:D:301:ARG:HB2	1.95	0.48
1:E:34:GLU:OE1	1:E:297:LYS:NZ	2.29	0.48
1:H:77:PRO:HG2	1:H:126:GLU:HA	1.94	0.48
1:I:216:ASP:OD2	1:I:219:ARG:NH1	2.45	0.48
1:F:135:MET:SD	1:F:153:GLY:HA3	2.54	0.48
1:I:253:ASN:HD21	1:I:256:GLY:HA3	1.79	0.48
1:A:56:MET:HE3	1:A:64:VAL:HG22	1.95	0.48
1:F:77:PRO:HG2	1:F:126:GLU:HA	1.95	0.48
1:C:119:GLU:N	1:C:119:GLU:OE1	2.43	0.48
1:C:27:ILE:HG22	1:C:29:PRO:HD3	1.95	0.48
1:J:269:VAL:HG23	1:J:322:THR:HB	1.96	0.48
1:L:299:TRP:CD1	1:L:300:LEU:HG	2.48	0.48
1:A:117:ARG:NH1	1:A:119:GLU:OE2	2.46	0.48
1:K:58:ASP:HB3	1:K:62:GLY:HA3	1.95	0.48
1:L:225:SER:HG	1:L:236:PHE:HE1	1.62	0.48
1:G:253:ASN:ND2	1:G:256:GLY:H	2.12	0.48
1:H:163:ARG:HG2	1:H:168:LEU:HD11	1.95	0.48
1:K:248:TYR:HB3	1:K:264:ILE:HD11	1.96	0.48
1:K:170:CYS:HA	1:L:170:CYS:HA	1.96	0.48
1:F:98:TYR:OH	1:F:192:MET:HB3	2.13	0.47
1:K:245:LEU:HD23	1:K:267:PRO:HD3	1.95	0.47
1:C:36:GLY:HA3	1:H:6:TYR:O	2.14	0.47
1:G:14:ALA:HB2	1:G:301:ARG:HB2	1.95	0.47
1:E:113:ALA:HB1	1:E:122:PHE:HB3	1.96	0.47
1:F:163:ARG:HG2	1:F:168:LEU:HD11	1.97	0.47
1:F:253:ASN:HD21	1:F:256:GLY:HA3	1.79	0.47
1:J:85:ASP:O	1:J:137:PRO:HD2	2.15	0.47
1:B:56:MET:SD	1:B:64:VAL:HG22	2.53	0.47
1:D:117:ARG:NH2	1:D:119:GLU:OE2	2.47	0.47
1:J:45:HIS:HD2	1:J:299:TRP:HB3	1.79	0.47
1:L:85:ASP:O	1:L:137:PRO:HD2	2.15	0.47
1:C:163:ARG:HG2	1:C:168:LEU:HD11	1.96	0.47
1:F:153:GLY:O	1:F:159:LEU:HB2	2.14	0.47
1:A:5:ARG:NH1	1:A:59:VAL:O	2.46	0.47
1:G:265:LEU:CD1	1:G:266:THR:H	2.08	0.47
1:G:303:LEU:N	4:G:508:HOH:O	2.37	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:266:THR:HG1	1:I:321:GLY:H	1.59	0.47
1:A:113:ALA:HB1	1:A:122:PHE:HB3	1.95	0.47
1:F:290:ASP:OD2	1:F:322:THR:HG21	2.15	0.47
1:G:263:VAL:O	1:G:263:VAL:HG12	2.13	0.47
1:K:233:LYS:HD2	1:K:250:THR:HG23	1.97	0.46
1:F:233:LYS:HE2	1:F:314:GLY:HA3	1.98	0.46
1:G:265:LEU:CD1	1:G:266:THR:N	2.73	0.46
1:G:42:ASN:ND2	1:G:44:ASP:OD2	2.48	0.46
1:H:138:CYS:SG	1:H:150:TYR:HB2	2.56	0.46
1:I:39:GLU:O	1:I:39:GLU:HG3	2.15	0.46
1:D:34:GLU:OE1	1:D:299:TRP:NE1	2.34	0.46
1:G:40:ASN:H	1:G:45:HIS:CD2	2.33	0.46
1:H:97:MET:HB3	1:H:113:ALA:HB3	1.98	0.46
1:K:85:ASP:O	1:K:137:PRO:HD2	2.15	0.46
1:A:266:THR:HG1	1:A:321:GLY:H	1.57	0.46
1:G:266:THR:OG1	1:G:321:GLY:N	2.48	0.46
1:J:117:ARG:HB2	1:J:120:GLY:H	1.81	0.46
1:B:245:LEU:HD23	1:B:267:PRO:HD3	1.98	0.46
1:C:246:ILE:HB	1:C:265:LEU:HB3	1.98	0.46
1:G:266:THR:HB	1:G:267:PRO:CD	2.46	0.46
1:A:14:ALA:O	1:A:30:SER:N	2.45	0.46
1:C:31:HIS:ND1	1:C:50:ASP:OD1	2.46	0.46
1:E:58:ASP:HB3	1:E:62:GLY:HA3	1.98	0.46
1:C:218:LYS:HG2	1:C:243:THR:HG21	1.98	0.45
1:E:105:ASN:OD1	1:F:109:ARG:NH1	2.46	0.45
1:K:135:MET:SD	1:K:153:GLY:HA3	2.56	0.45
1:G:5:ARG:NH2	1:G:62:GLY:O	2.39	0.45
1:L:246:ILE:HB	1:L:265:LEU:HB3	1.98	0.45
1:H:187:ARG:HB2	1:H:255:TYR:OH	2.16	0.45
1:H:222:PHE:HD2	1:H:223:GLU:HG2	1.82	0.45
1:F:7:LEU:HB3	1:F:64:VAL:HG21	1.98	0.45
1:L:222:PHE:HD2	1:L:223:GLU:HG2	1.82	0.45
1:B:25:LEU:HD22	1:B:59:VAL:HG22	1.97	0.45
1:F:53:VAL:HG13	1:F:67:HIS:HB2	1.98	0.45
1:A:85:ASP:O	1:A:137:PRO:HD2	2.16	0.45
1:C:222:PHE:HD2	1:C:223:GLU:HG2	1.82	0.45
1:F:242:ASP:O	1:F:244:HIS:CD2	2.70	0.45
1:A:245:LEU:HD23	1:A:267:PRO:HD3	1.99	0.44
1:C:219:ARG:HG2	1:C:260:TYR:CZ	2.52	0.44
1:C:78:TRP:CG	1:C:128:PRO:HD3	2.53	0.44
1:F:51:TYR:OH	1:F:84:TRP:O	2.28	0.44
1:D:222:PHE:HD2	1:D:223:GLU:HG2	1.82	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:78:TRP:CG	1:F:128:PRO:HD3	2.52	0.44
1:H:266:THR:HG1	1:H:321:GLY:H	1.61	0.44
1:G:45:HIS:CD2	1:G:45:HIS:H	2.34	0.44
1:D:268:VAL:HB	1:D:290:ASP:HB3	1.99	0.44
1:F:151:PHE:CE1	1:F:184:LYS:HB2	2.53	0.44
1:G:51:TYR:OH	1:G:84:TRP:O	2.25	0.44
1:L:242:ASP:OD1	1:L:242:ASP:N	2.48	0.44
1:H:19:HIS:NE2	1:H:87:ASP:OD1	2.40	0.44
1:F:317:GLN:HA	1:F:318:PRO:HD3	1.89	0.44
1:H:135:MET:SD	1:H:153:GLY:HA3	2.58	0.44
1:J:253:ASN:HD21	1:J:256:GLY:HA3	1.82	0.44
1:D:260:TYR:O	1:E:117:ARG:NH1	2.44	0.44
1:K:5:ARG:NH1	1:K:59:VAL:O	2.43	0.44
1:C:95:TYR:CE1	1:C:118:PRO:HG3	2.53	0.43
1:H:56:MET:HE3	1:H:64:VAL:HG22	1.99	0.43
1:I:135:MET:SD	1:I:153:GLY:HA3	2.58	0.43
1:H:219:ARG:HG2	1:H:260:TYR:CZ	2.53	0.43
1:J:227:MET:HE1	1:J:236:PHE:HB2	2.00	0.43
1:J:39:GLU:HG3	1:J:45:HIS:CE1	2.53	0.43
1:B:40:ASN:ND2	1:B:42:ASN:OD1	2.50	0.43
1:J:219:ARG:O	1:J:238:TYR:OH	2.30	0.43
1:L:184:LYS:NZ	1:L:201:ASP:OD1	2.35	0.43
1:F:151:PHE:CZ	1:F:184:LYS:HB2	2.53	0.43
1:G:244:HIS:ND1	1:G:271:TRP:O	2.44	0.43
1:B:113:ALA:HB1	1:B:122:PHE:HB3	1.99	0.43
1:C:78:TRP:HA	1:D:130:LYS:HD3	2.00	0.43
1:H:44:ASP:OD1	1:H:81:ARG:NH1	2.40	0.43
1:J:292:VAL:HG21	1:J:322:THR:HG21	1.99	0.43
1:D:139:ILE:HD12	1:D:188:LEU:HD11	2.00	0.43
1:G:7:LEU:HB3	1:G:64:VAL:HG21	2.01	0.43
1:I:245:LEU:HD23	1:I:267:PRO:HD3	2.00	0.43
1:K:130:LYS:NZ	4:K:511:HOH:O	2.45	0.43
1:J:117:ARG:NE	1:J:119:GLU:OE2	2.52	0.43
1:D:103:ASP:OD1	1:D:107:ILE:O	2.37	0.43
1:G:269:VAL:H	1:G:322:THR:HB	1.84	0.43
1:J:142:ASP:HB3	1:J:144:ASP:OD1	2.19	0.43
1:L:79:ALA:HA	1:L:101:LEU:HD21	2.01	0.43
1:B:78:TRP:CG	1:B:128:PRO:HD3	2.54	0.43
1:E:78:TRP:CG	1:E:128:PRO:HD3	2.54	0.43
1:J:119:GLU:H	1:J:119:GLU:CD	2.22	0.43
1:C:43:GLY:HA2	1:C:45:HIS:CE1	2.53	0.42
1:E:161:ARG:NH2	4:E:533:HOH:O	2.52	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:77:PRO:HG2	1:D:126:GLU:HA	2.00	0.42
1:F:240:THR:HG21	1:F:243:THR:HB	1.99	0.42
1:E:85:ASP:O	1:E:137:PRO:HD2	2.20	0.42
1:G:266:THR:HB	1:G:267:PRO:HD2	2.01	0.42
1:H:253:ASN:HB2	1:H:254:PRO:HD2	2.01	0.42
1:J:119:GLU:CD	1:J:119:GLU:N	2.73	0.42
1:A:58:ASP:O	1:A:62:GLY:HA3	2.19	0.42
1:B:242:ASP:N	1:B:242:ASP:OD1	2.51	0.42
1:C:294:SER:HB3	1:C:297:LYS:HB2	2.01	0.42
1:D:58:ASP:HB3	1:D:62:GLY:HA3	2.01	0.42
1:G:180:ALA:HB2	1:G:214:ALA:HA	2.01	0.42
1:B:135:MET:SD	1:B:153:GLY:HA3	2.59	0.42
1:D:78:TRP:CG	1:D:128:PRO:HD3	2.55	0.42
1:G:216:ASP:OD2	1:G:219:ARG:HD3	2.18	0.42
1:H:269:VAL:O	1:H:291:CYS:N	2.49	0.42
1:K:113:ALA:HA	1:K:124:PRO:HA	2.02	0.42
1:K:222:PHE:HD2	1:K:223:GLU:HG2	1.84	0.42
1:K:44:ASP:OD1	1:K:81:ARG:NH1	2.52	0.42
1:D:78:TRP:CH2	1:D:101:LEU:HB3	2.54	0.42
1:E:286:LEU:HD22	1:E:319:ILE:HD12	2.02	0.42
1:G:166:LYS:NZ	1:H:104:GLN:O	2.36	0.42
1:I:233:LYS:HB2	1:I:233:LYS:HE3	1.90	0.42
1:I:85:ASP:O	1:I:137:PRO:HD2	2.19	0.42
1:E:245:LEU:HD22	1:E:263:VAL:HG11	2.01	0.42
1:E:51:TYR:OH	1:E:84:TRP:O	2.22	0.42
1:A:158:GLN:HB3	1:A:161:ARG:HD2	2.01	0.42
1:E:130:LYS:HD3	1:F:78:TRP:HA	2.01	0.42
1:J:163:ARG:HG2	1:J:168:LEU:HD11	2.02	0.42
1:J:242:ASP:N	1:J:242:ASP:OD1	2.52	0.42
1:J:76:ILE:HA	1:J:77:PRO:HD3	1.93	0.42
1:A:78:TRP:CG	1:A:128:PRO:HD3	2.54	0.42
1:A:163:ARG:HG2	1:A:168:LEU:HD11	2.01	0.42
1:B:98:TYR:OH	1:B:192:MET:HB3	2.20	0.42
1:D:12:TYR:CE1	1:D:300:LEU:HB3	2.55	0.42
1:G:16:PRO:HB3	1:G:303:LEU:HD13	2.01	0.42
1:I:5:ARG:HH12	1:I:62:GLY:H	1.66	0.42
1:G:299:TRP:CD1	1:G:300:LEU:HG	2.55	0.41
1:I:27:ILE:HG21	1:I:303:LEU:HD21	2.02	0.41
1:I:78:TRP:CG	1:I:128:PRO:HD3	2.55	0.41
1:C:76:ILE:HA	1:C:77:PRO:HD3	1.92	0.41
1:F:219:ARG:O	1:F:238:TYR:OH	2.32	0.41
1:E:229:TYR:OH	1:E:232:GLY:HA2	2.20	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:59:VAL:HG11	1:F:285:TYR:CZ	2.55	0.41
1:L:78:TRP:CG	1:L:128:PRO:HD3	2.55	0.41
1:E:288:HIS:NE2	1:E:304:LYS:HD2	2.35	0.41
1:L:165:ASN:ND2	1:L:199:PRO:HB2	2.35	0.41
1:L:150:TYR:CD1	1:L:202:LEU:HD22	2.55	0.41
1:B:76:ILE:HB	1:B:79:ALA:HB2	2.02	0.41
1:D:244:HIS:HB2	1:D:267:PRO:HB3	2.01	0.41
1:D:235:TYR:OH	1:D:314:GLY:O	2.31	0.41
1:I:320:LYS:HB2	1:I:320:LYS:HE3	1.86	0.41
1:A:40:ASN:OD1	1:A:41:ASP:N	2.54	0.41
1:H:85:ASP:O	1:H:137:PRO:HD2	2.20	0.41
1:K:98:TYR:OH	1:K:192:MET:HB3	2.21	0.41
1:C:163:ARG:NH2	1:C:172:LEU:O	2.44	0.41
1:F:208:LYS:HE2	1:F:208:LYS:HB3	1.91	0.41
1:K:78:TRP:CG	1:K:128:PRO:HD3	2.56	0.41
1:H:245:LEU:HD22	1:H:263:VAL:HG11	2.03	0.41
1:F:233:LYS:NZ	1:F:235:TYR:OH	2.29	0.41
1:K:227:MET:HB2	1:K:227:MET:HE2	1.94	0.41
1:L:309:LYS:HE2	1:L:309:LYS:HB3	1.87	0.41
1:A:140:TRP:HA	1:A:141:PRO:HD3	1.93	0.41
1:D:150:TYR:CD1	1:D:202:LEU:HD22	2.56	0.41
1:E:227:MET:HB2	1:E:227:MET:HE2	1.95	0.41
1:G:183:PRO:HG2	1:G:202:LEU:HB3	2.03	0.41
1:B:224:ALA:O	1:B:238:TYR:HA	2.21	0.40
1:C:113:ALA:HA	1:C:124:PRO:HA	2.03	0.40
1:E:180:ALA:HB2	1:E:214:ALA:HA	2.03	0.40
1:F:58:ASP:HB3	1:F:62:GLY:HA3	2.04	0.40
1:H:12:TYR:HE1	1:H:294:SER:HB3	1.86	0.40
1:I:76:ILE:HB	1:I:79:ALA:HB2	2.02	0.40
1:A:222:PHE:HD2	1:A:223:GLU:HG2	1.86	0.40
1:B:117:ARG:HG3	1:B:120:GLY:H	1.87	0.40
1:H:198:GLU:HA	1:H:199:PRO:HD3	1.95	0.40
1:I:76:ILE:HA	1:I:77:PRO:HD3	1.93	0.40
1:K:289:HIS:NE2	1:K:301:ARG:HD2	2.36	0.40
1:K:51:TYR:OH	1:K:84:TRP:O	2.25	0.40
1:L:163:ARG:HG2	1:L:168:LEU:HD11	2.02	0.40
1:B:85:ASP:O	1:B:137:PRO:HD2	2.21	0.40
1:B:163:ARG:HG2	1:B:168:LEU:HD11	2.03	0.40
1:C:13:MET:HB2	1:C:303:LEU:HG	2.03	0.40
1:I:279:GLU:HB2	1:I:284:TRP:CE2	2.57	0.40
1:K:37:ILE:HG21	1:K:44:ASP:HB3	2.04	0.40
1:K:87:ASP:OD1	1:K:88:VAL:N	2.52	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:15:ASP:H	1:F:289:HIS:HD2	1.68	0.40
1:I:229:TYR:OH	1:I:232:GLY:HA2	2.20	0.40
1:B:25:LEU:HD21	1:B:287:PHE:CZ	2.56	0.40
1:L:198:GLU:HA	1:L:199:PRO:HD3	1.91	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	321/324 (99%)	302 (94%)	19 (6%)	0	100	100
1	B	321/324 (99%)	303 (94%)	18 (6%)	0	100	100
1	C	321/324 (99%)	303 (94%)	18 (6%)	0	100	100
1	D	320/324 (99%)	304 (95%)	16 (5%)	0	100	100
1	E	321/324 (99%)	304 (95%)	17 (5%)	0	100	100
1	F	321/324 (99%)	300 (94%)	21 (6%)	0	100	100
1	G	322/324 (99%)	298 (92%)	24 (8%)	0	100	100
1	H	322/324 (99%)	304 (94%)	18 (6%)	0	100	100
1	I	319/324 (98%)	296 (93%)	23 (7%)	0	100	100
1	J	321/324 (99%)	300 (94%)	21 (6%)	0	100	100
1	K	322/324 (99%)	302 (94%)	20 (6%)	0	100	100
1	L	320/324 (99%)	301 (94%)	19 (6%)	0	100	100
All	All	3851/3888 (99%)	3617 (94%)	234 (6%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	280/281 (100%)	276 (99%)	4 (1%)	78	95
1	B	280/281 (100%)	275 (98%)	5 (2%)	71	93
1	C	279/281 (99%)	276 (99%)	3 (1%)	84	96
1	D	280/281 (100%)	274 (98%)	6 (2%)	66	91
1	E	280/281 (100%)	278 (99%)	2 (1%)	91	98
1	F	279/281 (99%)	277 (99%)	2 (1%)	91	98
1	G	279/281 (99%)	275 (99%)	4 (1%)	78	95
1	H	280/281 (100%)	277 (99%)	3 (1%)	84	96
1	I	276/281 (98%)	273 (99%)	3 (1%)	84	96
1	J	279/281 (99%)	276 (99%)	3 (1%)	84	96
1	K	280/281 (100%)	273 (98%)	7 (2%)	60	89
1	L	279/281 (99%)	273 (98%)	6 (2%)	64	90
All	All	3351/3372 (99%)	3303 (99%)	48 (1%)	78	95

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	GLN
1	A	108	PHE
1	A	135	MET
1	A	220	ARG
1	B	108	PHE
1	B	135	MET
1	B	220	ARG
1	B	233	LYS
1	B	297	LYS
1	C	135	MET
1	C	220	ARG
1	C	303	LEU
1	D	11	ASP
1	D	49	LYS
1	D	108	PHE

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Mol	Chain	Res	Type
1	D	135	MET
1	D	220	ARG
1	D	303	LEU
1	E	135	MET
1	E	220	ARG
1	F	135	MET
1	F	220	ARG
1	G	108	PHE
1	G	135	MET
1	G	220	ARG
1	G	266	THR
1	H	135	MET
1	H	200	ARG
1	H	220	ARG
1	I	135	MET
1	I	220	ARG
1	I	231	ASN
1	J	108	PHE
1	J	135	MET
1	J	220	ARG
1	K	108	PHE
1	K	135	MET
1	K	220	ARG
1	K	266	THR
1	K	294	SER
1	K	322	THR
1	K	324	GLU
1	L	135	MET
1	L	220	ARG
1	L	269	VAL
1	L	290	ASP
1	L	301	ARG
1	L	322	THR

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

Mol	Chain	Res	Type
1	F	289	HIS
1	G	45	HIS
1	I	52	HIS
1	K	274	HIS
1	L	165	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 36 ligands modelled in this entry, 12 are monoatomic - leaving 24 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	$\# Z > 2$
3	SO4	A	402	-	4,4,4	0.23	0	6,6,6	0.11	0
3	SO4	A	403	-	4,4,4	0.22	0	6,6,6	0.07	0
3	SO4	B	402	-	4,4,4	0.22	0	6,6,6	0.11	0
3	SO4	B	403	-	4,4,4	0.23	0	6,6,6	0.08	0
3	SO4	C	402	-	4,4,4	0.23	0	6,6,6	0.06	0
3	SO4	C	403	-	4,4,4	0.22	0	6,6,6	0.07	0
3	SO4	D	402	-	4,4,4	0.23	0	6,6,6	0.11	0
3	SO4	D	403	-	4,4,4	0.23	0	6,6,6	0.10	0
3	SO4	E	402	-	4,4,4	0.22	0	6,6,6	0.07	0
3	SO4	E	403	-	4,4,4	0.24	0	6,6,6	0.08	0
3	SO4	F	402	-	4,4,4	0.23	0	6,6,6	0.08	0
3	SO4	F	403	-	4,4,4	0.24	0	6,6,6	0.09	0
3	SO4	G	402	-	4,4,4	0.23	0	6,6,6	0.07	0
3	SO4	G	403	-	4,4,4	0.23	0	6,6,6	0.08	0
3	SO4	H	402	-	4,4,4	0.23	0	6,6,6	0.10	0
3	SO4	H	403	-	4,4,4	0.22	0	6,6,6	0.08	0
3	SO4	I	402	-	4,4,4	0.24	0	6,6,6	0.06	0
3	SO4	I	403	-	4,4,4	0.23	0	6,6,6	0.07	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	J	402	-	4,4,4	0.23	0	6,6,6	0.06	0
3	SO4	J	403	-	4,4,4	0.23	0	6,6,6	0.08	0
3	SO4	K	402	-	4,4,4	0.24	0	6,6,6	0.09	0
3	SO4	K	403	-	4,4,4	0.22	0	6,6,6	0.08	0
3	SO4	L	402	-	4,4,4	0.24	0	6,6,6	0.08	0
3	SO4	L	403	-	4,4,4	0.23	0	6,6,6	0.06	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	402	-	-	0/0/0/0	0/0/0/0
3	SO4	A	403	-	-	0/0/0/0	0/0/0/0
3	SO4	B	402	-	-	0/0/0/0	0/0/0/0
3	SO4	B	403	-	-	0/0/0/0	0/0/0/0
3	SO4	C	402	-	-	0/0/0/0	0/0/0/0
3	SO4	C	403	-	-	0/0/0/0	0/0/0/0
3	SO4	D	402	-	-	0/0/0/0	0/0/0/0
3	SO4	D	403	-	-	0/0/0/0	0/0/0/0
3	SO4	E	402	-	-	0/0/0/0	0/0/0/0
3	SO4	E	403	-	-	0/0/0/0	0/0/0/0
3	SO4	F	402	-	-	0/0/0/0	0/0/0/0
3	SO4	F	403	-	-	0/0/0/0	0/0/0/0
3	SO4	G	402	-	-	0/0/0/0	0/0/0/0
3	SO4	G	403	-	-	0/0/0/0	0/0/0/0
3	SO4	H	402	-	-	0/0/0/0	0/0/0/0
3	SO4	H	403	-	-	0/0/0/0	0/0/0/0
3	SO4	I	402	-	-	0/0/0/0	0/0/0/0
3	SO4	I	403	-	-	0/0/0/0	0/0/0/0
3	SO4	J	402	-	-	0/0/0/0	0/0/0/0
3	SO4	J	403	-	-	0/0/0/0	0/0/0/0
3	SO4	K	402	-	-	0/0/0/0	0/0/0/0
3	SO4	K	403	-	-	0/0/0/0	0/0/0/0
3	SO4	L	402	-	-	0/0/0/0	0/0/0/0
3	SO4	L	403	-	-	0/0/0/0	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, 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	323/324 (99%)	-0.07	1 (0%) 91 95	43, 60, 88, 99	0
1	B	323/324 (99%)	0.07	9 (2%) 50 56	45, 67, 86, 105	0
1	C	323/324 (99%)	0.24	8 (2%) 54 61	50, 78, 112, 146	0
1	D	322/324 (99%)	-0.04	1 (0%) 91 95	43, 65, 91, 121	0
1	E	323/324 (99%)	0.07	6 (1%) 64 70	48, 66, 91, 120	0
1	F	323/324 (99%)	0.26	10 (3%) 47 52	49, 75, 104, 131	0
1	G	324/324 (100%)	0.22	5 (1%) 70 75	46, 78, 117, 183	0
1	H	324/324 (100%)	0.21	8 (2%) 54 61	51, 75, 105, 159	0
1	I	321/324 (99%)	0.35	17 (5%) 25 28	53, 76, 109, 148	0
1	J	323/324 (99%)	0.13	6 (1%) 64 70	50, 69, 99, 118	0
1	K	324/324 (100%)	0.29	10 (3%) 47 52	54, 77, 103, 140	0
1	L	322/324 (99%)	0.58	17 (5%) 25 28	58, 89, 123, 151	0
All	All	3875/3888 (99%)	0.19	98 (2%) 54 61	43, 72, 107, 183	0

All (98) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	323	ALA	6.0
1	H	324	GLU	4.8
1	F	120	GLY	4.1
1	E	36	GLY	3.9
1	B	136	ASP	3.7
1	I	23	ASP	3.5
1	G	138	CYS	3.4
1	B	138	CYS	3.4
1	L	138	CYS	3.4
1	K	225	SER	3.3
1	K	80	GLY	3.3

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Mol	Chain	Res	Type	RSRZ
1	I	275	HIS	3.2
1	H	225	SER	3.1
1	F	85	ASP	3.1
1	F	137	PRO	3.0
1	L	225	SER	3.0
1	L	42	ASN	3.0
1	K	120	GLY	3.0
1	J	175	GLU	3.0
1	H	322	THR	3.0
1	C	138	CYS	3.0
1	L	85	ASP	3.0
1	G	320	LYS	2.9
1	A	136	ASP	2.8
1	J	322	THR	2.8
1	K	177	ASP	2.8
1	E	138	CYS	2.7
1	I	86	SER	2.7
1	L	276	SER	2.7
1	F	136	ASP	2.7
1	E	37	ILE	2.7
1	I	274	HIS	2.7
1	L	210	LYS	2.7
1	I	224	ALA	2.6
1	I	320	LYS	2.6
1	I	17	ALA	2.6
1	B	225	SER	2.6
1	H	207	GLU	2.6
1	K	224	ALA	2.6
1	K	323	ALA	2.5
1	L	121	PRO	2.5
1	L	62	GLY	2.5
1	D	224	ALA	2.5
1	G	85	ASP	2.5
1	L	86	SER	2.5
1	B	16	PRO	2.5
1	C	225	SER	2.4
1	I	239	SER	2.4
1	I	4	PRO	2.4
1	J	177	ASP	2.4
1	E	10	GLY	2.4
1	H	120	GLY	2.4
1	L	318	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	L	66	ASP	2.4
1	F	276	SER	2.4
1	L	136	ASP	2.4
1	B	86	SER	2.4
1	H	209	GLY	2.3
1	K	38	PRO	2.3
1	K	15	ASP	2.3
1	F	1	MET	2.3
1	I	63	GLU	2.3
1	J	138	CYS	2.3
1	L	232	GLY	2.3
1	I	136	ASP	2.3
1	L	177	ASP	2.2
1	C	121	PRO	2.2
1	F	225	SER	2.2
1	L	144	ASP	2.2
1	B	85	ASP	2.2
1	I	85	ASP	2.2
1	I	225	SER	2.2
1	F	224	ALA	2.2
1	K	138	CYS	2.2
1	C	275	HIS	2.2
1	G	74	GLU	2.2
1	B	276	SER	2.2
1	K	207	GLU	2.2
1	I	34	GLU	2.2
1	C	276	SER	2.1
1	H	224	ALA	2.1
1	H	137	PRO	2.1
1	B	275	HIS	2.1
1	G	274	HIS	2.1
1	L	175	GLU	2.1
1	C	85	ASP	2.1
1	B	137	PRO	2.1
1	F	121	PRO	2.1
1	J	274	HIS	2.1
1	I	15	ASP	2.1
1	C	120	GLY	2.1
1	I	120	GLY	2.1
1	C	24	LYS	2.1
1	F	17	ALA	2.0
1	E	322	THR	2.0

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Mol	Chain	Res	Type	RSRZ
1	J	1	MET	2.0
1	E	61	GLN	2.0
1	L	176	GLY	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	SO4	I	403	5/5	0.45	11.18	162,162,163,163	0
3	SO4	G	403	5/5	0.39	6.85	148,149,149,149	0
3	SO4	K	403	5/5	0.47	5.50	161,161,161,161	0
2	CA	A	401	1/1	0.41	4.95	62,62,62,62	0
3	SO4	D	403	5/5	0.27	4.69	113,113,113,114	0
2	CA	E	401	1/1	0.41	4.59	70,70,70,70	0
3	SO4	L	403	5/5	0.41	4.36	152,153,153,153	0
3	SO4	H	403	5/5	0.28	4.15	113,113,114,114	0
3	SO4	G	402	5/5	0.51	3.93	126,126,126,126	0
3	SO4	A	402	5/5	0.52	3.84	109,109,109,109	0
2	CA	B	401	1/1	0.43	3.83	73,73,73,73	0
3	SO4	C	403	5/5	0.30	3.61	137,137,137,137	0
2	CA	C	401	1/1	0.40	3.39	77,77,77,77	0
3	SO4	E	402	5/5	0.37	3.22	105,105,105,105	0
2	CA	K	401	1/1	0.44	3.10	81,81,81,81	0
3	SO4	D	402	5/5	0.31	2.97	92,93,93,93	0
3	SO4	K	402	5/5	0.41	2.93	110,110,111,111	0
2	CA	I	401	1/1	0.49	2.89	78,78,78,78	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	CA	F	401	1/1	0.44	2.75	84,84,84,84	0
3	SO4	A	403	5/5	0.27	2.74	123,123,124,124	0
3	SO4	J	403	5/5	0.37	2.73	153,154,154,154	0
3	SO4	F	403	5/5	0.27	2.63	136,136,137,137	0
3	SO4	F	402	5/5	0.47	2.57	106,106,107,107	0
2	CA	H	401	1/1	0.41	2.19	71,71,71,71	0
2	CA	G	401	1/1	0.37	2.11	77,77,77,77	0
3	SO4	I	402	5/5	0.42	2.10	105,105,105,106	0
3	SO4	C	402	5/5	0.33	1.95	103,103,104,104	0
2	CA	J	401	1/1	0.33	1.94	70,70,70,70	0
2	CA	D	401	1/1	0.33	1.52	65,65,65,65	0
2	CA	L	401	1/1	0.37	1.10	80,80,80,80	0
3	SO4	E	403	5/5	0.23	1.07	131,131,131,131	0
3	SO4	J	402	5/5	0.29	0.96	101,101,102,102	0
3	SO4	H	402	5/5	0.27	0.92	94,94,94,95	0
3	SO4	B	402	5/5	0.29	0.50	85,85,86,86	0
3	SO4	L	402	5/5	0.28	0.45	111,111,111,111	0
3	SO4	B	403	5/5	0.17	0.43	116,116,116,116	0

6.5 Other polymers ⓘ

There are no such residues in this entry.