



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:31 AM UTC

PDB ID : 3TFZ / pdb_00003tfz
Title : Crystal structure of Zhui aromatase/cyclase from Streptomyces sp. R1128
Authors : Ames, B.D.; Lee, M.Y.; Moody, C.; Zhang, W.; Tang, Y.; Wong, S.K.; Tsai, S.C.
Deposited on : 2011-08-16
Resolution : 2.39 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

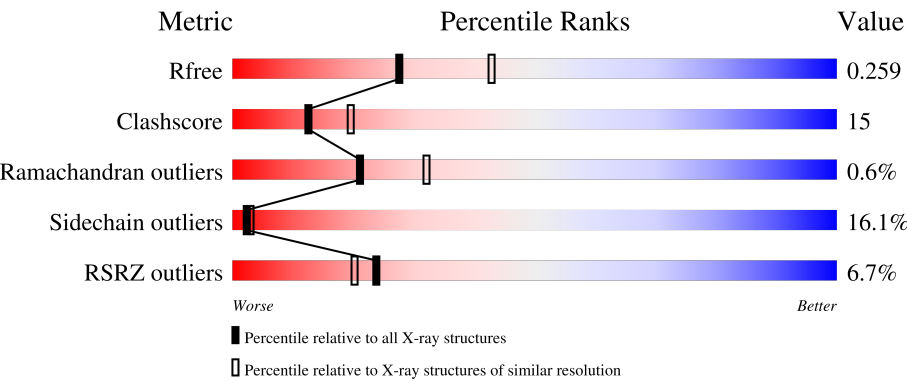
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry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

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.39 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	172	63%27%6%.
1	B	172	3%67%24%...
1	C	172	6%51%31%11%.5%
1	D	172	16%49%32%9%5%5%
1	E	172	6%62%23%10%.

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Mol	Chain	Length	Quality of chain
1	F	172	5% 60% 26% 6% 6%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7945 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Cyclase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	168	Total	C	N	O	S	0	0	0
			1300	811	237	249	3			
1	B	165	Total	C	N	O	S	0	0	0
			1282	800	234	245	3			
1	C	164	Total	C	N	O	S	0	0	0
			1275	795	233	244	3			
1	D	164	Total	C	N	O	S	0	0	0
			1275	795	233	244	3			
1	E	165	Total	C	N	O	S	0	0	0
			1281	798	234	246	3			
1	F	161	Total	C	N	O	S	0	0	0
			1262	788	230	241	3			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q9F6D3
A	-1	SER	-	expression tag	UNP Q9F6D3
A	0	HIS	-	expression tag	UNP Q9F6D3
B	-2	GLY	-	expression tag	UNP Q9F6D3
B	-1	SER	-	expression tag	UNP Q9F6D3
B	0	HIS	-	expression tag	UNP Q9F6D3
C	-2	GLY	-	expression tag	UNP Q9F6D3
C	-1	SER	-	expression tag	UNP Q9F6D3
C	0	HIS	-	expression tag	UNP Q9F6D3
D	-2	GLY	-	expression tag	UNP Q9F6D3
D	-1	SER	-	expression tag	UNP Q9F6D3
D	0	HIS	-	expression tag	UNP Q9F6D3
E	-2	GLY	-	expression tag	UNP Q9F6D3
E	-1	SER	-	expression tag	UNP Q9F6D3
E	0	HIS	-	expression tag	UNP Q9F6D3
F	-2	GLY	-	expression tag	UNP Q9F6D3
F	-1	SER	-	expression tag	UNP Q9F6D3

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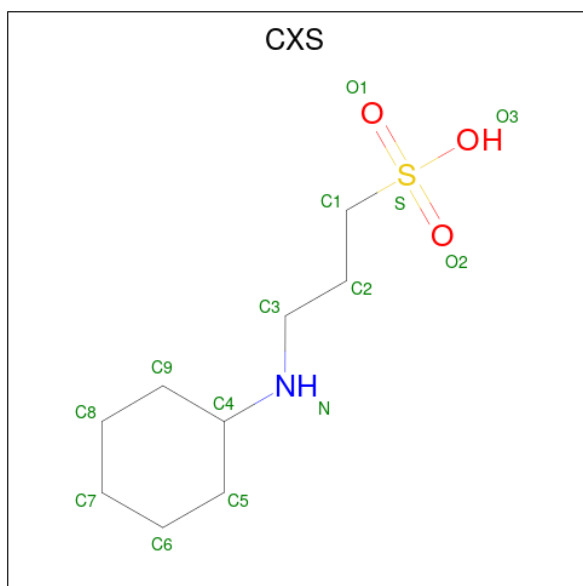
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Chain	Residue	Modelled	Actual	Comment	Reference
F	0	HIS	-	expression tag	UNP Q9F6D3

- Molecule 2 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total K 1 1	0	0
2	B	1	Total K 1 1	0	0
2	E	1	Total K 1 1	0	0
2	F	1	Total K 1 1	0	0

- Molecule 3 is 3-CYCLOHEXYL-1-PROPYLSULFONIC ACID (CCD ID: CXS) (formula: C₉H₁₉NO₃S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O S 14 9 1 3 1	0	0
3	B	1	Total C N O S 14 9 1 3 1	0	0
3	C	1	Total C N O S 14 9 1 3 1	0	0
3	D	1	Total C N O S 14 9 1 3 1	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	D	1	Total	C	N	O	S	0	0
			14	9	1	3	1		
3	E	1	Total	C	N	O	S	0	0
			14	9	1	3	1		
3	F	1	Total	C	N	O	S	0	0
			14	9	1	3	1		

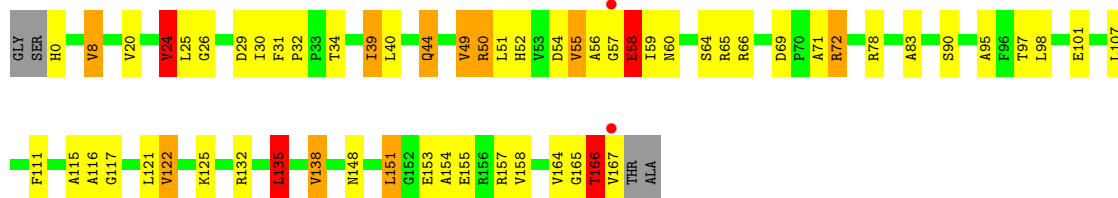
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	64	Total	O	0	0
			64	64		
4	B	36	Total	O	0	0
			36	36		
4	C	17	Total	O	0	0
			17	17		
4	D	17	Total	O	0	0
			17	17		
4	E	21	Total	O	0	0
			21	21		
4	F	13	Total	O	0	0
			13	13		

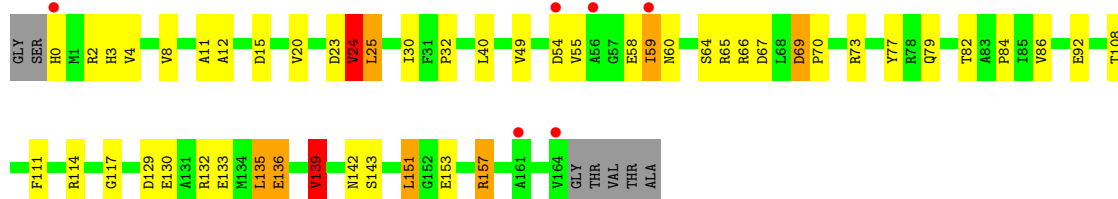
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

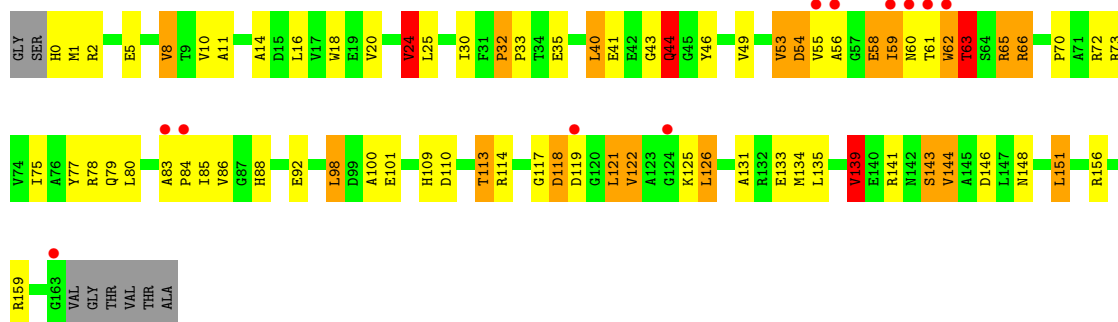
• Molecule 1: Cyclase



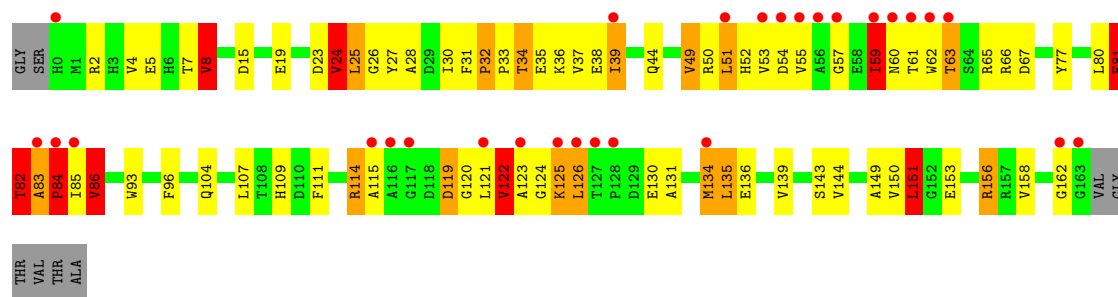
• Molecule 1: Cyclase



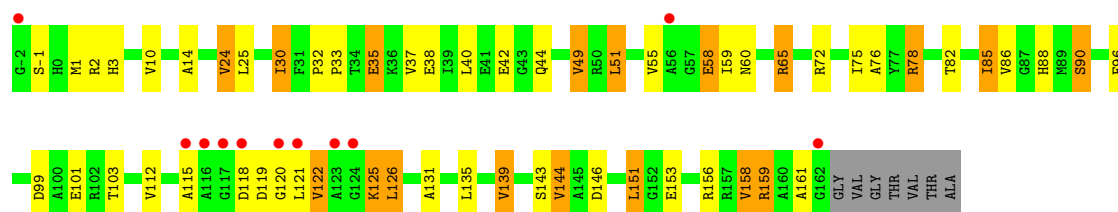
• Molecule 1: Cyclase



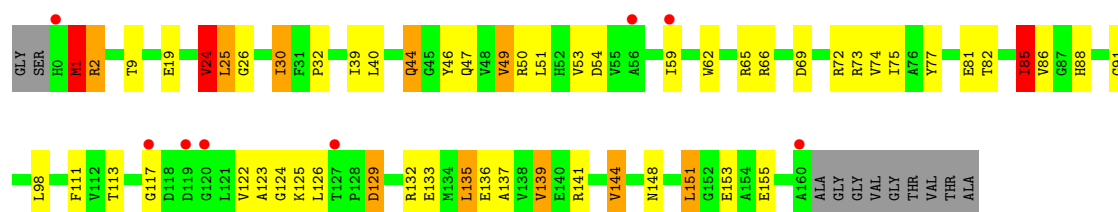
• Molecule 1: Cyclase



• Molecule 1: Cyclase



• Molecule 1: Cyclase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	50.02Å 135.73Å 175.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.39 50.00 – 2.39	Depositor EDS
% Data completeness (in resolution range)	99.3 (50.00-2.39) 99.4 (50.00-2.39)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.52 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.193 , 0.261 0.192 , 0.259	Depositor DCC
R_{free} test set	2430 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	38.4	Xtriage
Anisotropy	0.324	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 46.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7945	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.96% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CXS, K

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	1.86	23/1323 (1.7%)	1.68	24/1803 (1.3%)
1	B	1.74	10/1305 (0.8%)	1.52	15/1778 (0.8%)
1	C	1.52	10/1298 (0.8%)	1.52	21/1768 (1.2%)
1	D	1.52	10/1298 (0.8%)	1.50	16/1768 (0.9%)
1	E	1.59	10/1304 (0.8%)	1.54	20/1776 (1.1%)
1	F	1.48	7/1285 (0.5%)	1.49	14/1751 (0.8%)
All	All	1.63	70/7813 (0.9%)	1.54	110/10644 (1.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	2

All (70) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	76	ALA	CA-CB	11.84	1.69	1.53
1	D	86	VAL	C-O	11.34	1.36	1.24
1	B	59	ILE	CA-CB	8.77	1.65	1.54
1	B	11	ALA	CA-CB	8.72	1.63	1.52
1	A	26	GLY	C-O	-8.57	1.13	1.23
1	F	9	THR	CA-CB	8.04	1.64	1.53
1	B	23	ASP	CA-C	7.77	1.62	1.53
1	C	11	ALA	CA-CB	7.72	1.62	1.52
1	A	49	VAL	CA-CB	7.22	1.64	1.54
1	C	14	ALA	CA-CB	7.04	1.64	1.53
1	E	51	LEU	CA-C	-6.84	1.44	1.52
1	C	20	VAL	CA-CB	6.67	1.62	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	20	VAL	CA-C	6.65	1.61	1.52
1	C	32	PRO	CA-C	6.61	1.58	1.52
1	B	3	HIS	CA-C	6.41	1.60	1.52
1	B	32	PRO	CA-C	6.31	1.58	1.52
1	A	155	GLU	CA-C	6.29	1.60	1.52
1	F	74	VAL	CA-CB	6.25	1.62	1.54
1	B	92	GLU	C-O	-6.24	1.16	1.23
1	A	154	ALA	CA-CB	6.22	1.63	1.53
1	A	122	VAL	CA-CB	6.17	1.61	1.54
1	F	32	PRO	CA-C	6.11	1.57	1.52
1	D	86	VAL	C-N	6.08	1.42	1.33
1	E	10	VAL	CA-CB	6.06	1.61	1.54
1	B	4	VAL	C-O	-6.05	1.18	1.24
1	C	66	ARG	CA-C	5.98	1.59	1.52
1	F	49	VAL	CA-CB	5.94	1.63	1.55
1	A	20	VAL	CA-C	5.89	1.60	1.52
1	A	116	ALA	CA-CB	5.86	1.63	1.53
1	A	39	ILE	CA-C	-5.85	1.45	1.52
1	C	72	ARG	CZ-NH1	5.78	1.40	1.32
1	E	139	VAL	CA-CB	5.78	1.62	1.54
1	A	101	GLU	CA-C	-5.76	1.44	1.52
1	F	139	VAL	CA-CB	5.74	1.61	1.54
1	C	63	THR	CA-CB	5.70	1.61	1.53
1	D	124	GLY	C-O	5.68	1.31	1.23
1	E	30	ILE	CA-CB	5.68	1.61	1.54
1	A	95	ALA	C-O	-5.61	1.17	1.24
1	F	85	ILE	CA-CB	5.61	1.61	1.54
1	A	31	PHE	CA-C	-5.58	1.45	1.52
1	A	115	ALA	CA-C	5.57	1.59	1.52
1	A	121	LEU	C-O	5.54	1.30	1.23
1	C	122	VAL	CA-CB	5.47	1.62	1.54
1	B	86	VAL	CA-CB	5.42	1.60	1.54
1	D	126	LEU	C-O	5.41	1.30	1.23
1	A	164	VAL	CA-C	5.40	1.58	1.52
1	A	90	SER	C-O	5.39	1.30	1.23
1	F	47	GLN	C-O	-5.34	1.17	1.23
1	E	78	ARG	CD-NE	-5.34	1.38	1.46
1	A	72	ARG	CA-C	5.32	1.59	1.52
1	B	108	THR	N-CA	5.30	1.52	1.45
1	D	7	THR	CA-CB	5.28	1.63	1.53
1	A	107	LEU	CA-C	5.28	1.58	1.52
1	A	78	ARG	NE-CZ	5.27	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	49	VAL	C-O	-5.25	1.18	1.24
1	A	148	ASN	CA-C	5.23	1.59	1.52
1	E	146	ASP	CA-C	5.19	1.59	1.52
1	A	135	LEU	C-O	-5.17	1.18	1.24
1	A	158	VAL	CA-CB	5.13	1.60	1.54
1	A	97	THR	C-O	5.11	1.30	1.24
1	C	53	VAL	CA-C	5.11	1.59	1.52
1	E	99	ASP	CA-C	-5.10	1.47	1.53
1	D	5	GLU	CA-C	-5.08	1.46	1.52
1	D	4	VAL	N-CA	5.05	1.51	1.46
1	A	64	SER	CA-C	-5.04	1.46	1.52
1	D	32	PRO	CA-C	5.03	1.56	1.52
1	C	144	VAL	CA-CB	5.03	1.60	1.54
1	D	150	VAL	CA-CB	5.02	1.60	1.54
1	E	3	HIS	N-CA	-5.01	1.40	1.46
1	D	149	ALA	CA-C	5.00	1.59	1.52

All (110) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	125	LYS	CB-CA-C	-10.42	95.06	110.06
1	A	58	GLU	N-CA-C	10.05	122.10	108.23
1	A	72	ARG	NE-CZ-NH2	-9.86	110.32	119.20
1	C	24	VAL	CB-CA-C	-9.42	99.70	112.04
1	D	25	LEU	N-CA-C	-9.00	98.63	110.43
1	D	125	LYS	N-CA-C	8.98	127.38	113.28
1	E	144	VAL	N-CA-CB	8.65	119.70	110.62
1	B	24	VAL	CB-CA-C	-8.49	99.01	112.16
1	E	78	ARG	NE-CZ-NH2	-8.28	111.75	119.20
1	A	72	ARG	NE-CZ-NH1	8.10	129.60	121.50
1	F	24	VAL	CB-CA-C	-7.83	100.03	112.16
1	E	24	VAL	CB-CA-C	-7.66	100.29	112.16
1	D	126	LEU	N-CA-CB	7.64	124.81	111.13
1	E	144	VAL	CB-CA-C	7.60	121.42	111.70
1	D	24	VAL	N-CA-C	7.45	117.57	110.42
1	F	47	GLN	CB-CA-C	7.31	120.91	110.24
1	C	118	ASP	N-CA-C	-7.28	102.80	114.09
1	F	1	MET	N-CA-C	-7.18	97.23	108.14
1	A	72	ARG	CD-NE-CZ	7.14	134.39	124.40
1	B	69	ASP	CA-C-N	7.06	127.66	119.47
1	B	69	ASP	C-N-CA	7.06	127.66	119.47
1	C	54	ASP	N-CA-C	6.94	120.80	109.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	59	ILE	N-CA-C	6.80	117.23	108.12
1	C	148	ASN	N-CA-C	6.78	119.50	111.71
1	C	32	PRO	N-CA-C	6.62	118.78	110.70
1	D	115	ALA	N-CA-C	6.62	118.40	110.19
1	C	30	ILE	N-CA-C	6.60	119.56	112.96
1	E	65	ARG	NE-CZ-NH2	-6.57	113.29	119.20
1	F	148	ASN	N-CA-C	6.53	118.39	111.28
1	B	82	THR	N-CA-C	6.48	120.07	110.48
1	A	24	VAL	CB-CA-C	-6.46	102.62	112.05
1	E	49	VAL	CB-CA-C	6.41	120.36	110.50
1	A	166	THR	CB-CA-C	-6.35	99.81	114.41
1	A	72	ARG	CB-CA-C	6.27	122.42	109.95
1	B	59	ILE	CB-CA-C	6.26	119.29	111.15
1	B	12	ALA	N-CA-C	-6.24	98.11	108.23
1	A	132	ARG	NE-CZ-NH2	-6.21	113.61	119.20
1	A	32	PRO	CA-C-N	-6.20	113.38	119.64
1	A	32	PRO	C-N-CA	-6.20	113.38	119.64
1	B	20	VAL	N-CA-C	6.16	117.64	110.62
1	D	151	LEU	N-CA-C	6.10	117.60	111.07
1	A	30	ILE	CA-CB-CG2	6.04	120.77	110.50
1	A	8	VAL	CG1-CB-CG2	6.00	123.99	110.80
1	A	122	VAL	CB-CA-C	5.99	119.18	110.98
1	A	34	THR	N-CA-C	5.92	117.68	109.15
1	C	86	VAL	N-CA-C	5.89	117.07	108.58
1	E	146	ASP	CA-C-N	5.88	128.09	120.44
1	E	146	ASP	C-N-CA	5.88	128.09	120.44
1	E	118	ASP	CB-CA-C	-5.88	101.59	110.06
1	A	165	GLY	N-CA-C	5.87	118.23	111.36
1	A	138	VAL	N-CA-C	-5.84	105.39	111.58
1	F	117	GLY	N-CA-C	5.82	126.97	113.18
1	E	151	LEU	CA-CB-CG	5.82	136.65	116.30
1	C	139	VAL	N-CA-CB	5.78	119.25	110.58
1	C	146	ASP	N-CA-C	5.75	118.01	111.11
1	E	78	ARG	CB-CA-C	-5.72	102.33	110.34
1	E	78	ARG	CD-NE-CZ	5.71	132.39	124.40
1	C	24	VAL	N-CA-CB	5.71	117.60	110.47
1	B	139	VAL	N-CA-C	5.68	116.41	110.62
1	D	8	VAL	N-CA-CB	-5.66	100.86	111.62
1	A	71	ALA	N-CA-C	-5.65	104.74	111.69
1	C	65	ARG	CG-CD-NE	-5.62	99.65	112.00
1	A	132	ARG	N-CA-C	5.61	117.08	111.07
1	C	62	TRP	N-CA-C	5.58	117.55	109.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	132	ARG	CG-CD-NE	-5.57	99.75	112.00
1	E	24	VAL	CA-CB-CG2	5.55	119.84	110.40
1	B	151	LEU	CA-CB-CG	5.54	135.68	116.30
1	E	65	ARG	NE-CZ-NH1	5.53	127.03	121.50
1	A	122	VAL	N-CA-CB	5.52	117.67	111.21
1	F	24	VAL	N-CA-C	5.51	117.94	111.05
1	D	124	GLY	N-CA-C	-5.49	100.17	113.18
1	C	24	VAL	N-CA-C	5.47	116.86	110.62
1	C	61	THR	N-CA-C	5.47	117.82	108.90
1	E	115	ALA	N-CA-C	5.47	118.28	110.24
1	E	90	SER	N-CA-CB	-5.45	103.11	111.56
1	E	161	ALA	N-CA-C	5.43	119.66	113.19
1	A	117	GLY	N-CA-C	-5.43	104.83	112.60
1	A	8	VAL	CA-CB-CG1	5.40	119.59	110.40
1	E	118	ASP	N-CA-C	5.39	121.75	113.28
1	D	24	VAL	N-CA-CB	5.39	116.86	110.55
1	B	32	PRO	N-CA-C	5.37	117.25	110.70
1	C	121	LEU	N-CA-C	5.37	117.56	109.23
1	B	117	GLY	N-CA-C	-5.34	104.40	112.41
1	F	53	VAL	CA-C-N	-5.34	114.66	122.19
1	F	53	VAL	C-N-CA	-5.34	114.66	122.19
1	A	151	LEU	CA-CB-CG	5.33	134.94	116.30
1	C	44	GLN	CA-C-N	-5.32	110.98	121.41
1	C	44	GLN	C-N-CA	-5.32	110.98	121.41
1	C	126	LEU	N-CA-C	5.32	117.80	110.35
1	D	81	GLU	CA-C-N	5.31	131.26	121.70
1	D	81	GLU	C-N-CA	5.31	131.26	121.70
1	F	139	VAL	N-CA-CB	5.29	118.51	110.58
1	A	98	LEU	N-CA-C	-5.28	105.66	112.68
1	C	41	GLU	N-CA-C	5.26	116.51	108.52
1	F	46	TYR	N-CA-C	5.24	116.81	108.79
1	D	57	GLY	N-CA-C	-5.23	107.79	114.85
1	C	98	LEU	N-CA-C	-5.22	106.59	113.12
1	F	91	GLY	N-CA-C	5.20	117.03	110.38
1	B	30	ILE	N-CA-C	5.18	120.12	109.34
1	E	143	SER	CB-CA-C	-5.17	102.71	110.92
1	C	101	GLU	N-CA-C	5.16	119.12	112.41
1	B	136	GLU	N-CA-C	5.16	116.90	111.28
1	F	144	VAL	CB-CA-C	5.12	119.05	112.14
1	D	122	VAL	N-CA-C	5.10	115.32	110.53
1	E	24	VAL	CG1-CB-CG2	5.10	122.01	110.80
1	B	55	VAL	N-CA-C	-5.08	98.74	106.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	19	GLU	N-CA-C	-5.05	105.77	111.28
1	B	136	GLU	CA-CB-CG	-5.03	104.05	114.10
1	F	54	ASP	N-CA-CB	5.03	118.45	110.16
1	D	151	LEU	CA-CB-CG	5.02	133.87	116.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	123	ALA	Peptide
1	D	34	THR	Peptide

5.2 Too-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 hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1300	0	1282	30	0
1	B	1282	0	1263	15	0
1	C	1275	0	1254	49	0
1	D	1275	0	1254	87	0
1	E	1281	0	1259	36	0
1	F	1262	0	1243	32	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	14	0	19	1	0
3	B	14	0	19	1	0
3	C	14	0	19	2	0
3	D	28	0	38	2	0
3	E	14	0	19	1	0
3	F	14	0	18	2	0
4	A	64	0	0	2	0
4	B	36	0	0	1	0
4	C	17	0	0	0	0
4	D	17	0	0	0	0
4	E	21	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	F	13	0	0	3	0
All	All	7945	0	7687	236	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 15.

All (236) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:81:GLU:CA	1:D:82:THR:HG23	1.36	1.48
1:C:1:MET:HE1	1:C:88:HIS:CD2	1.68	1.26
1:D:55:VAL:CG2	1:D:60:ASN:HB2	1.65	1.25
1:D:81:GLU:N	1:D:82:THR:HG23	1.53	1.22
1:D:81:GLU:HA	1:D:82:THR:CG2	1.76	1.15
1:D:55:VAL:HG21	1:D:60:ASN:HB2	1.26	1.14
1:D:81:GLU:O	1:D:81:GLU:HG2	1.49	1.10
1:E:159:ARG:HH11	1:E:159:ARG:HG2	1.16	1.08
1:D:81:GLU:CA	1:D:82:THR:CG2	2.29	1.07
1:D:84:PRO:HG3	1:D:114:ARG:CB	1.89	1.01
1:A:166:THR:HG22	1:A:167:VAL:H	1.26	1.01
1:C:63:THR:HG22	1:C:80:LEU:HB2	1.41	0.99
1:E:159:ARG:HG2	1:E:159:ARG:NH1	1.70	0.97
1:D:81:GLU:HA	1:D:82:THR:HG23	1.00	0.96
1:D:65:ARG:NH1	1:D:67:ASP:OD2	2.00	0.95
1:E:159:ARG:HH11	1:E:159:ARG:CG	1.79	0.95
1:D:81:GLU:N	1:D:82:THR:CG2	2.30	0.94
1:D:84:PRO:HA	1:D:86:VAL:N	1.84	0.93
1:D:55:VAL:CG2	1:D:60:ASN:CB	2.47	0.91
1:C:1:MET:CE	1:C:88:HIS:CD2	2.54	0.91
1:F:85:ILE:HD12	1:F:122:VAL:HG12	1.51	0.90
1:D:84:PRO:HG3	1:D:114:ARG:HB2	1.52	0.90
1:E:119:ASP:N	1:E:120:GLY:HA2	1.90	0.86
1:D:84:PRO:HA	1:D:86:VAL:H	1.38	0.86
1:D:80:LEU:C	1:D:82:THR:CG2	2.57	0.78
1:C:83:ALA:HB1	1:C:84:PRO:CD	2.13	0.78
1:D:53:VAL:HG12	1:D:54:ASP:H	1.47	0.78
1:D:55:VAL:HG21	1:D:60:ASN:CB	2.11	0.77
1:D:80:LEU:C	1:D:82:THR:HG23	2.09	0.77
1:D:84:PRO:HD3	1:D:114:ARG:HG3	1.66	0.77
1:D:81:GLU:O	1:D:81:GLU:CG	2.30	0.76
1:C:10:VAL:HG22	1:C:151:LEU:HB2	1.67	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:83:ALA:HB1	1:C:84:PRO:HD3	1.68	0.74
1:A:50:ARG:HG2	1:A:50:ARG:HH11	1.53	0.73
1:D:28:ALA:HB2	1:D:37:VAL:HG23	1.69	0.73
1:D:63:THR:HG22	1:D:80:LEU:HB2	1.71	0.73
1:D:34:THR:HG21	1:D:51:LEU:HD23	1.71	0.72
1:E:88:HIS:NE2	1:E:90:SER:HB2	2.04	0.72
1:D:2:ARG:HB3	1:D:2:ARG:NH1	2.04	0.70
1:D:2:ARG:HB3	1:D:2:ARG:HH11	1.56	0.70
1:F:85:ILE:CD1	1:F:122:VAL:HG12	2.21	0.70
1:D:119:ASP:N	1:D:120:GLY:HA2	2.05	0.70
1:D:36:LYS:HB3	1:D:52:HIS:CD2	2.27	0.69
1:D:81:GLU:HA	1:D:82:THR:CB	2.22	0.69
1:F:137:ALA:O	1:F:141:ARG:HG3	1.94	0.68
1:F:26:GLY:O	1:F:30:ILE:HG23	1.94	0.67
1:A:44:GLN:CD	1:A:44:GLN:H	2.02	0.67
1:D:84:PRO:HG3	1:D:114:ARG:CG	2.25	0.66
1:D:55:VAL:HG23	1:D:60:ASN:HB2	1.73	0.66
1:D:139:VAL:O	1:D:143:SER:HB2	1.95	0.66
1:A:50:ARG:HG2	1:A:50:ARG:NH1	2.09	0.65
1:D:84:PRO:CD	1:D:114:ARG:HG3	2.25	0.65
1:D:82:THR:OG1	1:D:83:ALA:N	2.30	0.64
1:D:49:VAL:HG13	1:D:51:LEU:HD12	1.80	0.64
1:D:122:VAL:HG11	1:D:131:ALA:HB2	1.78	0.64
1:C:24:VAL:HG13	1:C:66:ARG:CZ	2.28	0.63
1:D:24:VAL:O	1:D:27:TYR:HB2	1.98	0.63
1:D:34:THR:CG2	1:D:51:LEU:HD23	2.28	0.63
1:B:111:PHE:CD2	1:B:135:LEU:HD13	2.34	0.62
1:A:153:GLU:OE2	1:E:78:ARG:NH2	2.33	0.62
1:C:63:THR:CG2	1:C:80:LEU:HB2	2.24	0.62
1:D:125:LYS:O	1:D:126:LEU:HD23	2.00	0.62
1:E:119:ASP:H	1:E:120:GLY:HA2	1.65	0.61
1:E:126:LEU:HD23	1:E:131:ALA:HA	1.82	0.61
1:F:123:ALA:O	1:F:125:LYS:HG2	2.01	0.61
1:B:111:PHE:HD2	1:B:135:LEU:HD13	1.66	0.61
1:C:55:VAL:O	1:C:58:GLU:HG2	2.01	0.60
1:C:98:LEU:HD13	1:D:96:PHE:HZ	1.65	0.60
1:D:30:ILE:O	1:D:32:PRO:HD3	2.01	0.60
1:E:58:GLU:HG2	1:E:60:ASN:HD21	1.66	0.60
1:C:139:VAL:O	1:C:143:SER:HB2	2.02	0.60
1:B:64:SER:HB3	1:B:79:GLN:HA	1.84	0.59
1:E:30:ILE:O	1:E:32:PRO:HD3	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:ARG:HH11	1:A:50:ARG:CG	2.14	0.59
1:E:159:ARG:NH1	1:E:159:ARG:CG	2.44	0.59
1:F:129:ASP:HB2	4:F:185:HOH:O	2.03	0.59
1:D:84:PRO:CG	1:D:114:ARG:HG3	2.33	0.58
1:A:153:GLU:CD	1:E:78:ARG:HH22	2.12	0.58
1:E:85:ILE:HG22	1:E:122:VAL:HG22	1.86	0.58
1:F:141:ARG:NH2	4:F:175:HOH:O	2.36	0.58
1:B:15:ASP:OD1	1:B:73:ARG:NH1	2.36	0.58
1:D:84:PRO:HG3	1:D:114:ARG:HG3	1.85	0.58
1:D:84:PRO:HA	1:D:85:ILE:C	2.27	0.58
1:D:130:GLU:O	1:D:134:MET:HB2	2.04	0.57
1:C:62:TRP:HB2	1:C:79:GLN:OE1	2.04	0.57
1:A:166:THR:HG22	1:A:167:VAL:N	2.08	0.57
1:E:153:GLU:OE2	1:E:156:ARG:NH1	2.37	0.57
1:E:72:ARG:NH1	1:F:98:LEU:O	2.38	0.56
1:C:8:VAL:HG13	1:C:151:LEU:HD23	1.87	0.56
1:D:59:ILE:HG12	1:D:59:ILE:O	2.05	0.56
1:F:24:VAL:HG13	1:F:66:ARG:CZ	2.35	0.56
1:D:55:VAL:HG22	1:D:60:ASN:CB	2.34	0.56
1:A:50:ARG:HD2	1:A:51:LEU:N	2.20	0.56
1:C:0:HIS:HB3	1:E:119:ASP:O	2.06	0.55
1:D:125:LYS:O	1:D:125:LYS:CG	2.54	0.55
1:F:85:ILE:HD12	1:F:122:VAL:CG1	2.32	0.55
1:C:54:ASP:O	1:C:141:ARG:NH2	2.40	0.55
1:D:51:LEU:O	1:D:61:THR:HA	2.07	0.55
1:B:65:ARG:NH1	1:B:67:ASP:OD1	2.39	0.55
1:A:44:GLN:H	1:A:44:GLN:NE2	2.05	0.54
1:D:85:ILE:O	1:D:85:ILE:HG22	2.07	0.54
1:E:35:GLU:OE1	1:E:59:ILE:HD11	2.08	0.53
1:F:151:LEU:HD13	1:F:155:GLU:CD	2.33	0.53
1:A:50:ARG:NH1	4:A:215:HOH:O	2.39	0.53
1:C:32:PRO:HD2	1:C:33:PRO:HD2	1.90	0.53
1:E:125:LYS:HD3	1:E:126:LEU:HD13	1.91	0.53
1:B:135:LEU:O	1:B:139:VAL:HG13	2.10	0.52
1:D:122:VAL:HG13	1:D:126:LEU:HB2	1.91	0.52
1:D:55:VAL:HG21	1:D:60:ASN:HD22	1.73	0.52
1:B:129:ASP:O	1:B:133:GLU:HG2	2.10	0.51
1:D:85:ILE:O	1:D:85:ILE:CG2	2.57	0.51
1:C:63:THR:HG23	1:C:80:LEU:HD12	1.93	0.51
1:D:84:PRO:CA	1:D:86:VAL:H	2.17	0.51
1:C:53:VAL:HB	1:C:62:TRP:CH2	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:98:LEU:HD13	1:D:96:PHE:CZ	2.46	0.50
1:E:85:ILE:HG22	1:E:122:VAL:CG2	2.41	0.50
1:D:158:VAL:O	1:D:162:GLY:N	2.45	0.50
1:A:52:HIS:HB3	1:A:59:ILE:HD11	1.92	0.50
1:D:82:THR:O	1:D:83:ALA:O	2.30	0.50
1:F:111:PHE:HD2	1:F:135:LEU:HD13	1.76	0.50
1:C:43:GLY:HA3	1:C:46:TYR:CE1	2.47	0.50
1:C:156:ARG:HG2	1:C:159:ARG:HH22	1.77	0.50
1:D:50:ARG:HA	1:D:62:TRP:O	2.12	0.49
1:A:29:ASP:HB3	1:E:78:ARG:HD2	1.93	0.49
1:B:25:LEU:HB2	4:B:203:HOH:O	2.12	0.49
1:C:32:PRO:CD	1:C:33:PRO:HD2	2.42	0.49
1:C:59:ILE:HG13	1:C:59:ILE:O	2.11	0.49
1:C:109:HIS:NE2	3:C:170:CXS:H4	2.28	0.49
1:B:24:VAL:HG13	1:B:66:ARG:CZ	2.42	0.49
1:F:1:MET:SD	1:F:88:HIS:HD2	2.35	0.49
1:C:40:LEU:N	1:C:40:LEU:CD1	2.76	0.49
1:C:18:TRP:CE2	1:C:70:PRO:HB3	2.48	0.49
1:F:113:THR:HG21	1:F:132:ARG:HA	1.93	0.48
1:A:111:PHE:HD2	1:A:135:LEU:HD13	1.78	0.48
1:C:65:ARG:O	1:C:77:TYR:HA	2.13	0.48
1:D:34:THR:HG21	1:D:51:LEU:CD2	2.42	0.48
1:D:34:THR:HG22	1:D:35:GLU:H	1.79	0.48
1:A:153:GLU:O	1:A:157:ARG:HG3	2.13	0.48
1:B:132:ARG:O	1:B:136:GLU:HG3	2.13	0.48
1:D:122:VAL:HG11	1:D:131:ALA:CB	2.43	0.48
1:E:82:THR:HB	1:E:86:VAL:HB	1.96	0.48
1:A:65:ARG:NH1	1:C:78:ARG:NH2	2.62	0.48
1:A:111:PHE:CD2	1:A:135:LEU:HD13	2.49	0.48
1:D:119:ASP:HB2	1:D:121:LEU:H	1.79	0.48
1:F:62:TRP:HB2	1:F:81:GLU:O	2.14	0.48
1:F:132:ARG:O	1:F:136:GLU:HG3	2.13	0.48
1:D:8:VAL:HG13	1:D:151:LEU:HD23	1.96	0.47
1:A:56:ALA:O	1:A:57:GLY:C	2.56	0.47
1:D:27:TYR:O	1:D:31:PHE:HB2	2.14	0.47
1:C:56:ALA:HB3	1:C:134:MET:CE	2.44	0.47
1:C:156:ARG:HD2	1:C:156:ARG:O	2.15	0.47
1:D:125:LYS:O	1:D:125:LYS:HG2	2.15	0.47
1:F:44:GLN:HE21	1:F:44:GLN:HA	1.79	0.47
1:F:85:ILE:CD1	1:F:122:VAL:CG1	2.92	0.47
1:C:43:GLY:O	1:C:44:GLN:C	2.58	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:83:ALA:CB	1:C:84:PRO:CD	2.85	0.47
1:D:63:THR:CG2	1:D:80:LEU:HB2	2.42	0.47
1:E:119:ASP:N	1:E:120:GLY:CA	2.72	0.47
1:F:73:ARG:HD3	4:F:174:HOH:O	2.14	0.47
1:F:82:THR:HB	1:F:86:VAL:HB	1.97	0.47
1:C:0:HIS:CB	1:E:119:ASP:O	2.62	0.47
1:A:60:ASN:OD1	1:A:83:ALA:HB1	2.14	0.47
1:B:153:GLU:O	1:B:157:ARG:HG2	2.14	0.47
1:F:25:LEU:CD2	1:F:39:ILE:HD12	2.45	0.46
1:D:23:ASP:O	1:D:25:LEU:O	2.33	0.46
1:B:84:PRO:O	1:B:114:ARG:NE	2.46	0.46
1:E:14:ALA:HA	1:E:103:THR:OG1	2.15	0.46
1:F:69:ASP:OD1	1:F:72:ARG:HG3	2.15	0.46
1:C:56:ALA:HB3	1:C:134:MET:HE2	1.98	0.46
1:F:2:ARG:HD3	1:F:136:GLU:OE2	2.15	0.46
1:A:65:ARG:NH1	1:C:78:ARG:HH22	2.13	0.45
1:A:65:ARG:HH12	1:C:78:ARG:NH2	2.14	0.45
1:D:2:ARG:HH11	1:D:2:ARG:CB	2.28	0.45
1:D:153:GLU:OE2	1:D:156:ARG:NH1	2.49	0.45
1:F:1:MET:SD	1:F:88:HIS:CD2	3.09	0.45
1:C:125:LYS:HG3	1:C:126:LEU:H	1.82	0.45
1:E:35:GLU:HG3	1:E:59:ILE:HD13	1.98	0.45
1:C:75:ILE:O	1:C:92:GLU:HA	2.16	0.45
1:D:25:LEU:HD11	1:D:39:ILE:HG13	1.98	0.45
1:C:8:VAL:CG1	1:C:151:LEU:HD23	2.46	0.44
1:D:49:VAL:HG13	1:D:51:LEU:CD1	2.46	0.44
1:D:84:PRO:HG3	1:D:114:ARG:HB3	1.87	0.44
1:D:84:PRO:HB3	1:D:85:ILE:HA	1.99	0.44
1:D:53:VAL:HG12	1:D:54:ASP:N	2.26	0.44
1:D:122:VAL:CG1	1:D:126:LEU:HB2	2.46	0.44
1:E:125:LYS:HB2	1:E:126:LEU:HD13	2.00	0.44
1:D:15:ASP:O	1:D:19:GLU:HG2	2.17	0.44
1:E:37:VAL:HG22	1:E:51:LEU:HD23	1.99	0.44
1:E:58:GLU:CG	1:E:60:ASN:HD21	2.29	0.44
1:E:88:HIS:HB3	1:E:112:VAL:HG12	1.99	0.44
1:F:113:THR:HG22	1:F:135:LEU:HD12	1.99	0.44
1:A:39:ILE:CG2	1:E:-1:SER:HB3	2.48	0.43
1:D:93:TRP:CD2	1:D:107:LEU:HD13	2.53	0.43
1:D:62:TRP:HE1	3:D:170:CXS:H31	1.82	0.43
1:D:139:VAL:O	1:D:143:SER:CB	2.64	0.43
1:F:123:ALA:C	1:F:125:LYS:H	2.26	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:ASP:HA	1:B:70:PRO:HD2	1.77	0.43
1:C:2:ARG:O	1:C:110:ASP:HA	2.18	0.43
1:A:59:ILE:HG23	1:A:59:ILE:O	2.19	0.43
1:D:26:GLY:O	1:D:30:ILE:HG12	2.18	0.43
1:D:33:PRO:HB2	1:D:53:VAL:CG1	2.48	0.43
1:D:65:ARG:O	1:D:77:TYR:HA	2.19	0.43
1:D:111:PHE:HD2	1:D:135:LEU:HD13	1.83	0.43
1:A:39:ILE:HG21	1:E:-1:SER:HB3	2.01	0.43
1:C:113:THR:HG21	1:C:131:ALA:HB1	2.00	0.43
1:D:85:ILE:HD13	1:D:126:LEU:HD12	2.01	0.43
1:A:24:VAL:HG13	1:A:66:ARG:CZ	2.49	0.42
1:C:119:ASP:HB3	1:C:121:LEU:HG	2.02	0.42
1:D:24:VAL:HG13	1:D:66:ARG:CZ	2.50	0.42
1:C:119:ASP:CB	1:C:121:LEU:HG	2.49	0.42
1:E:32:PRO:HB2	1:E:33:PRO:HD3	2.01	0.42
1:F:77:TYR:CZ	3:F:171:CXS:H21	2.55	0.42
1:C:1:MET:HE1	1:C:88:HIS:NE2	2.26	0.42
1:F:111:PHE:CD2	1:F:135:LEU:HD13	2.55	0.42
1:B:129:ASP:OD1	1:B:132:ARG:NH1	2.53	0.42
1:F:153:GLU:OE1	1:F:153:GLU:HA	2.20	0.41
1:B:77:TYR:HB2	3:B:171:CXS:O1	2.21	0.41
3:F:171:CXS:H31	3:F:171:CXS:H91	1.78	0.41
1:A:0:HIS:HB3	4:A:222:HOH:O	2.21	0.41
1:A:65:ARG:HH12	1:C:78:ARG:HH22	1.68	0.41
1:C:40:LEU:N	1:C:40:LEU:HD12	2.35	0.41
1:C:156:ARG:HD2	1:C:156:ARG:C	2.45	0.41
1:E:51:LEU:HD13	3:E:171:CXS:O3	2.20	0.41
1:D:109:HIS:NE2	3:D:170:CXS:H51	2.36	0.41
1:E:75:ILE:HD13	1:E:75:ILE:HG21	1.77	0.41
1:E:158:VAL:HG12	1:E:159:ARG:N	2.35	0.41
1:C:1:MET:HE1	1:C:88:HIS:CG	2.40	0.41
1:F:25:LEU:HD22	1:F:39:ILE:CD1	2.51	0.40
1:A:50:ARG:HD2	1:A:51:LEU:H	1.85	0.40
1:A:55:VAL:N	1:A:58:GLU:O	2.50	0.40
1:A:69:ASP:CG	1:A:72:ARG:HB3	2.46	0.40
3:A:171:CXS:H31	3:A:171:CXS:H52	1.97	0.40
1:E:96:PHE:HZ	1:F:98:LEU:HD13	1.85	0.40
1:D:96:PHE:HB2	1:D:104:GLN:HB3	2.04	0.40
1:C:66:ARG:NH1	3:C:170:CXS:O1	2.53	0.40
1:F:123:ALA:C	1:F:125:LYS:N	2.80	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	166/172 (96%)	159 (96%)	7 (4%)	0	100	100
1	B	163/172 (95%)	157 (96%)	6 (4%)	0	100	100
1	C	162/172 (94%)	148 (91%)	12 (7%)	2 (1%)	10	16
1	D	162/172 (94%)	139 (86%)	20 (12%)	3 (2%)	6	8
1	E	163/172 (95%)	153 (94%)	10 (6%)	0	100	100
1	F	159/172 (92%)	154 (97%)	4 (2%)	1 (1%)	21	32
All	All	975/1032 (94%)	910 (93%)	59 (6%)	6 (1%)	21	32

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	100	ALA
1	D	82	THR
1	F	124	GLY
1	D	83	ALA
1	C	117	GLY
1	D	84	PRO

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	134/136 (98%)	118 (88%)	16 (12%)	5	7
1	B	132/136 (97%)	114 (86%)	18 (14%)	3	5

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	131/136 (96%)	106 (81%)	25 (19%)	1	2
1	D	131/136 (96%)	109 (83%)	22 (17%)	2	3
1	E	132/136 (97%)	107 (81%)	25 (19%)	1	2
1	F	131/136 (96%)	110 (84%)	21 (16%)	2	3
All	All	791/816 (97%)	664 (84%)	127 (16%)	2	3

All (127) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	VAL
1	A	24	VAL
1	A	25	LEU
1	A	40	LEU
1	A	44	GLN
1	A	49	VAL
1	A	50	ARG
1	A	54	ASP
1	A	55	VAL
1	A	58	GLU
1	A	122	VAL
1	A	125	LYS
1	A	135	LEU
1	A	138	VAL
1	A	151	LEU
1	A	166	THR
1	B	0	HIS
1	B	2	ARG
1	B	8	VAL
1	B	24	VAL
1	B	25	LEU
1	B	40	LEU
1	B	49	VAL
1	B	54	ASP
1	B	58	GLU
1	B	59	ILE
1	B	60	ASN
1	B	130	GLU
1	B	135	LEU
1	B	139	VAL
1	B	142	ASN
1	B	143	SER

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Mol	Chain	Res	Type
1	B	151	LEU
1	B	157	ARG
1	C	5	GLU
1	C	8	VAL
1	C	16	LEU
1	C	24	VAL
1	C	25	LEU
1	C	35	GLU
1	C	40	LEU
1	C	44	GLN
1	C	49	VAL
1	C	58	GLU
1	C	59	ILE
1	C	60	ASN
1	C	63	THR
1	C	73	ARG
1	C	85	ILE
1	C	113	THR
1	C	114	ARG
1	C	118	ASP
1	C	122	VAL
1	C	133	GLU
1	C	135	LEU
1	C	139	VAL
1	C	143	SER
1	C	144	VAL
1	C	151	LEU
1	D	8	VAL
1	D	24	VAL
1	D	38	GLU
1	D	39	ILE
1	D	44	GLN
1	D	49	VAL
1	D	51	LEU
1	D	59	ILE
1	D	63	THR
1	D	81	GLU
1	D	82	THR
1	D	84	PRO
1	D	86	VAL
1	D	114	ARG
1	D	119	ASP

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Mol	Chain	Res	Type
1	D	122	VAL
1	D	134	MET
1	D	135	LEU
1	D	136	GLU
1	D	144	VAL
1	D	151	LEU
1	D	156	ARG
1	E	1	MET
1	E	2	ARG
1	E	24	VAL
1	E	25	LEU
1	E	35	GLU
1	E	38	GLU
1	E	40	LEU
1	E	42	GLU
1	E	44	GLN
1	E	49	VAL
1	E	55	VAL
1	E	58	GLU
1	E	65	ARG
1	E	85	ILE
1	E	101	GLU
1	E	121	LEU
1	E	122	VAL
1	E	125	LYS
1	E	126	LEU
1	E	135	LEU
1	E	139	VAL
1	E	144	VAL
1	E	151	LEU
1	E	158	VAL
1	E	159	ARG
1	F	1	MET
1	F	2	ARG
1	F	24	VAL
1	F	25	LEU
1	F	30	ILE
1	F	40	LEU
1	F	44	GLN
1	F	49	VAL
1	F	50	ARG
1	F	51	LEU

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Mol	Chain	Res	Type
1	F	59	ILE
1	F	65	ARG
1	F	75	ILE
1	F	85	ILE
1	F	126	LEU
1	F	129	ASP
1	F	133	GLU
1	F	135	LEU
1	F	139	VAL
1	F	144	VAL
1	F	151	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	0	HIS
1	A	44	GLN
1	B	6	HIS
1	B	142	ASN
1	C	44	GLN
1	D	6	HIS
1	E	6	HIS
1	E	52	HIS
1	E	60	ASN
1	F	6	HIS
1	F	44	GLN
1	F	52	HIS
1	F	60	ASN
1	F	142	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 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	CXS	A	171	-	14,14,14	1.68	1 (7%)	18,18,18	2.59	6 (33%)
3	CXS	C	170	-	14,14,14	1.91	1 (7%)	18,18,18	1.98	5 (27%)
3	CXS	D	170	-	14,14,14	1.69	1 (7%)	18,18,18	2.06	5 (27%)
3	CXS	B	171	-	14,14,14	1.51	2 (14%)	18,18,18	2.18	7 (38%)
3	CXS	D	171	-	14,14,14	1.59	2 (14%)	18,18,18	2.69	5 (27%)
3	CXS	F	171	-	14,14,14	1.79	1 (7%)	18,18,18	2.80	5 (27%)
3	CXS	E	171	-	14,14,14	1.59	2 (14%)	18,18,18	2.35	6 (33%)

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	CXS	A	171	-	-	4/8/16/16	0/1/1/1
3	CXS	C	170	-	-	3/8/16/16	0/1/1/1
3	CXS	D	170	-	-	4/8/16/16	0/1/1/1
3	CXS	B	171	-	-	4/8/16/16	0/1/1/1
3	CXS	D	171	-	-	5/8/16/16	0/1/1/1
3	CXS	F	171	-	-	6/8/16/16	0/1/1/1
3	CXS	E	171	-	-	4/8/16/16	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	170	CXS	C1-S	-6.34	1.68	1.77
3	F	171	CXS	C1-S	-5.98	1.69	1.77
3	A	171	CXS	C1-S	-5.57	1.69	1.77
3	D	170	CXS	C1-S	-5.19	1.70	1.77
3	E	171	CXS	C1-S	-4.16	1.71	1.77
3	D	171	CXS	C1-S	-4.02	1.72	1.77
3	B	171	CXS	C1-S	-4.02	1.72	1.77
3	E	171	CXS	O2-S	3.04	1.53	1.45
3	D	171	CXS	C5-C4	2.57	1.57	1.52
3	B	171	CXS	O1-S	2.51	1.52	1.45

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	171	CXS	O2-S-C1	8.35	119.35	106.73
3	A	171	CXS	O1-S-C1	7.31	117.77	106.73
3	F	171	CXS	C2-C1-S	6.27	122.86	113.25
3	D	171	CXS	O1-S-C1	5.99	115.78	106.73
3	D	171	CXS	C3-N-C4	5.72	125.19	114.18
3	E	171	CXS	O1-S-C1	5.60	115.19	106.73
3	D	170	CXS	C3-N-C4	5.58	124.93	114.18
3	C	170	CXS	C3-N-C4	5.47	124.72	114.18
3	E	171	CXS	C3-N-C4	5.09	123.98	114.18
3	B	171	CXS	O1-S-C1	4.78	113.96	106.73
3	A	171	CXS	C9-C4-C5	-4.41	103.19	110.80
3	D	171	CXS	O3-S-O1	-4.38	100.44	111.40
3	D	171	CXS	C8-C9-C4	4.33	118.86	111.09
3	A	171	CXS	C3-N-C4	4.02	121.92	114.18
3	E	171	CXS	C9-C4-C5	-3.78	104.28	110.80
3	B	171	CXS	O3-S-C1	3.70	113.25	106.00
3	D	171	CXS	C7-C8-C9	3.56	118.73	111.42
3	D	170	CXS	C8-C9-C4	3.43	117.24	111.09
3	B	171	CXS	O2-S-C1	-3.42	101.56	106.73
3	B	171	CXS	O2-S-O1	-3.38	102.84	113.82
3	F	171	CXS	O2-S-O1	-3.27	103.18	113.82
3	A	171	CXS	C8-C7-C6	-2.88	102.64	111.19
3	B	171	CXS	C3-N-C4	2.78	119.53	114.18
3	F	171	CXS	C8-C9-C4	2.74	116.01	111.09
3	C	170	CXS	O3-S-C1	2.70	111.29	106.00
3	E	171	CXS	C8-C9-C4	2.64	115.83	111.09
3	A	171	CXS	O3-S-O1	-2.61	104.86	111.40
3	E	171	CXS	O3-S-O1	-2.59	104.93	111.40
3	C	170	CXS	O1-S-C1	2.38	110.33	106.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	170	CXS	C2-C1-S	2.38	116.89	113.25
3	E	171	CXS	C7-C8-C9	2.35	116.24	111.42
3	D	170	CXS	O1-S-C1	2.21	110.07	106.73
3	C	170	CXS	O2-S-C1	2.19	110.04	106.73
3	C	170	CXS	O3-S-O2	-2.17	105.96	111.40
3	F	171	CXS	C7-C8-C9	-2.13	107.05	111.42
3	B	171	CXS	C7-C6-C5	2.12	115.78	111.42
3	D	170	CXS	O2-S-O1	-2.10	106.99	113.82
3	B	171	CXS	C9-C4-C5	-2.09	107.18	110.80
3	A	171	CXS	C7-C6-C5	2.09	115.72	111.42

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	171	CXS	S-C1-C2-C3
3	A	171	CXS	C5-C4-N-C3
3	B	171	CXS	C2-C1-S-O2
3	B	171	CXS	C2-C1-S-O3
3	D	171	CXS	C1-C2-C3-N
3	D	171	CXS	C9-C4-N-C3
3	F	171	CXS	C1-C2-C3-N
3	F	171	CXS	C9-C4-N-C3
3	D	170	CXS	C2-C1-S-O3
3	F	171	CXS	C2-C1-S-O3
3	E	171	CXS	S-C1-C2-C3
3	C	170	CXS	C1-C2-C3-N
3	E	171	CXS	C2-C3-N-C4
3	B	171	CXS	C2-C1-S-O1
3	D	170	CXS	C2-C1-S-O1
3	D	170	CXS	C2-C1-S-O2
3	F	171	CXS	C2-C1-S-O1
3	F	171	CXS	C2-C1-S-O2
3	A	171	CXS	C2-C3-N-C4
3	C	170	CXS	C2-C3-N-C4
3	D	170	CXS	C2-C3-N-C4
3	F	171	CXS	C5-C4-N-C3
3	C	170	CXS	S-C1-C2-C3
3	D	171	CXS	C2-C1-S-O2
3	E	171	CXS	C2-C1-S-O2
3	B	171	CXS	C2-C3-N-C4
3	E	171	CXS	C1-C2-C3-N

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Mol	Chain	Res	Type	Atoms
3	A	171	CXS	C9-C4-N-C3
3	D	171	CXS	C5-C4-N-C3
3	D	171	CXS	C2-C1-S-O3

There are no ring outliers.

6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	171	CXS	1	0
3	C	170	CXS	2	0
3	D	170	CXS	2	0
3	B	171	CXS	1	0
3	F	171	CXS	2	0
3	E	171	CXS	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	168/172 (97%)	-0.51	2 (1%) 76 73	12, 24, 46, 69	0
1	B	165/172 (95%)	-0.20	6 (3%) 46 42	13, 28, 63, 88	0
1	C	164/172 (95%)	0.35	11 (6%) 24 20	22, 48, 100, 131	0
1	D	164/172 (95%)	0.60	28 (17%) 4 3	19, 54, 123, 156	0
1	E	165/172 (95%)	-0.11	11 (6%) 24 20	17, 34, 82, 118	0
1	F	161/172 (93%)	-0.04	8 (4%) 34 30	19, 41, 93, 128	0
All	All	987/1032 (95%)	0.01	66 (6%) 24 20	12, 36, 97, 156	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	163	GLY	5.4
1	B	161	ALA	5.0
1	B	164	VAL	4.6
1	F	120	GLY	4.5
1	D	59	ILE	4.1
1	D	55	VAL	4.1
1	D	57	GLY	4.0
1	D	62	TRP	3.9
1	C	56	ALA	3.9
1	D	39	ILE	3.9
1	B	56	ALA	3.6
1	D	115	ALA	3.6
1	D	56	ALA	3.5
1	A	167	VAL	3.4
1	C	163	GLY	3.4
1	C	59	ILE	3.3
1	D	0	HIS	3.3
1	C	55	VAL	3.3
1	D	162	GLY	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	126	LEU	3.2
1	D	51	LEU	3.2
1	F	117	GLY	3.2
1	D	117	GLY	3.1
1	C	60	ASN	3.1
1	D	123	ALA	3.1
1	D	85	ILE	3.0
1	C	119	ASP	3.0
1	D	83	ALA	2.9
1	F	160	ALA	2.9
1	E	162	GLY	2.9
1	E	120	GLY	2.8
1	D	60	ASN	2.7
1	E	117	GLY	2.7
1	C	61	THR	2.7
1	F	56	ALA	2.7
1	F	59	ILE	2.6
1	E	124	GLY	2.6
1	B	0	HIS	2.6
1	D	121	LEU	2.6
1	E	123	ALA	2.6
1	D	61	THR	2.6
1	C	83	ALA	2.5
1	C	124	GLY	2.5
1	E	-2	GLY	2.5
1	A	57	GLY	2.4
1	D	53	VAL	2.4
1	C	62	TRP	2.4
1	E	121	LEU	2.4
1	B	59	ILE	2.4
1	C	84	PRO	2.3
1	D	84	PRO	2.3
1	E	116	ALA	2.3
1	F	0	HIS	2.3
1	E	118	ASP	2.2
1	D	134	MET	2.2
1	D	127	THR	2.2
1	B	54	ASP	2.2
1	D	125	LYS	2.2
1	D	54	ASP	2.1
1	D	128	PRO	2.1
1	F	119	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	63	THR	2.1
1	F	127	THR	2.0
1	D	116	ALA	2.0
1	E	56	ALA	2.0
1	E	115	ALA	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	CXS	B	171	14/14	0.71	0.23	57,60,80,82	0
3	CXS	E	171	14/14	0.80	0.18	39,45,51,57	0
3	CXS	D	170	14/14	0.81	0.19	76,81,85,85	0
3	CXS	D	171	14/14	0.90	0.13	38,41,49,49	0
3	CXS	C	170	14/14	0.91	0.12	55,57,61,62	0
3	CXS	F	171	14/14	0.92	0.11	44,47,51,52	0
2	K	F	170	1/1	0.96	0.04	53,53,53,53	0
2	K	B	170	1/1	0.97	0.08	56,56,56,56	0
3	CXS	A	171	14/14	0.97	0.09	27,36,41,41	0
2	K	A	170	1/1	0.98	0.04	37,37,37,37	0
2	K	E	170	1/1	0.98	0.07	36,36,36,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.