



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 01:09 AM GMT

PDB ID : 2FG6
Title : N-succinyl-L-ornithinetranscarbamylase from *B. fragilis* complexed with sulfate and N-succinyl-L-norvaline
Authors : Shi, D.; Yu, X.; Malamy, M.H.; Allewell, N.M.; Mendel, T.
Deposited on : 2005-12-21
Resolution : 2.80 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>

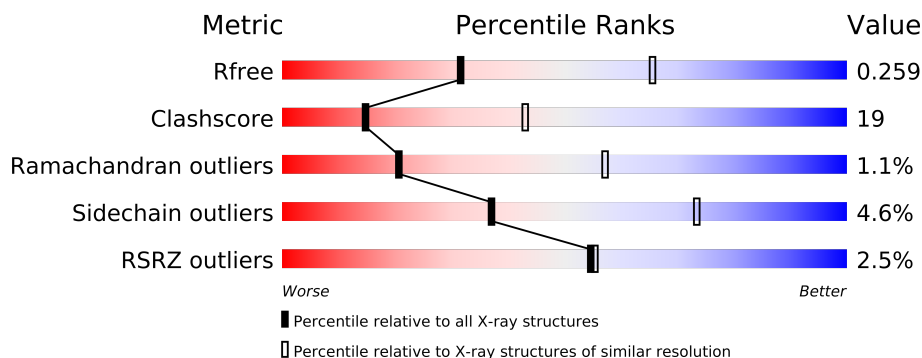
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 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 2.80 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	1799 (2.80-2.80)
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (2.80-2.80)

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	C	338	
1	D	338	
1	E	338	
1	X	338	
1	Y	338	
1	Z	338	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15978 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 putative ornithine carbamoyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	X	322	Total	C	N	O	S	0	0	0
			2593	1651	448	478	16			
1	Y	322	Total	C	N	O	S	0	0	0
			2593	1651	448	478	16			
1	Z	320	Total	C	N	O	S	0	0	0
			2578	1643	443	476	16			
1	C	321	Total	C	N	O	S	0	0	0
			2582	1645	444	477	16			
1	D	322	Total	C	N	O	S	0	0	0
			2593	1651	448	478	16			
1	E	322	Total	C	N	O	S	0	0	0
			2593	1651	448	478	16			

There are 126 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	-19	MET	-	EXPRESSION TAG	UNP Q5LI27
X	-18	GLY	-	EXPRESSION TAG	UNP Q5LI27
X	-17	SER	-	EXPRESSION TAG	UNP Q5LI27
X	-16	SER	-	EXPRESSION TAG	UNP Q5LI27
X	-15	HIS	-	EXPRESSION TAG	UNP Q5LI27
X	-14	HIS	-	EXPRESSION TAG	UNP Q5LI27
X	-13	HIS	-	EXPRESSION TAG	UNP Q5LI27
X	-12	HIS	-	EXPRESSION TAG	UNP Q5LI27
X	-11	HIS	-	EXPRESSION TAG	UNP Q5LI27
X	-10	HIS	-	EXPRESSION TAG	UNP Q5LI27
X	-9	SER	-	EXPRESSION TAG	UNP Q5LI27
X	-8	SER	-	EXPRESSION TAG	UNP Q5LI27
X	-7	GLY	-	EXPRESSION TAG	UNP Q5LI27
X	-6	LEU	-	EXPRESSION TAG	UNP Q5LI27
X	-5	VAL	-	EXPRESSION TAG	UNP Q5LI27
X	-4	PRO	-	EXPRESSION TAG	UNP Q5LI27
X	-3	ARG	-	EXPRESSION TAG	UNP Q5LI27

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Chain	Residue	Modelled	Actual	Comment	Reference
X	-2	GLY	-	EXPRESSION TAG	UNP Q5LI27
X	-1	SER	-	EXPRESSION TAG	UNP Q5LI27
X	0	HIS	-	EXPRESSION TAG	UNP Q5LI27
X	242	LEU	THR	ENGINEERED	UNP Q5LI27
Y	-19	MET	-	EXPRESSION TAG	UNP Q5LI27
Y	-18	GLY	-	EXPRESSION TAG	UNP Q5LI27
Y	-17	SER	-	EXPRESSION TAG	UNP Q5LI27
Y	-16	SER	-	EXPRESSION TAG	UNP Q5LI27
Y	-15	HIS	-	EXPRESSION TAG	UNP Q5LI27
Y	-14	HIS	-	EXPRESSION TAG	UNP Q5LI27
Y	-13	HIS	-	EXPRESSION TAG	UNP Q5LI27
Y	-12	HIS	-	EXPRESSION TAG	UNP Q5LI27
Y	-11	HIS	-	EXPRESSION TAG	UNP Q5LI27
Y	-10	HIS	-	EXPRESSION TAG	UNP Q5LI27
Y	-9	SER	-	EXPRESSION TAG	UNP Q5LI27
Y	-8	SER	-	EXPRESSION TAG	UNP Q5LI27
Y	-7	GLY	-	EXPRESSION TAG	UNP Q5LI27
Y	-6	LEU	-	EXPRESSION TAG	UNP Q5LI27
Y	-5	VAL	-	EXPRESSION TAG	UNP Q5LI27
Y	-4	PRO	-	EXPRESSION TAG	UNP Q5LI27
Y	-3	ARG	-	EXPRESSION TAG	UNP Q5LI27
Y	-2	GLY	-	EXPRESSION TAG	UNP Q5LI27
Y	-1	SER	-	EXPRESSION TAG	UNP Q5LI27
Y	0	HIS	-	EXPRESSION TAG	UNP Q5LI27
Y	242	LEU	THR	ENGINEERED	UNP Q5LI27
Z	-19	MET	-	EXPRESSION TAG	UNP Q5LI27
Z	-18	GLY	-	EXPRESSION TAG	UNP Q5LI27
Z	-17	SER	-	EXPRESSION TAG	UNP Q5LI27
Z	-16	SER	-	EXPRESSION TAG	UNP Q5LI27
Z	-15	HIS	-	EXPRESSION TAG	UNP Q5LI27
Z	-14	HIS	-	EXPRESSION TAG	UNP Q5LI27
Z	-13	HIS	-	EXPRESSION TAG	UNP Q5LI27
Z	-12	HIS	-	EXPRESSION TAG	UNP Q5LI27
Z	-11	HIS	-	EXPRESSION TAG	UNP Q5LI27
Z	-10	HIS	-	EXPRESSION TAG	UNP Q5LI27
Z	-9	SER	-	EXPRESSION TAG	UNP Q5LI27
Z	-8	SER	-	EXPRESSION TAG	UNP Q5LI27
Z	-7	GLY	-	EXPRESSION TAG	UNP Q5LI27
Z	-6	LEU	-	EXPRESSION TAG	UNP Q5LI27
Z	-5	VAL	-	EXPRESSION TAG	UNP Q5LI27
Z	-4	PRO	-	EXPRESSION TAG	UNP Q5LI27
Z	-3	ARG	-	EXPRESSION TAG	UNP Q5LI27

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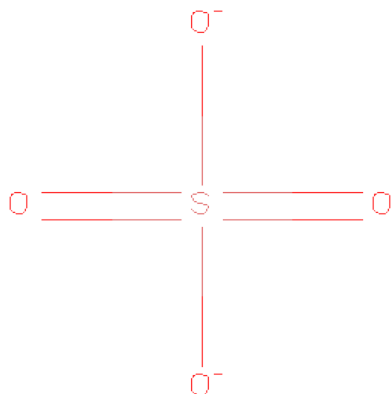
Chain	Residue	Modelled	Actual	Comment	Reference
Z	-2	GLY	-	EXPRESSION TAG	UNP Q5LI27
Z	-1	SER	-	EXPRESSION TAG	UNP Q5LI27
Z	0	HIS	-	EXPRESSION TAG	UNP Q5LI27
Z	242	LEU	THR	ENGINEERED	UNP Q5LI27
C	-19	MET	-	EXPRESSION TAG	UNP Q5LI27
C	-18	GLY	-	EXPRESSION TAG	UNP Q5LI27
C	-17	SER	-	EXPRESSION TAG	UNP Q5LI27
C	-16	SER	-	EXPRESSION TAG	UNP Q5LI27
C	-15	HIS	-	EXPRESSION TAG	UNP Q5LI27
C	-14	HIS	-	EXPRESSION TAG	UNP Q5LI27
C	-13	HIS	-	EXPRESSION TAG	UNP Q5LI27
C	-12	HIS	-	EXPRESSION TAG	UNP Q5LI27
C	-11	HIS	-	EXPRESSION TAG	UNP Q5LI27
C	-10	HIS	-	EXPRESSION TAG	UNP Q5LI27
C	-9	SER	-	EXPRESSION TAG	UNP Q5LI27
C	-8	SER	-	EXPRESSION TAG	UNP Q5LI27
C	-7	GLY	-	EXPRESSION TAG	UNP Q5LI27
C	-6	LEU	-	EXPRESSION TAG	UNP Q5LI27
C	-5	VAL	-	EXPRESSION TAG	UNP Q5LI27
C	-4	PRO	-	EXPRESSION TAG	UNP Q5LI27
C	-3	ARG	-	EXPRESSION TAG	UNP Q5LI27
C	-2	GLY	-	EXPRESSION TAG	UNP Q5LI27
C	-1	SER	-	EXPRESSION TAG	UNP Q5LI27
C	0	HIS	-	EXPRESSION TAG	UNP Q5LI27
C	242	LEU	THR	ENGINEERED	UNP Q5LI27
D	-19	MET	-	EXPRESSION TAG	UNP Q5LI27
D	-18	GLY	-	EXPRESSION TAG	UNP Q5LI27
D	-17	SER	-	EXPRESSION TAG	UNP Q5LI27
D	-16	SER	-	EXPRESSION TAG	UNP Q5LI27
D	-15	HIS	-	EXPRESSION TAG	UNP Q5LI27
D	-14	HIS	-	EXPRESSION TAG	UNP Q5LI27
D	-13	HIS	-	EXPRESSION TAG	UNP Q5LI27
D	-12	HIS	-	EXPRESSION TAG	UNP Q5LI27
D	-11	HIS	-	EXPRESSION TAG	UNP Q5LI27
D	-10	HIS	-	EXPRESSION TAG	UNP Q5LI27
D	-9	SER	-	EXPRESSION TAG	UNP Q5LI27
D	-8	SER	-	EXPRESSION TAG	UNP Q5LI27
D	-7	GLY	-	EXPRESSION TAG	UNP Q5LI27
D	-6	LEU	-	EXPRESSION TAG	UNP Q5LI27
D	-5	VAL	-	EXPRESSION TAG	UNP Q5LI27
D	-4	PRO	-	EXPRESSION TAG	UNP Q5LI27
D	-3	ARG	-	EXPRESSION TAG	UNP Q5LI27

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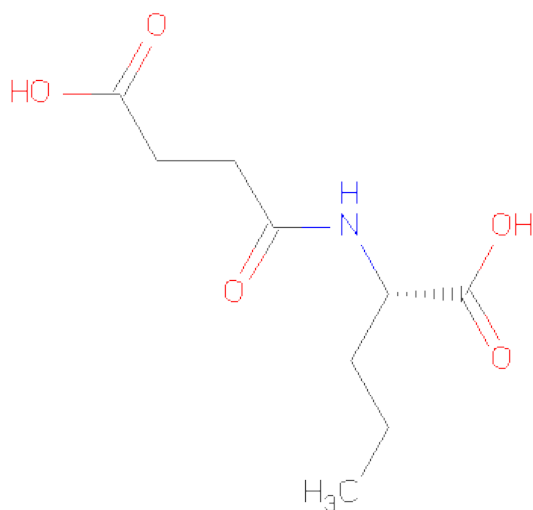
Chain	Residue	Modelled	Actual	Comment	Reference
D	-2	GLY	-	EXPRESSION TAG	UNP Q5LI27
D	-1	SER	-	EXPRESSION TAG	UNP Q5LI27
D	0	HIS	-	EXPRESSION TAG	UNP Q5LI27
D	242	LEU	THR	ENGINEERED	UNP Q5LI27
E	-19	MET	-	EXPRESSION TAG	UNP Q5LI27
E	-18	GLY	-	EXPRESSION TAG	UNP Q5LI27
E	-17	SER	-	EXPRESSION TAG	UNP Q5LI27
E	-16	SER	-	EXPRESSION TAG	UNP Q5LI27
E	-15	HIS	-	EXPRESSION TAG	UNP Q5LI27
E	-14	HIS	-	EXPRESSION TAG	UNP Q5LI27
E	-13	HIS	-	EXPRESSION TAG	UNP Q5LI27
E	-12	HIS	-	EXPRESSION TAG	UNP Q5LI27
E	-11	HIS	-	EXPRESSION TAG	UNP Q5LI27
E	-10	HIS	-	EXPRESSION TAG	UNP Q5LI27
E	-9	SER	-	EXPRESSION TAG	UNP Q5LI27
E	-8	SER	-	EXPRESSION TAG	UNP Q5LI27
E	-7	GLY	-	EXPRESSION TAG	UNP Q5LI27
E	-6	LEU	-	EXPRESSION TAG	UNP Q5LI27
E	-5	VAL	-	EXPRESSION TAG	UNP Q5LI27
E	-4	PRO	-	EXPRESSION TAG	UNP Q5LI27
E	-3	ARG	-	EXPRESSION TAG	UNP Q5LI27
E	-2	GLY	-	EXPRESSION TAG	UNP Q5LI27
E	-1	SER	-	EXPRESSION TAG	UNP Q5LI27
E	0	HIS	-	EXPRESSION TAG	UNP Q5LI27
E	242	LEU	THR	ENGINEERED	UNP Q5LI27

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	X	1	Total	O	S	0	0
			5	4	1		
2	Y	1	Total	O	S	0	0
			5	4	1		
2	Z	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	X	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is N-(3-CARBOXYPROPANOYL)-L-NORVALINE (three-letter code: SN0) (formula: C₉H₁₅NO₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	X	1	Total	C	N	O	0	0
			15	9	1	5		
3	Y	1	Total	C	N	O	0	0
			15	9	1	5		
3	Z	1	Total	C	N	O	0	0
			15	9	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	N	O	0	0
			15	9	1	5		
3	D	1	Total	C	N	O	0	0
			15	9	1	5		
3	E	1	Total	C	N	O	0	0
			15	9	1	5		

- Molecule 4 is water.

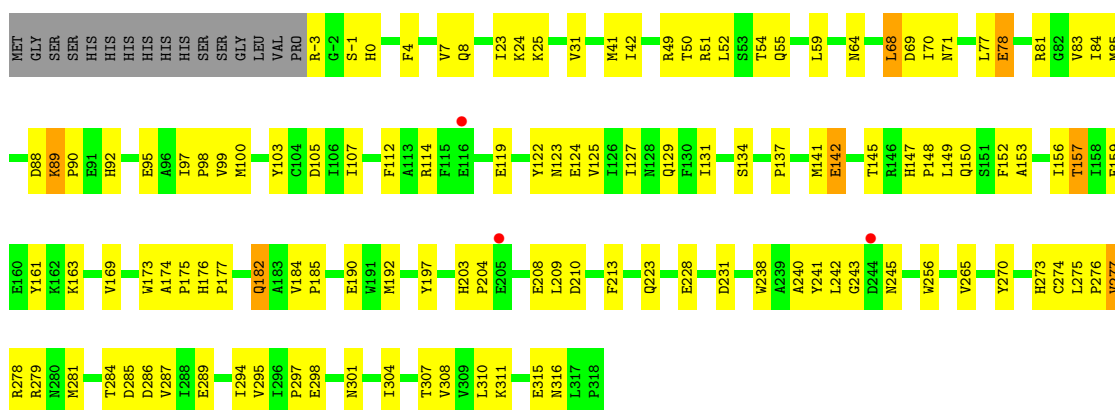
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	27	Total	O	0	0
			27	27		
4	D	25	Total	O	0	0
			25	25		
4	E	38	Total	O	0	0
			38	38		
4	X	62	Total	O	0	0
			62	62		
4	Y	109	Total	O	0	0
			109	109		
4	Z	55	Total	O	0	0
			55	55		

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.

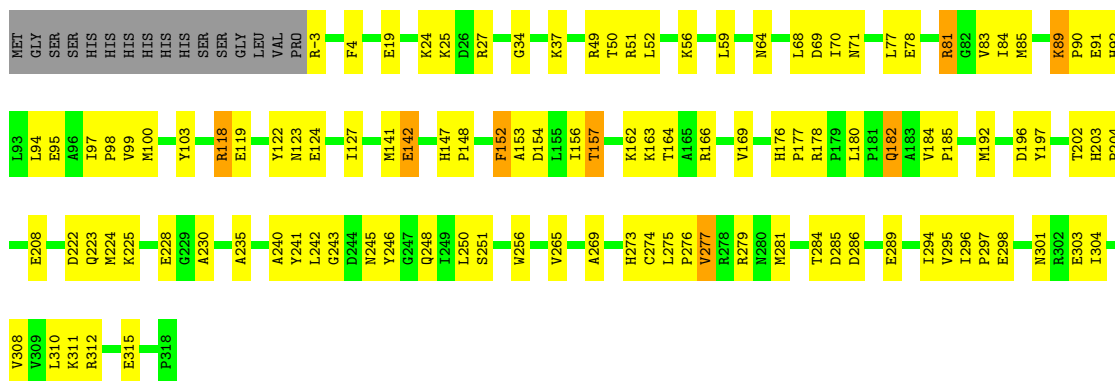
- Molecule 1: putative ornithine carbamoyltransferase

Chain X: 



- Molecule 1: putative ornithine carbamoyltransferase

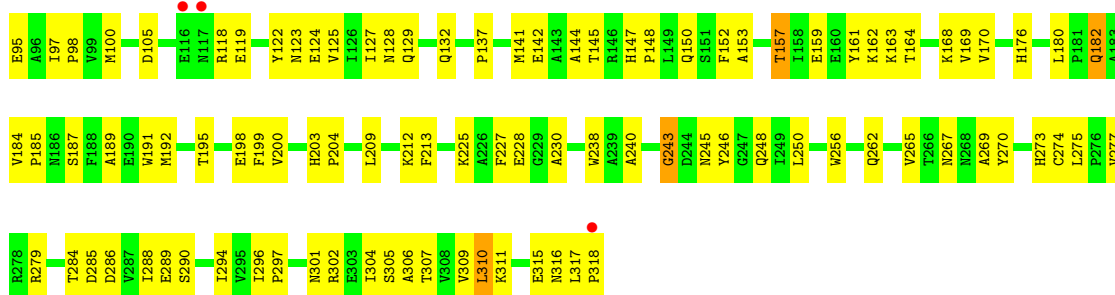
Chain Y: 



- Molecule 1: putative ornithine carbamoyltransferase

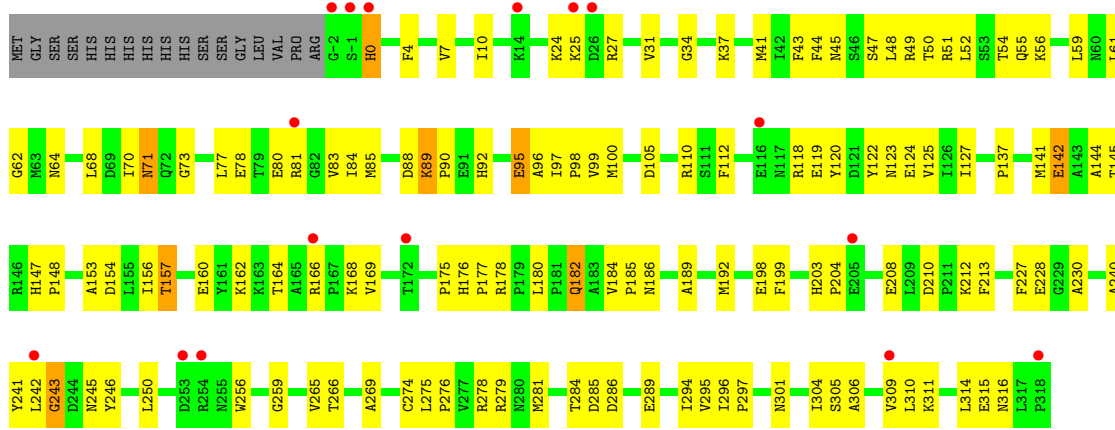
Chain Z: 





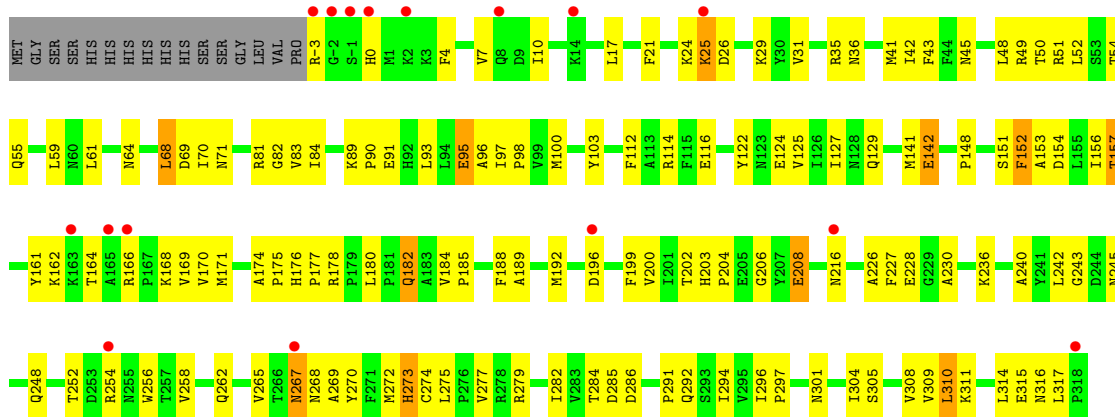
• Molecule 1: putative ornithine carbamoyltransferase

Chain C:



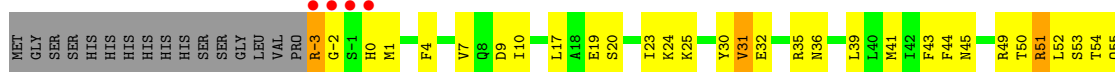
• Molecule 1: putative ornithine carbamoyltransferase

Chain D:



• Molecule 1: putative ornithine carbamoyltransferase

Chain E:





4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	155.70Å 155.70Å 120.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.99 – 2.80 19.98 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.5 (19.99-2.80) 99.6 (19.98-2.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.87 (at 2.79Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.216 , 0.260 0.216 , 0.259	Depositor DCC
R_{free} test set	3585 reflections (5.09%)	DCC
Wilson B-factor (Å ²)	51.6	Xtriage
Anisotropy	0.314	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 27.4	EDS
Estimated twinning fraction	0.012 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 70401 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	15978	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.23% 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: SO4, SN0

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	C	0.34	0/2642	0.60	0/3579
1	D	0.36	0/2653	0.61	0/3593
1	E	0.38	0/2653	0.63	0/3593
1	X	0.39	0/2653	0.65	0/3593
1	Y	0.42	0/2653	0.68	1/3593 (0.0%)
1	Z	0.40	0/2638	0.66	0/3574
All	All	0.38	0/15892	0.64	1/21525 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Y	68	LEU	CA-CB-CG	5.31	127.51	115.30

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	C	2582	0	2551	118	0
1	D	2593	0	2564	115	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2593	0	2564	107	0
1	X	2593	0	2564	100	0
1	Y	2593	0	2564	96	0
1	Z	2578	0	2548	93	0
2	C	10	0	0	1	0
2	D	5	0	0	0	0
2	E	5	0	0	0	0
2	X	10	0	0	0	0
2	Y	5	0	0	0	0
2	Z	5	0	0	0	0
3	C	15	0	13	2	0
3	D	15	0	13	1	0
3	E	15	0	13	2	0
3	X	15	0	13	0	0
3	Y	15	0	13	1	0
3	Z	15	0	13	0	0
4	C	27	0	0	4	0
4	D	25	0	0	8	0
4	E	38	0	0	4	0
4	X	62	0	0	5	0
4	Y	109	0	0	1	0
4	Z	55	0	0	4	0
All	All	15978	0	15433	597	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 19.

All (597) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:Z:169:VAL:HG11	1:Z:192:MET:HE2	1.45	0.98
1:Y:153:ALA:O	1:Y:157:THR:HG22	1.67	0.94
1:X:24:LYS:HZ3	1:X:301:ASN:HD21	1.14	0.92
1:X:24:LYS:NZ	1:X:301:ASN:HD21	1.68	0.90
1:X:153:ALA:O	1:X:157:THR:HG22	1.72	0.89
1:D:153:ALA:O	1:D:157:THR:HG22	1.74	0.88
1:C:0:HIS:HA	1:C:316:ASN:ND2	1.90	0.86
1:C:97:ILE:HB	1:C:98:PRO:HD3	1.56	0.86
1:E:153:ALA:O	1:E:157:THR:HG22	1.77	0.84
1:Z:153:ALA:O	1:Z:157:THR:HG22	1.77	0.83
1:X:24:LYS:HG2	1:X:304:ILE:HG12	1.61	0.82
1:Y:24:LYS:NZ	1:Y:301:ASN:HD21	1.79	0.80
1:X:0:HIS:HA	1:X:316:ASN:ND2	1.95	0.80

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:41:MET:HE3	1:E:54:THR:HB	1.63	0.80
1:Y:24:LYS:HG2	1:Y:304:ILE:HG12	1.64	0.80
1:C:169:VAL:HG11	1:C:192:MET:HE2	1.64	0.80
1:C:119:GLU:HG2	1:C:123:ASN:HD21	1.47	0.79
1:Z:0:HIS:HA	1:Z:316:ASN:ND2	1.97	0.79
1:C:96:ALA:O	1:C:100:MET:HG3	1.83	0.78
1:D:169:VAL:HG11	1:D:192:MET:HE2	1.65	0.78
1:D:0:HIS:HA	1:D:316:ASN:ND2	1.98	0.78
1:C:119:GLU:HG2	1:C:123:ASN:ND2	1.98	0.78
1:X:311:LYS:O	1:X:315:GLU:HG3	1.84	0.78
1:Z:50:THR:HG21	1:Z:141:MET:HB3	1.64	0.78
1:D:124:GLU:HB3	1:D:127:ILE:CG2	2.15	0.77
1:D:24:LYS:HG2	1:D:304:ILE:HG12	1.66	0.77
1:Y:169:VAL:HG11	1:Y:192:MET:HE2	1.67	0.76
1:Z:24:LYS:NZ	1:Z:301:ASN:HD21	1.83	0.76
1:C:41:MET:HE3	1:C:54:THR:HB	1.68	0.75
1:X:284:THR:HG22	1:X:286:ASP:H	1.52	0.75
1:Y:24:LYS:HZ3	1:Y:301:ASN:HD21	1.30	0.75
1:C:284:THR:HG22	1:C:286:ASP:H	1.51	0.75
1:C:153:ALA:O	1:C:157:THR:HG22	1.87	0.75
1:E:0:HIS:HA	1:E:316:ASN:ND2	2.01	0.74
1:D:0:HIS:HA	1:D:316:ASN:HD22	1.52	0.74
1:D:97:ILE:HB	1:D:98:PRO:HD3	1.70	0.73
1:E:35:ARG:HG2	1:E:36:ASN:HD22	1.54	0.72
1:X:95:GLU:O	1:X:98:PRO:HD2	1.89	0.72
1:C:124:GLU:HB3	1:C:127:ILE:CG2	2.20	0.72
1:E:273:HIS:ND1	1:E:277:VAL:HG22	2.04	0.72
1:E:124:GLU:HB3	1:E:127:ILE:CG2	2.20	0.72
1:Y:284:THR:HG22	1:Y:286:ASP:H	1.53	0.72
1:E:294:ILE:C	1:E:297:PRO:HD2	2.11	0.70
1:E:169:VAL:HG11	1:E:192:MET:HE2	1.71	0.70
1:Y:124:GLU:HB3	1:Y:127:ILE:CG2	2.20	0.70
1:D:204:PRO:HG3	1:D:256:TRP:CE2	2.27	0.70
1:Y:284:THR:HG22	1:Y:286:ASP:N	2.07	0.69
1:X:122:TYR:OH	1:X:182:GLN:HG3	1.92	0.69
1:D:176:HIS:CD2	1:D:180:LEU:HD11	2.27	0.69
1:Z:169:VAL:HG11	1:Z:192:MET:CE	2.22	0.69
1:D:184:VAL:HB	1:D:185:PRO:HD3	1.76	0.68
1:X:-3:ARG:NH1	1:X:8:GLN:HB3	2.08	0.68
1:E:95:GLU:O	1:E:99:VAL:HG23	1.94	0.68
1:D:91:GLU:HG3	4:D:368:HOH:O	1.94	0.67
1:X:274:CYS:O	1:X:275:LEU:HB2	1.93	0.67

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:50:THR:HG21	1:C:141:MET:HB3	1.76	0.67
1:C:49:ARG:HH21	1:C:276:PRO:HG3	1.59	0.67
1:Z:311:LYS:O	1:Z:315:GLU:HG3	1.95	0.67
1:C:59:LEU:HD13	1:E:64:ASN:HB3	1.77	0.67
1:Y:192:MET:HE3	1:Y:197:TYR:HD1	1.60	0.66
1:C:24:LYS:HG2	1:C:304:ILE:HG12	1.76	0.66
1:Z:24:LYS:HG2	1:Z:304:ILE:HG12	1.76	0.66
1:X:124:GLU:HB3	1:X:127:ILE:CG2	2.26	0.66
1:X:285:ASP:O	1:X:289:GLU:HG3	1.95	0.66
1:E:-3:ARG:HH11	1:E:-3:ARG:HB3	1.60	0.65
1:C:61:LEU:HD13	1:C:314:LEU:HD12	1.79	0.65
1:Y:122:TYR:OH	1:Y:182:GLN:HG3	1.95	0.64
1:Z:24:LYS:HZ2	1:Z:301:ASN:HD21	1.44	0.64
1:Y:285:ASP:O	1:Y:289:GLU:HG3	1.96	0.64
1:Y:311:LYS:O	1:Y:315:GLU:HG3	1.97	0.64
1:Z:119:GLU:HG2	1:Z:123:ASN:ND2	2.13	0.64
1:D:208:GLU:CG	1:D:242:LEU:HD11	2.28	0.63
1:Z:70:ILE:O	1:Z:71:ASN:HB2	1.99	0.63
1:X:169:VAL:HG11	1:X:192:MET:HE2	1.80	0.63
1:X:24:LYS:NZ	1:X:301:ASN:ND2	2.45	0.63
1:Y:204:PRO:HG3	1:Y:256:TRP:CE2	2.34	0.63
1:E:24:LYS:HG2	1:E:304:ILE:HG12	1.81	0.63
1:E:279:ARG:HD3	1:E:285:ASP:OD2	1.97	0.62
1:Y:70:ILE:O	1:Y:71:ASN:HB2	1.99	0.62
1:Y:228:GLU:HA	1:Y:265:VAL:HG12	1.81	0.62
1:Y:208:GLU:HG2	1:Y:242:LEU:HD11	1.80	0.62
1:X:265:VAL:O	1:X:265:VAL:HG12	2.00	0.62
1:E:291:PRO:HG2	1:E:292:GLN:NE2	2.14	0.62
1:Z:274:CYS:O	1:Z:275:LEU:HB2	1.99	0.62
1:D:245:ASN:O	1:D:248:GLN:HB2	1.99	0.62
1:Z:89:LYS:HD2	1:Z:89:LYS:N	2.15	0.62
1:D:95:GLU:O	1:D:98:PRO:HD2	1.99	0.62
1:Z:97:ILE:HB	1:Z:98:PRO:HD3	1.82	0.62
1:E:88:ASP:HB3	1:E:89:LYS:HD2	1.82	0.62
1:Y:70:ILE:HD13	1:Y:77:LEU:HD11	1.82	0.62
1:Z:284:THR:HG21	4:Z:383:HOH:O	2.00	0.62
1:D:125:VAL:O	1:D:129:GLN:HG3	2.00	0.61
1:Z:95:GLU:O	1:Z:98:PRO:HD2	2.01	0.61
1:D:311:LYS:O	1:D:315:GLU:HG3	2.01	0.61
1:Y:97:ILE:HB	1:Y:98:PRO:HD3	1.81	0.61
1:E:70:ILE:O	1:E:71:ASN:HB2	2.01	0.61
1:D:142:GLU:OE1	3:D:353:SN0:HD3	2.01	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:34:GLY:O	1:C:37:LYS:HB2	2.00	0.61
1:E:97:ILE:HB	1:E:98:PRO:HD3	1.83	0.60
1:D:208:GLU:HG2	1:D:242:LEU:HD11	1.81	0.60
1:X:284:THR:HG22	1:X:286:ASP:N	2.15	0.60
1:Z:279:ARG:HD3	1:Z:285:ASP:OD2	2.01	0.60
1:Y:279:ARG:HD3	1:Y:285:ASP:OD2	2.01	0.60
1:E:291:PRO:HG2	1:E:292:GLN:HE22	1.66	0.60
1:E:89:LYS:N	1:E:89:LYS:HD2	2.17	0.60
1:E:56:LYS:NZ	1:E:303:GLU:OE2	2.33	0.60
1:C:95:GLU:O	1:C:99:VAL:HG23	2.02	0.60
1:D:70:ILE:O	1:D:71:ASN:HB2	2.02	0.59
1:E:35:ARG:HG2	1:E:36:ASN:ND2	2.17	0.59
1:Z:204:PRO:HG3	1:Z:256:TRP:CE2	2.36	0.59
1:D:51:ARG:NH2	1:D:69:ASP:OD2	2.35	0.59
1:D:61:LEU:HD13	1:D:314:LEU:HD12	1.83	0.59
1:D:279:ARG:HD3	1:D:285:ASP:OD2	2.02	0.59
1:C:176:HIS:CD2	1:C:180:LEU:HD11	2.37	0.59
1:D:51:ARG:O	1:D:55:GLN:HG3	2.02	0.59
1:D:304:ILE:O	1:D:308:VAL:HG23	2.04	0.58
1:X:192:MET:HE3	1:X:197:TYR:HD1	1.67	0.58
1:Y:49:ARG:NH2	1:Y:276:PRO:HG3	2.17	0.58
1:D:96:ALA:O	1:D:100:MET:HG3	2.03	0.58
1:E:124:GLU:HB3	1:E:127:ILE:HG22	1.86	0.58
1:Y:124:GLU:HB3	1:Y:127:ILE:HG22	1.85	0.58
1:X:41:MET:HE3	1:X:54:THR:HB	1.85	0.58
1:X:0:HIS:HA	1:X:316:ASN:HD21	1.68	0.58
1:Z:275:LEU:HD12	1:Z:277:VAL:CG2	2.34	0.58
1:C:279:ARG:HD3	1:C:285:ASP:OD2	2.04	0.58
1:Y:95:GLU:O	1:Y:98:PRO:HD2	2.03	0.58
1:E:294:ILE:O	1:E:297:PRO:HD2	2.03	0.58
1:C:284:THR:HG21	4:C:359:HOH:O	2.03	0.58
1:C:296:ILE:HB	1:C:297:PRO:HD3	1.86	0.57
1:X:284:THR:HB	1:X:287:VAL:HG23	1.85	0.57
1:C:70:ILE:O	1:C:71:ASN:HB2	2.05	0.57
1:E:19:GLU:O	1:E:23:ILE:HG13	2.04	0.57
1:Z:124:GLU:HB3	1:Z:127:ILE:CG2	2.34	0.57
1:D:169:VAL:HG11	1:D:192:MET:CE	2.33	0.57
1:E:284:THR:HG22	1:E:286:ASP:H	1.70	0.57
1:Y:176:HIS:CD2	1:Y:180:LEU:HD11	2.39	0.57
1:Y:208:GLU:CG	1:Y:242:LEU:HD11	2.34	0.57
1:C:7:VAL:O	1:C:10:ILE:HG22	2.04	0.57
1:E:96:ALA:O	1:E:100:MET:HG3	2.05	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:X:157:THR:HG21	1:X:298:GLU:HG3	1.86	0.57
1:D:64:ASN:HB3	1:E:59:LEU:HD13	1.86	0.57
1:Y:184:VAL:HB	1:Y:185:PRO:HD3	1.86	0.57
1:C:0:HIS:HA	1:C:316:ASN:HD22	1.64	0.56
1:Y:203:HIS:HD2	1:Y:204:PRO:O	1.88	0.56
1:Z:182:GLN:C	1:Z:185:PRO:HD2	2.25	0.56
1:X:70:ILE:HD11	1:X:100:MET:HE1	1.87	0.56
1:X:114:ARG:HD3	4:X:396:HOH:O	2.05	0.56
1:C:70:ILE:HD13	1:C:77:LEU:HD11	1.86	0.56
1:Z:305:SER:O	1:Z:309:VAL:HG23	2.06	0.56
1:Y:192:MET:CE	1:Y:197:TYR:HD1	2.19	0.56
1:E:203:HIS:HD2	1:E:204:PRO:O	1.88	0.56
1:X:4:PHE:CE2	1:X:148:PRO:HB2	2.39	0.56
1:C:241:TYR:CD2	1:C:242:LEU:HD23	2.41	0.55
1:X:192:MET:CE	1:X:197:TYR:HD1	2.19	0.55
1:D:50:THR:HG22	1:D:141:MET:HE2	1.87	0.55
1:C:184:VAL:HB	1:C:185:PRO:HD3	1.88	0.55
1:X:24:LYS:HZ3	1:X:301:ASN:ND2	1.94	0.55
1:C:47:SER:HB2	1:C:110:ARG:NH1	2.22	0.55
1:Z:285:ASP:O	1:Z:289:GLU:HG3	2.06	0.55
1:Z:267:ASN:O	4:Z:372:HOH:O	2.18	0.55
1:Y:24:LYS:NZ	1:Y:301:ASN:ND2	2.54	0.55
1:C:285:ASP:O	1:C:289:GLU:HG3	2.06	0.55
1:E:230:ALA:O	1:E:269:ALA:HB2	2.06	0.55
1:D:166:ARG:HG2	1:D:196:ASP:OD2	2.06	0.55
1:E:70:ILE:HG22	1:E:71:ASN:N	2.22	0.55
1:Z:161:TYR:HB3	1:Z:270:TYR:CE1	2.42	0.55
1:X:88:ASP:HB3	1:X:89:LYS:HD2	1.88	0.55
1:Z:176:HIS:CD2	1:Z:180:LEU:HD11	2.42	0.55
1:D:45:ASN:ND2	1:D:112:PHE:HB2	2.22	0.55
1:Y:274:CYS:O	1:Y:275:LEU:HB2	2.06	0.54
1:D:0:HIS:CA	1:D:316:ASN:ND2	2.69	0.54
1:D:50:THR:HG21	1:D:141:MET:HB3	1.88	0.54
1:X:182:GLN:C	1:X:185:PRO:HD2	2.28	0.54
1:Z:4:PHE:CE2	1:Z:148:PRO:HB2	2.42	0.54
1:E:182:GLN:C	1:E:185:PRO:HD2	2.28	0.54
1:Z:184:VAL:HB	1:Z:185:PRO:HD3	1.88	0.54
1:E:49:ARG:HH21	1:E:276:PRO:HG3	1.73	0.54
1:E:105:ASP:O	1:E:137:PRO:HD2	2.07	0.54
1:D:305:SER:O	1:D:309:VAL:HG23	2.08	0.54
1:X:210:ASP:HA	1:X:241:TYR:CE2	2.43	0.54
1:Y:49:ARG:HH21	1:Y:276:PRO:HG3	1.71	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:-3:ARG:HH11	1:E:-3:ARG:CB	2.19	0.54
1:X:7:VAL:HG12	1:X:190:GLU:OE2	2.07	0.54
1:E:306:ALA:O	1:E:310:LEU:HB2	2.08	0.54
1:Y:119:GLU:HG2	1:Y:123:ASN:ND2	2.22	0.54
1:C:56:LYS:HE2	4:E:378:HOH:O	2.07	0.54
1:X:294:ILE:C	1:X:297:PRO:HD2	2.28	0.54
1:C:245:ASN:HB3	1:C:250:LEU:HD21	1.90	0.53
1:C:119:GLU:HB2	4:C:371:HOH:O	2.09	0.53
1:C:305:SER:O	1:C:309:VAL:HG23	2.08	0.53
1:X:97:ILE:HB	1:X:98:PRO:HD3	1.91	0.53
1:D:122:TYR:OH	1:D:182:GLN:HG3	2.08	0.53
1:C:230:ALA:O	1:C:269:ALA:HB2	2.09	0.53
1:D:189:ALA:HA	1:D:199:PHE:CE2	2.43	0.53
1:Y:154:ASP:O	1:Y:157:THR:HG23	2.09	0.53
1:X:51:ARG:O	1:X:55:GLN:HG3	2.08	0.53
1:D:161:TYR:HB3	1:D:270:TYR:CE1	2.44	0.53
1:D:242:LEU:HG	4:D:363:HOH:O	2.07	0.53
1:D:265:VAL:HG12	1:D:265:VAL:O	2.09	0.53
1:Y:50:THR:HG21	1:Y:141:MET:HB3	1.89	0.53
1:Z:0:HIS:HA	1:Z:316:ASN:HD22	1.74	0.53
1:D:24:LYS:CG	1:D:304:ILE:HG12	2.39	0.53
1:D:284:THR:HG21	4:D:355:HOH:O	2.09	0.53
1:C:120:TYR:O	1:C:125:VAL:HG23	2.08	0.53
1:D:93:LEU:HG	1:D:97:ILE:HD11	1.91	0.52
1:Y:27:ARG:HD2	1:Y:303:GLU:OE2	2.09	0.52
1:X:70:ILE:O	1:X:71:ASN:HB2	2.09	0.52
1:Y:275:LEU:HD11	1:Y:295:VAL:HG13	1.91	0.52
1:D:4:PHE:CE2	1:D:148:PRO:HB2	2.44	0.52
1:E:235:ALA:O	1:E:273:HIS:CD2	2.63	0.52
1:C:4:PHE:CD2	1:C:309:VAL:HG11	2.45	0.52
1:C:203:HIS:HD2	1:C:204:PRO:O	1.93	0.52
1:C:24:LYS:NZ	1:C:301:ASN:HD21	2.08	0.52
1:C:83:VAL:HG12	1:C:84:ILE:N	2.24	0.52
1:C:45:ASN:ND2	1:C:112:PHE:HB2	2.25	0.52
1:Y:24:LYS:HG2	1:Y:304:ILE:CG1	2.39	0.52
1:E:273:HIS:CE1	1:E:277:VAL:HG22	2.45	0.52
1:D:68:LEU:HD22	1:D:69:ASP:N	2.24	0.52
1:X:89:LYS:N	1:X:89:LYS:HD2	2.25	0.52
1:Y:273:HIS:O	1:Y:275:LEU:HD13	2.09	0.52
1:D:230:ALA:O	1:D:269:ALA:HB2	2.09	0.52
1:C:240:ALA:HB2	1:C:250:LEU:HG	1.92	0.52
1:E:188:PHE:O	1:E:192:MET:HG2	2.10	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:77:LEU:O	1:C:89:LYS:HG2	2.10	0.52
1:Z:83:VAL:HG12	1:Z:84:ILE:N	2.25	0.52
1:Z:228:GLU:HG2	4:Z:401:HOH:O	2.10	0.52
1:D:273:HIS:CD2	1:D:277:VAL:HG13	2.45	0.52
1:D:291:PRO:HG2	1:D:292:GLN:HE22	1.75	0.52
1:D:291:PRO:HG2	1:D:292:GLN:NE2	2.25	0.51
1:C:122:TYR:OH	1:C:182:GLN:HG3	2.10	0.51
1:C:189:ALA:HA	1:C:199:PHE:CE2	2.45	0.51
1:X:294:ILE:O	1:X:297:PRO:HD2	2.11	0.51
1:X:175:PRO:HG2	1:X:281:MET:O	2.10	0.51
1:E:210:ASP:HB3	1:E:213:PHE:HD1	1.74	0.51
1:C:51:ARG:O	1:C:55:GLN:HG3	2.11	0.51
1:D:90:PRO:HB3	3:E:355:SN0:C4	2.40	0.51
1:X:24:LYS:CG	1:X:304:ILE:HG12	2.37	0.51
1:C:284:THR:HG22	1:C:286:ASP:N	2.23	0.51
1:C:24:LYS:CG	1:C:304:ILE:HG12	2.40	0.51
1:D:41:MET:HE3	1:D:54:THR:HB	1.93	0.51
1:D:170:VAL:HG22	1:D:200:VAL:HB	1.91	0.51
1:E:208:GLU:HG2	1:E:242:LEU:HD11	1.93	0.51
1:D:98:PRO:HB3	1:E:296:ILE:HD13	1.93	0.51
1:C:274:CYS:O	1:C:275:LEU:HB2	2.11	0.51
1:Y:157:THR:HG21	1:Y:298:GLU:HG3	1.92	0.51
1:D:98:PRO:CB	1:E:296:ILE:HD13	2.41	0.51
1:Y:70:ILE:HD11	1:Y:100:MET:HE1	1.92	0.51
1:Z:124:GLU:OE1	1:Z:144:ALA:HB3	2.11	0.51
1:E:125:VAL:O	1:E:129:GLN:HG3	2.10	0.51
1:X:49:ARG:HH21	1:X:276:PRO:HG3	1.76	0.51
1:C:285:ASP:HB3	1:E:81:ARG:HH12	1.75	0.50
1:X:83:VAL:HG12	1:X:84:ILE:N	2.25	0.50
1:X:240:ALA:HB3	1:X:245:ASN:O	2.10	0.50
1:Y:308:VAL:O	1:Y:312:ARG:HG2	2.11	0.50
1:X:23:ILE:HD13	1:X:307:THR:HG22	1.91	0.50
1:X:122:TYR:HA	1:X:145:THR:HG21	1.93	0.50
1:Z:118:ARG:NH2	1:Z:243:GLY:N	2.60	0.50
1:C:210:ASP:HB3	1:C:213:PHE:HD1	1.75	0.50
1:C:105:ASP:O	1:C:137:PRO:HD2	2.12	0.50
1:X:112:PHE:HA	4:X:369:HOH:O	2.10	0.50
1:C:124:GLU:HB3	1:C:127:ILE:HG21	1.92	0.50
1:C:59:LEU:HD13	1:E:64:ASN:CB	2.41	0.50
1:Y:208:GLU:CD	1:Y:208:GLU:H	2.14	0.50
1:X:70:ILE:HD13	1:X:77:LEU:HD11	1.92	0.50
1:C:168:LYS:HA	1:C:198:GLU:HB3	1.94	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:49:ARG:HH22	1:E:91:GLU:CD	2.15	0.50
1:D:208:GLU:HG3	1:D:242:LEU:HD11	1.94	0.50
1:E:212:LYS:HE3	1:E:213:PHE:CE1	2.47	0.50
1:E:118:ARG:HG3	1:E:246:TYR:CD1	2.47	0.50
1:X:184:VAL:HB	1:X:185:PRO:HD3	1.92	0.50
1:X:68:LEU:HD22	1:X:69:ASP:N	2.27	0.50
1:C:51:ARG:HD3	1:E:68:LEU:HD23	1.94	0.50
1:C:169:VAL:HG11	1:C:192:MET:CE	2.38	0.50
1:C:284:THR:HG22	1:C:285:ASP:N	2.27	0.50
1:X:127:ILE:O	1:X:131:ILE:HG13	2.12	0.50
1:Y:70:ILE:O	1:Y:71:ASN:CB	2.60	0.50
1:D:208:GLU:HG2	1:D:242:LEU:CD1	2.41	0.49
1:X:209:LEU:HD21	1:X:238:TRP:CE2	2.47	0.49
1:Y:162:LYS:HD2	1:Y:164:THR:O	2.11	0.49
1:X:95:GLU:O	1:X:99:VAL:HG23	2.12	0.49
1:D:89:LYS:N	1:D:89:LYS:HD2	2.26	0.49
1:C:228:GLU:HG3	1:C:265:VAL:CG1	2.41	0.49
1:Y:81:ARG:HH12	1:Z:285:ASP:HB3	1.78	0.49
1:Z:153:ALA:O	1:Z:157:THR:CG2	2.56	0.49
1:X:279:ARG:HD3	1:X:285:ASP:OD2	2.13	0.49
1:C:70:ILE:O	1:C:71:ASN:CB	2.60	0.49
1:Y:240:ALA:HB3	1:Y:245:ASN:O	2.11	0.49
1:Z:284:THR:HG22	1:Z:285:ASP:N	2.28	0.49
1:E:308:VAL:O	1:E:312:ARG:HG2	2.13	0.49
1:Y:95:GLU:O	1:Y:99:VAL:HG23	2.12	0.49
1:D:314:LEU:HD23	1:D:317:LEU:HD12	1.94	0.49
1:X:210:ASP:HB3	1:X:213:PHE:HD1	1.77	0.49
1:E:0:HIS:HA	1:E:316:ASN:HD21	1.75	0.49
1:E:0:HIS:HA	1:E:316:ASN:HD22	1.77	0.49
1:E:285:ASP:O	1:E:289:GLU:HG3	2.12	0.49
1:E:39:LEU:HD22	1:E:310:LEU:HD21	1.95	0.48
1:D:228:GLU:HA	1:D:265:VAL:HG12	1.95	0.48
1:X:203:HIS:HD2	1:X:204:PRO:O	1.96	0.48
1:D:4:PHE:CZ	1:D:148:PRO:HB2	2.48	0.48
1:C:97:ILE:CB	1:C:98:PRO:HD3	2.37	0.48
1:E:122:TYR:OH	1:E:182:GLN:HG3	2.13	0.48
1:E:132:GLN:HG2	4:E:362:HOH:O	2.13	0.48
1:Z:125:VAL:O	1:Z:129:GLN:HG3	2.13	0.48
1:Z:24:LYS:HZ3	1:Z:301:ASN:HD21	1.56	0.48
1:E:168:LYS:HA	1:E:198:GLU:HB3	1.95	0.48
1:E:1:MET:HE1	1:E:10:ILE:HD12	1.96	0.48
1:Z:24:LYS:HZ2	1:Z:301:ASN:ND2	2.10	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:X:273:HIS:O	1:X:275:LEU:HD13	2.12	0.48
1:D:70:ILE:HG22	1:D:71:ASN:N	2.29	0.48
1:E:51:ARG:O	1:E:55:GLN:HG3	2.14	0.48
1:D:35:ARG:HG2	1:D:36:ASN:ND2	2.29	0.48
1:Z:286:ASP:O	1:Z:290:SER:HB3	2.13	0.48
1:C:204:PRO:HG3	1:C:256:TRP:CE2	2.49	0.48
1:D:188:PHE:O	1:D:192:MET:HG2	2.13	0.48
1:E:70:ILE:O	1:E:71:ASN:CB	2.62	0.48
1:C:294:ILE:C	1:C:297:PRO:HD2	2.34	0.48
1:Z:228:GLU:HA	1:Z:265:VAL:HG12	1.95	0.48
1:E:208:GLU:CG	1:E:242:LEU:HD11	2.43	0.48
1:E:4:PHE:CE2	1:E:309:VAL:HG21	2.49	0.48
1:E:43:PHE:CZ	1:E:141:MET:HE1	2.49	0.48
1:D:82:GLY:HA2	1:E:285:ASP:OD1	2.14	0.48
1:D:202:THR:HG23	1:D:202:THR:O	2.13	0.48
1:X:103:TYR:O	1:Y:56:LYS:HD3	2.14	0.48
1:Z:209:LEU:HD21	1:Z:238:TRP:CE2	2.49	0.47
1:X:161:TYR:HB3	1:X:270:TYR:CZ	2.49	0.47
1:C:118:ARG:HA	1:C:246:TYR:CE2	2.50	0.47
1:D:216:ASN:ND2	4:D:372:HOH:O	2.46	0.47
1:D:83:VAL:HG12	1:D:84:ILE:N	2.28	0.47
1:X:51:ARG:NH2	1:X:69:ASP:OD1	2.47	0.47
1:Z:17:LEU:HD13	1:Z:159:GLU:HG3	1.96	0.47
1:Z:168:LYS:HA	1:Z:198:GLU:HB3	1.95	0.47
1:Z:23:ILE:HD13	1:Z:307:THR:HG22	1.97	0.47
1:X:95:GLU:C	1:X:98:PRO:HD2	2.34	0.47
1:Z:203:HIS:HD2	1:Z:204:PRO:O	1.97	0.47
1:X:70:ILE:HD11	1:X:100:MET:CE	2.45	0.47
1:C:4:PHE:CE2	1:C:309:VAL:HG21	2.49	0.47
1:Z:61:LEU:CD1	1:Z:310:LEU:HB3	2.44	0.47
1:D:24:LYS:NZ	1:D:301:ASN:HD21	2.12	0.47
1:C:59:LEU:CD1	1:E:64:ASN:HB3	2.45	0.47
1:X:127:ILE:HG22	4:X:397:HOH:O	2.15	0.47
1:X:204:PRO:HG3	1:X:256:TRP:CE2	2.49	0.47
1:Z:230:ALA:O	1:Z:269:ALA:HB2	2.15	0.47
1:E:50:THR:HG21	1:E:141:MET:HB3	1.97	0.47
1:Z:294:ILE:C	1:Z:297:PRO:HD2	2.34	0.47
1:C:49:ARG:NH2	1:C:276:PRO:HG3	2.26	0.47
1:E:7:VAL:O	1:E:10:ILE:HG22	2.15	0.47
1:Y:294:ILE:C	1:Y:297:PRO:HD2	2.36	0.47
1:D:296:ILE:HB	1:D:297:PRO:HD3	1.97	0.47
1:X:98:PRO:CB	1:Y:296:ILE:HD13	2.45	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:X:150:GLN:O	1:X:153:ALA:HB3	2.15	0.46
1:C:64:ASN:HB3	1:D:59:LEU:HD13	1.97	0.46
1:X:208:GLU:CG	1:X:242:LEU:HD11	2.46	0.46
1:E:154:ASP:HA	1:E:157:THR:CG2	2.45	0.46
1:E:192:MET:CE	1:E:197:TYR:HD1	2.28	0.46
1:Z:70:ILE:O	1:Z:71:ASN:CB	2.61	0.46
1:E:265:VAL:HG12	1:E:265:VAL:O	2.16	0.46
1:C:127:ILE:HG22	4:C:376:HOH:O	2.16	0.46
1:Y:177:PRO:HD3	1:Y:281:MET:CE	2.46	0.46
1:D:152:PHE:CE1	1:D:156:ILE:HD11	2.50	0.46
1:C:306:ALA:O	1:C:310:LEU:HB2	2.15	0.46
1:Y:70:ILE:HD11	1:Y:100:MET:CE	2.45	0.46
1:C:80:GLU:O	1:C:92:HIS:NE2	2.45	0.46
1:X:105:ASP:O	1:X:137:PRO:HD2	2.14	0.46
1:D:282:ILE:HG13	4:D:366:HOH:O	2.15	0.46
1:D:70:ILE:O	1:D:71:ASN:CB	2.64	0.46
1:D:26:ASP:CG	1:D:29:LYS:HB2	2.35	0.46
1:C:27:ARG:HB2	4:C:366:HOH:O	2.16	0.46
1:X:152:PHE:O	1:X:156:ILE:HG13	2.16	0.46
1:E:127:ILE:HG23	1:E:128:ASN:N	2.31	0.46
1:D:203:HIS:HD2	1:D:204:PRO:O	1.98	0.46
1:C:275:LEU:HD23	1:E:99:VAL:HG21	1.97	0.46
1:Z:273:HIS:O	1:Z:275:LEU:HD13	2.16	0.46
1:Z:4:PHE:CZ	1:Z:148:PRO:HB2	2.50	0.46
1:E:45:ASN:ND2	1:E:112:PHE:HB2	2.31	0.46
3:C:351:SN0:C4	1:E:90:PRO:HB3	2.46	0.46
1:E:240:ALA:HB2	1:E:250:LEU:HG	1.98	0.46
1:Z:227:PHE:CD2	1:Z:262:GLN:HB3	2.50	0.46
1:X:107:ILE:HD12	1:X:134:SER:HB2	1.98	0.46
1:Y:202:THR:HG23	1:Y:202:THR:O	2.15	0.46
1:X:149:LEU:HB2	4:X:364:HOH:O	2.16	0.46
1:D:116:GLU:HB2	4:D:357:HOH:O	2.15	0.45
1:Y:265:VAL:HG12	1:Y:265:VAL:O	2.16	0.45
1:X:64:ASN:HB3	1:Y:59:LEU:HD13	1.97	0.45
1:D:7:VAL:O	1:D:10:ILE:HG22	2.16	0.45
1:X:50:THR:HG21	1:X:141:MET:HB3	1.97	0.45
1:E:0:HIS:CE1	4:E:377:HOH:O	2.70	0.45
1:X:159:GLU:HG2	1:X:197:TYR:OH	2.16	0.45
1:D:162:LYS:HD2	1:D:164:THR:O	2.15	0.45
1:Y:89:LYS:HA	1:Y:90:PRO:HD3	1.84	0.45
1:D:43:PHE:CZ	1:D:141:MET:HE1	2.52	0.45
1:D:274:CYS:O	1:D:275:LEU:HB2	2.16	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:Z:88:ASP:HB3	1:Z:89:LYS:HD2	1.99	0.45
1:Y:166:ARG:HG2	1:Y:196:ASP:OD2	2.16	0.45
1:C:44:PHE:CZ	1:C:70:ILE:HD12	2.51	0.45
1:X:141:MET:O	1:X:142:GLU:CB	2.65	0.45
1:C:175:PRO:HG2	1:C:281:MET:O	2.16	0.45
1:Y:223:GLN:HB2	1:Y:256:TRP:CZ3	2.52	0.45
1:Z:118:ARG:HG3	1:Z:246:TYR:CD1	2.52	0.45
1:X:98:PRO:HB2	1:Y:296:ILE:HD13	1.99	0.45
1:E:95:GLU:O	1:E:98:PRO:HD2	2.17	0.45
1:Y:122:TYR:OH	1:Y:182:GLN:CG	2.64	0.45
1:X:208:GLU:HG2	1:X:242:LEU:HD11	1.98	0.45
1:Y:222:ASP:HB3	1:Y:225:LYS:HB2	1.98	0.45
1:C:311:LYS:O	1:C:315:GLU:HG3	2.17	0.45
1:Z:187:SER:O	1:Z:191:TRP:HD1	2.00	0.45
1:D:284:THR:HG22	1:D:286:ASP:H	1.82	0.44
1:C:294:ILE:O	1:C:297:PRO:HD2	2.18	0.44
1:Y:91:GLU:HG3	4:Y:366:HOH:O	2.17	0.44
1:Z:119:GLU:HG2	1:Z:123:ASN:HD21	1.82	0.44
1:X:23:ILE:HD12	1:X:308:VAL:HG22	1.97	0.44
1:D:294:ILE:C	1:D:297:PRO:HD2	2.38	0.44
1:Z:189:ALA:HA	1:Z:199:PHE:CE2	2.51	0.44
1:X:125:VAL:O	1:X:129:GLN:HG3	2.18	0.44
1:Y:24:LYS:HZ2	1:Y:301:ASN:ND2	2.15	0.44
1:D:184:VAL:CG1	1:D:236:LYS:HG2	2.47	0.44
1:C:49:ARG:HB2	2:C:352:SO4:O4	2.18	0.44
1:C:176:HIS:NE2	1:C:278:ARG:HD2	2.33	0.44
1:Z:124:GLU:O	1:Z:128:ASN:HB2	2.17	0.44
1:C:241:TYR:HD2	1:C:242:LEU:HD23	1.80	0.44
1:C:228:GLU:HA	1:C:265:VAL:HG12	2.00	0.44
1:C:228:GLU:HG3	1:C:265:VAL:HG11	1.98	0.44
1:Y:230:ALA:O	1:Y:269:ALA:HB2	2.17	0.44
1:Y:118:ARG:HA	1:Y:246:TYR:CE2	2.53	0.44
1:E:192:MET:HE3	1:E:197:TYR:HD1	1.81	0.44
1:Y:78:GLU:HG3	1:Y:85:MET:HG2	1.99	0.44
1:E:41:MET:HE3	1:E:54:THR:CB	2.42	0.44
1:Y:192:MET:CE	1:Y:197:TYR:CD1	3.00	0.44
1:C:276:PRO:HB3	1:E:91:GLU:HB2	1.99	0.44
1:X:275:LEU:HD12	1:X:277:VAL:HG22	2.00	0.44
1:Z:225:LYS:HB3	1:Z:225:LYS:HE2	1.86	0.44
1:D:267:ASN:O	1:D:268:ASN:C	2.55	0.44
1:Z:122:TYR:HA	1:Z:145:THR:HG21	1.98	0.44
1:C:49:ARG:CZ	1:C:275:LEU:HB3	2.47	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:68:LEU:HD23	1:D:51:ARG:HD3	1.99	0.44
1:Y:34:GLY:O	1:Y:37:LYS:HB2	2.16	0.44
1:Z:212:LYS:HE3	1:Z:213:PHE:CZ	2.52	0.44
1:Z:105:ASP:O	1:Z:137:PRO:HD2	2.17	0.44
1:X:-3:ARG:O	1:X:-1:SER:N	2.43	0.44
1:X:70:ILE:HG22	1:X:71:ASN:N	2.32	0.44
1:Y:275:LEU:HD12	1:Y:277:VAL:CG2	2.47	0.44
1:C:4:PHE:CE2	1:C:148:PRO:HB2	2.53	0.44
1:C:88:ASP:HB3	1:C:89:LYS:HD2	1.98	0.44
1:E:-2:GLY:O	1:E:312:ARG:HD2	2.18	0.44
1:E:51:ARG:NH2	1:E:69:ASP:OD2	2.51	0.44
1:Z:51:ARG:O	1:Z:55:GLN:HG3	2.17	0.44
1:D:175:PRO:HG3	1:D:254:ARG:HA	2.00	0.44
1:Z:95:GLU:C	1:Z:98:PRO:HD2	2.38	0.44
1:D:170:VAL:HG21	1:D:226:ALA:O	2.18	0.44
1:X:176:HIS:CG	1:X:177:PRO:HD2	2.53	0.44
1:Y:154:ASP:HA	1:Y:157:THR:CG2	2.47	0.43
1:Y:182:GLN:HE21	1:Y:182:GLN:HB3	1.54	0.43
1:Z:89:LYS:CD	1:Z:89:LYS:N	2.81	0.43
1:Z:7:VAL:O	1:Z:10:ILE:HG22	2.18	0.43
1:E:179:PRO:HA	4:E:381:HOH:O	2.18	0.43
1:X:-3:ARG:HH12	1:X:8:GLN:HB3	1.79	0.43
1:Y:51:ARG:NH2	1:Y:69:ASP:OD2	2.51	0.43
1:C:162:LYS:HD2	1:C:164:THR:O	2.18	0.43
1:E:83:VAL:HG12	1:E:84:ILE:N	2.33	0.43
1:D:284:THR:HG22	1:D:285:ASP:N	2.33	0.43
1:C:89:LYS:HA	1:C:90:PRO:HD3	1.70	0.43
1:Y:245:ASN:HA	1:Y:248:GLN:HG3	1.99	0.43
1:D:45:ASN:HD21	1:D:112:PHE:HB2	1.82	0.43
1:E:44:PHE:HB2	1:E:110:ARG:O	2.18	0.43
1:D:124:GLU:HB3	1:D:127:ILE:HG23	1.96	0.43
1:C:41:MET:HG2	1:C:141:MET:HE3	2.00	0.43
1:X:279:ARG:NE	1:Z:95:GLU:OE2	2.43	0.43
1:D:310:LEU:HD12	1:D:310:LEU:HA	1.82	0.43
1:D:272:MET:O	1:D:273:HIS:HB2	2.17	0.43
1:Y:83:VAL:HG12	1:Y:84:ILE:N	2.34	0.43
1:D:61:LEU:HD12	1:D:310:LEU:HB3	2.01	0.43
1:Z:17:LEU:O	1:Z:21:PHE:HD1	2.02	0.43
1:D:227:PHE:CD2	1:D:262:GLN:HB3	2.53	0.43
1:E:119:GLU:HG2	1:E:123:ASN:ND2	2.34	0.43
1:Y:152:PHE:O	1:Y:156:ILE:HG13	2.19	0.43
1:C:310:LEU:HA	1:C:310:LEU:HD12	1.89	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:Z:70:ILE:HG22	1:Z:71:ASN:N	2.32	0.43
1:Z:284:THR:HG22	1:Z:286:ASP:H	1.82	0.43
1:Y:89:LYS:HD2	1:Y:89:LYS:N	2.34	0.43
1:Y:103:TYR:CE2	1:Z:53:SER:HB3	2.53	0.43
1:Y:240:ALA:HB2	1:Y:250:LEU:HG	2.00	0.43
1:Z:250:LEU:O	1:D:252:THR:HG23	2.18	0.43
1:C:43:PHE:CZ	1:C:141:MET:HE1	2.53	0.43
1:D:93:LEU:O	1:D:97:ILE:HG13	2.19	0.43
1:D:61:LEU:CD1	1:D:310:LEU:HB3	2.49	0.43
1:C:212:LYS:HE3	1:C:213:PHE:CE1	2.54	0.43
1:Y:294:ILE:O	1:Y:297:PRO:HD2	2.19	0.43
1:E:174:ALA:HB1	1:E:282:ILE:HA	1.99	0.43
1:Z:245:ASN:O	1:Z:248:GLN:HB2	2.19	0.43
1:C:154:ASP:O	1:C:157:THR:HG23	2.19	0.42
1:Z:185:PRO:HG2	1:Z:238:TRP:HB3	2.01	0.42
1:E:310:LEU:HD12	1:E:310:LEU:HA	1.81	0.42
1:C:186:ASN:HB3	1:C:213:PHE:CG	2.54	0.42
1:E:202:THR:HG23	1:E:202:THR:O	2.19	0.42
1:E:227:PHE:O	1:E:266:THR:HG22	2.19	0.42
1:C:70:ILE:HG22	1:C:71:ASN:N	2.34	0.42
1:C:89:LYS:HD2	1:C:89:LYS:N	2.34	0.42
1:X:70:ILE:O	1:X:71:ASN:CB	2.67	0.42
1:E:182:GLN:HB3	1:E:182:GLN:HE21	1.66	0.42
1:D:206:GLY:N	1:D:208:GLU:OE2	2.52	0.42
1:C:176:HIS:ND1	1:C:177:PRO:HD2	2.34	0.42
1:X:41:MET:HB3	1:X:41:MET:HE2	1.92	0.42
1:E:141:MET:O	1:E:142:GLU:CB	2.67	0.42
1:D:93:LEU:HG	1:D:97:ILE:CD1	2.50	0.42
1:X:228:GLU:HA	1:X:265:VAL:HG12	2.01	0.42
1:Y:98:PRO:CB	1:Z:296:ILE:HD13	2.50	0.42
1:Y:98:PRO:HB3	1:Z:296:ILE:HD13	2.01	0.42
1:Y:176:HIS:CG	1:Y:177:PRO:HD2	2.55	0.42
1:E:204:PRO:HG3	1:E:256:TRP:CE2	2.55	0.42
1:C:97:ILE:HB	1:C:98:PRO:CD	2.39	0.42
1:X:122:TYR:OH	1:X:182:GLN:CG	2.66	0.42
1:C:95:GLU:CD	1:D:279:ARG:HH21	2.22	0.42
1:Z:150:GLN:OE1	1:Z:302:ARG:NH1	2.53	0.42
1:C:156:ILE:O	1:C:160:GLU:HB2	2.19	0.42
1:X:83:VAL:CG1	1:X:84:ILE:N	2.83	0.42
1:C:265:VAL:HG12	1:C:265:VAL:O	2.18	0.42
1:Z:170:VAL:HG22	1:Z:200:VAL:HB	2.01	0.42
1:D:176:HIS:ND1	1:D:177:PRO:HD2	2.35	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:Z:285:ASP:O	1:Z:288:ILE:HG22	2.20	0.42
1:C:182:GLN:HE21	1:C:182:GLN:HB3	1.60	0.42
1:D:21:PHE:O	1:D:25:LYS:HE2	2.19	0.42
1:E:166:ARG:HG2	1:E:196:ASP:OD2	2.20	0.42
1:E:17:LEU:O	1:E:20:SER:HB2	2.20	0.42
1:Z:296:ILE:HB	1:Z:297:PRO:HD3	2.02	0.42
1:X:223:GLN:HB2	1:X:256:TRP:CZ3	2.55	0.42
1:C:118:ARG:NH2	1:C:243:GLY:N	2.68	0.42
1:X:119:GLU:HG2	1:X:123:ASN:ND2	2.35	0.42
1:Y:64:ASN:HB3	1:Z:59:LEU:HD13	2.02	0.42
1:X:173:TRP:CG	1:X:174:ALA:N	2.88	0.42
1:C:78:GLU:HG3	1:C:85:MET:HG2	2.02	0.41
1:D:254:ARG:HD3	4:D:379:HOH:O	2.20	0.41
1:E:-3:ARG:HH12	1:E:9:ASP:HA	1.86	0.41
1:Z:184:VAL:N	1:Z:185:PRO:CD	2.83	0.41
1:X:59:LEU:HD13	1:Z:64:ASN:HB3	2.01	0.41
1:X:157:THR:HB	1:X:301:ASN:OD1	2.19	0.41
1:C:61:LEU:HD12	1:C:310:LEU:HB3	2.03	0.41
1:Y:228:GLU:HG3	1:Y:265:VAL:CG1	2.49	0.41
1:X:78:GLU:HB3	1:X:92:HIS:CD2	2.56	0.41
1:C:31:VAL:O	1:C:62:GLY:HA3	2.19	0.41
1:D:151:SER:O	1:D:154:ASP:HB2	2.21	0.41
1:C:141:MET:O	1:C:142:GLU:CB	2.67	0.41
1:X:98:PRO:HB2	1:Y:296:ILE:CD1	2.51	0.41
1:C:24:LYS:HZ2	1:C:301:ASN:HD21	1.68	0.41
1:Y:224:MET:HE3	1:Y:265:VAL:HG21	2.03	0.41
1:E:89:LYS:CD	1:E:89:LYS:N	2.84	0.41
1:Y:95:GLU:HG2	4:Z:384:HOH:O	2.19	0.41
1:C:122:TYR:CD2	1:C:145:THR:HG21	2.56	0.41
1:Z:306:ALA:O	1:Z:310:LEU:HB2	2.19	0.41
1:Z:191:TRP:O	1:Z:195:THR:HG23	2.20	0.41
1:E:30:TYR:C	1:E:32:GLU:H	2.23	0.41
1:Y:208:GLU:HG2	1:Y:242:LEU:CD1	2.48	0.41
1:D:49:ARG:CZ	1:D:275:LEU:HB3	2.51	0.41
1:X:90:PRO:HB3	3:Y:347:SN0:C4	2.50	0.41
1:Y:4:PHE:CE2	1:Y:148:PRO:HB2	2.56	0.41
1:D:154:ASP:HA	1:D:157:THR:CG2	2.50	0.41
1:E:296:ILE:N	1:E:297:PRO:CD	2.83	0.41
1:Y:241:TYR:CD2	1:Y:242:LEU:HD23	2.55	0.41
1:Y:19:GLU:OE2	1:Y:312:ARG:NE	2.44	0.41
1:E:225:LYS:HB3	1:E:225:LYS:HE2	1.82	0.41
1:X:-3:ARG:C	1:X:-1:SER:H	2.22	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:Z:70:ILE:HD11	1:Z:100:MET:HE1	2.02	0.41
1:Z:162:LYS:HD2	1:Z:164:THR:O	2.20	0.41
1:E:274:CYS:O	1:E:275:LEU:O	2.38	0.41
1:D:171:MET:HG3	1:D:188:PHE:CD2	2.56	0.41
1:C:50:THR:HG22	1:C:141:MET:HE2	2.03	0.41
1:Z:294:ILE:O	1:Z:297:PRO:HD2	2.21	0.41
1:C:176:HIS:CG	1:C:177:PRO:HD2	2.56	0.41
1:Y:103:TYR:O	1:Z:56:LYS:HD3	2.21	0.41
1:X:231:ASP:HB3	4:X:403:HOH:O	2.21	0.41
1:Z:304:ILE:HD12	1:Z:304:ILE:N	2.36	0.41
1:D:89:LYS:HA	1:D:90:PRO:HD3	1.90	0.41
1:D:17:LEU:O	1:D:21:PHE:HD1	2.03	0.41
1:E:31:VAL:HG22	1:E:31:VAL:O	2.21	0.41
1:Y:70:ILE:HG22	1:Y:71:ASN:N	2.36	0.40
1:Y:235:ALA:O	1:Y:273:HIS:CD2	2.75	0.40
1:D:114:ARG:HB2	4:D:357:HOH:O	2.21	0.40
1:E:41:MET:CE	1:E:141:MET:HE3	2.50	0.40
1:D:41:MET:CE	1:D:141:MET:HE1	2.51	0.40
1:D:174:ALA:HB1	1:D:282:ILE:HA	2.03	0.40
1:E:76:LYS:O	1:E:90:PRO:HD2	2.21	0.40
1:C:227:PHE:O	1:C:266:THR:HG22	2.22	0.40
1:C:73:GLY:O	1:D:51:ARG:NH1	2.53	0.40
1:D:90:PRO:HB3	3:E:355:SN0:H32	2.03	0.40
1:X:176:HIS:NE2	1:X:278:ARG:HD2	2.36	0.40
1:Y:103:TYR:CZ	1:Z:53:SER:HB3	2.56	0.40
1:Y:92:HIS:CE1	1:Y:94:LEU:HG	2.56	0.40
1:Z:317:LEU:O	1:Z:318:PRO:O	2.39	0.40
1:C:259:GLY:HA2	1:C:284:THR:OG1	2.22	0.40
1:C:124:GLU:OE1	1:C:144:ALA:HB3	2.22	0.40
1:X:42:ILE:HD13	1:X:100:MET:SD	2.61	0.40
1:Y:141:MET:O	1:Y:142:GLU:CB	2.69	0.40
1:C:45:ASN:HD21	1:C:112:PHE:HB2	1.86	0.40
1:E:24:LYS:NZ	1:E:301:ASN:HD21	2.20	0.40
1:D:42:ILE:HD13	1:D:100:MET:HE1	2.03	0.40
1:D:168:LYS:HD3	1:D:230:ALA:HA	2.02	0.40
1:C:118:ARG:NH2	1:C:243:GLY:H	2.19	0.40
3:C:351:SN0:OD2	1:E:90:PRO:HB3	2.21	0.40
1:D:103:TYR:CZ	1:E:53:SER:HB3	2.57	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	C	319/338 (94%)	290 (91%)	26 (8%)	3 (1%)	25	63
1	D	320/338 (95%)	289 (90%)	25 (8%)	6 (2%)	12	37
1	E	320/338 (95%)	292 (91%)	24 (8%)	4 (1%)	18	51
1	X	320/338 (95%)	290 (91%)	27 (8%)	3 (1%)	25	63
1	Y	320/338 (95%)	298 (93%)	20 (6%)	2 (1%)	33	72
1	Z	318/338 (94%)	287 (90%)	28 (9%)	3 (1%)	25	63
All	All	1917/2028 (94%)	1746 (91%)	150 (8%)	21 (1%)	21	57

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	X	142	GLU
1	Y	142	GLU
1	Z	142	GLU
1	C	142	GLU
1	D	142	GLU
1	E	142	GLU
1	X	243	GLY
1	Z	240	ALA
1	Z	243	GLY
1	C	243	GLY
1	E	243	GLY
1	D	243	GLY
1	C	71	ASN
1	D	240	ALA
1	D	273	HIS
1	X	85	MET
1	D	31	VAL
1	Y	243	GLY
1	E	275	LEU
1	D	258	VAL
1	E	31	VAL

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	C	279/294 (95%)	265 (95%)	14 (5%)	34	70
1	D	280/294 (95%)	266 (95%)	14 (5%)	34	70
1	E	280/294 (95%)	270 (96%)	10 (4%)	47	82
1	X	280/294 (95%)	266 (95%)	14 (5%)	34	70
1	Y	280/294 (95%)	265 (95%)	15 (5%)	31	66
1	Z	279/294 (95%)	269 (96%)	10 (4%)	47	82
All	All	1678/1764 (95%)	1601 (95%)	77 (5%)	37	73

All (77) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	X	25	LYS
1	X	31	VAL
1	X	52	LEU
1	X	68	LEU
1	X	78	GLU
1	X	81	ARG
1	X	89	LYS
1	X	147	HIS
1	X	157	THR
1	X	163	LYS
1	X	182	GLN
1	X	277	VAL
1	X	295	VAL
1	X	310	LEU
1	Y	-3	ARG
1	Y	25	LYS
1	Y	52	LEU
1	Y	81	ARG
1	Y	89	LYS
1	Y	118	ARG
1	Y	147	HIS
1	Y	152	PHE
1	Y	157	THR

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Mol	Chain	Res	Type
1	Y	163	LYS
1	Y	178	ARG
1	Y	182	GLN
1	Y	251	SER
1	Y	277	VAL
1	Y	310	LEU
1	Z	25	LYS
1	Z	48	LEU
1	Z	52	LEU
1	Z	132	GLN
1	Z	147	HIS
1	Z	152	PHE
1	Z	157	THR
1	Z	163	LYS
1	Z	182	GLN
1	Z	310	LEU
1	C	0	HIS
1	C	25	LYS
1	C	48	LEU
1	C	52	LEU
1	C	81	ARG
1	C	89	LYS
1	C	95	GLU
1	C	147	HIS
1	C	157	THR
1	C	166	ARG
1	C	178	ARG
1	C	182	GLN
1	C	208	GLU
1	C	295	VAL
1	D	-3	ARG
1	D	25	LYS
1	D	48	LEU
1	D	52	LEU
1	D	68	LEU
1	D	81	ARG
1	D	95	GLU
1	D	152	PHE
1	D	157	THR
1	D	178	ARG
1	D	182	GLN
1	D	208	GLU

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Mol	Chain	Res	Type
1	D	267	ASN
1	D	310	LEU
1	E	-3	ARG
1	E	25	LYS
1	E	51	ARG
1	E	52	LEU
1	E	132	GLN
1	E	163	LYS
1	E	182	GLN
1	E	208	GLU
1	E	268	ASN
1	E	295	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (49) such sidechains are listed below:

Mol	Chain	Res	Type
1	X	0	HIS
1	X	64	ASN
1	X	182	GLN
1	X	203	HIS
1	X	267	ASN
1	X	268	ASN
1	X	301	ASN
1	X	316	ASN
1	Y	64	ASN
1	Y	182	GLN
1	Y	203	HIS
1	Y	248	GLN
1	Y	267	ASN
1	Y	268	ASN
1	Y	301	ASN
1	Z	64	ASN
1	Z	123	ASN
1	Z	182	GLN
1	Z	203	HIS
1	Z	255	ASN
1	Z	267	ASN
1	Z	268	ASN
1	Z	301	ASN
1	Z	316	ASN
1	C	64	ASN
1	C	123	ASN

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Mol	Chain	Res	Type
1	C	182	GLN
1	C	203	HIS
1	C	267	ASN
1	C	268	ASN
1	C	301	ASN
1	C	316	ASN
1	D	8	GLN
1	D	64	ASN
1	D	182	GLN
1	D	203	HIS
1	D	267	ASN
1	D	268	ASN
1	D	292	GLN
1	D	301	ASN
1	D	316	ASN
1	E	0	HIS
1	E	182	GLN
1	E	203	HIS
1	E	267	ASN
1	E	268	ASN
1	E	292	GLN
1	E	301	ASN
1	E	316	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 ⓘ

14 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	SN0	C	351	-	14,14,14	0.69	0	17,17,17	1.14	2 (11%)
2	SO4	C	352	-	4,4,4	0.18	0	6,6,6	0.09	0
2	SO4	C	358	-	4,4,4	0.32	0	6,6,6	0.16	0
3	SN0	D	353	-	14,14,14	0.77	0	17,17,17	1.18	2 (11%)
2	SO4	D	354	-	4,4,4	0.37	0	6,6,6	0.28	0
3	SN0	E	355	-	14,14,14	0.67	0	17,17,17	0.93	0
2	SO4	E	356	-	4,4,4	0.21	0	6,6,6	0.19	0
3	SN0	X	345	-	14,14,14	0.64	0	17,17,17	1.26	2 (11%)
2	SO4	X	346	-	4,4,4	0.21	0	6,6,6	0.19	0
2	SO4	X	357	-	4,4,4	0.11	0	6,6,6	0.08	0
3	SN0	Y	347	-	14,14,14	0.68	0	17,17,17	1.16	3 (17%)
2	SO4	Y	348	-	4,4,4	0.24	0	6,6,6	0.25	0
3	SN0	Z	349	-	14,14,14	0.85	0	17,17,17	1.31	1 (5%)
2	SO4	Z	350	-	4,4,4	0.33	0	6,6,6	0.14	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	SN0	C	351	-	-	0/16/16/16	0/0/0/0
2	SO4	C	352	-	-	0/0/0/0	0/0/0/0
2	SO4	C	358	-	-	0/0/0/0	0/0/0/0
3	SN0	D	353	-	-	0/16/16/16	0/0/0/0
2	SO4	D	354	-	-	0/0/0/0	0/0/0/0
3	SN0	E	355	-	-	0/16/16/16	0/0/0/0
2	SO4	E	356	-	-	0/0/0/0	0/0/0/0
3	SN0	X	345	-	-	0/16/16/16	0/0/0/0
2	SO4	X	346	-	-	0/0/0/0	0/0/0/0
2	SO4	X	357	-	-	0/0/0/0	0/0/0/0
3	SN0	Y	347	-	-	0/16/16/16	0/0/0/0
2	SO4	Y	348	-	-	0/0/0/0	0/0/0/0
3	SN0	Z	349	-	-	0/16/16/16	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	Z	350	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Z	349	SN0	C-CA-N1	3.99	120.82	110.53
3	D	353	SN0	CB-CA-N1	3.21	116.95	110.83
3	C	351	SN0	CB-CA-N1	2.64	115.86	110.83
3	Y	347	SN0	OD2-C4-C3	-2.52	114.35	123.03
3	X	345	SN0	CB-CA-N1	2.47	115.53	110.83
3	Y	347	SN0	OD1-C4-C3	2.30	122.34	114.22
3	D	353	SN0	C-CA-N1	2.25	116.32	110.53
3	C	351	SN0	C-CA-N1	2.14	116.05	110.53
3	Y	347	SN0	C-CA-N1	2.07	115.88	110.53
3	X	345	SN0	C2-C3-C4	-2.01	109.80	113.53

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	C	321/338 (94%)	0.11	16 (4%) 28 28	54, 72, 92, 108	0
1	D	322/338 (95%)	0.18	16 (4%) 28 28	46, 71, 96, 117	0
1	E	322/338 (95%)	-0.16	9 (2%) 50 52	38, 58, 78, 101	0
1	X	322/338 (95%)	-0.41	3 (0%) 81 81	28, 48, 71, 84	0
1	Y	322/338 (95%)	-0.64	0 100 100	22, 35, 51, 78	0
1	Z	320/338 (94%)	-0.45	5 (1%) 68 69	27, 44, 62, 82	0
All	All	1929/2028 (95%)	-0.23	49 (2%) 54 55	22, 55, 86, 117	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	-3	ARG	5.0
1	C	318	PRO	4.7
1	D	-2	GLY	4.5
1	C	-1	SER	4.3
1	D	25	LYS	4.3
1	C	254	ARG	4.2
1	D	318	PRO	3.8
1	E	-2	GLY	3.7
1	D	166	ARG	3.5
1	C	116	GLU	3.5
1	Z	-1	SER	3.5
1	C	-2	GLY	3.4
1	D	0	HIS	3.4
1	Z	318	PRO	3.3
1	C	166	ARG	3.1
1	E	-3	ARG	3.1
1	X	205	GLU	3.0
1	E	318	PRO	3.0
1	C	0	HIS	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	254	ARG	2.9
1	D	267	ASN	2.8
1	C	172	THR	2.8
1	E	291	PRO	2.7
1	D	14	LYS	2.7
1	E	0	HIS	2.7
1	C	26	ASP	2.6
1	X	244	ASP	2.6
1	E	163	LYS	2.6
1	C	242	LEU	2.5
1	Z	0	HIS	2.5
1	C	25	LYS	2.5
1	E	71	ASN	2.5
1	D	163	LYS	2.4
1	C	309	VAL	2.4
1	D	2	LYS	2.4
1	C	205	GLU	2.4
1	D	-1	SER	2.4
1	X	116	GLU	2.3
1	C	253	ASP	2.3
1	E	-1	SER	2.3
1	Z	117	ASN	2.2
1	C	14	LYS	2.2
1	D	8	GLN	2.2
1	C	81	ARG	2.2
1	D	196	ASP	2.1
1	D	216	ASN	2.1
1	D	165	ALA	2.1
1	E	260	ASP	2.0
1	Z	116	GLU	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	SN0	D	353	15/15	0.21	1.14	41,52,57,57	0
3	SN0	X	345	15/15	0.20	1.13	46,51,54,56	0
2	SO4	Y	348	5/5	0.14	0.43	32,35,35,36	0
3	SN0	E	355	15/15	0.16	0.23	46,51,54,56	0
3	SN0	C	351	15/15	0.17	0.04	68,69,71,71	0
3	SN0	Y	347	15/15	0.13	-0.05	30,35,41,41	0
2	SO4	Z	350	5/5	0.14	-0.10	32,36,38,38	0
2	SO4	D	354	5/5	0.18	-0.17	57,60,62,63	0
2	SO4	E	356	5/5	0.14	-0.30	49,50,53,55	0
3	SN0	Z	349	15/15	0.12	-0.87	37,39,41,41	0
2	SO4	C	352	5/5	0.12	-1.51	59,60,60,62	0
2	SO4	X	346	5/5	0.10	-1.60	42,44,44,44	0
2	SO4	X	357	5/5	0.09	-3.08	53,55,56,58	0
2	SO4	C	358	5/5	0.17	-3.24	83,83,85,85	0

6.5 Other polymers ⓘ

There are no such residues in this entry.