



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 07:38 pm BST

PDB ID : 3NEY
Title : Crystal structure of the kinase domain of MPP1/p55
Authors : Shen, Y.; Tong, Y.; Zhong, N.; Guan, X.; Tempel, W.; MacKenzie, F.; Arrow-smith, C.H.; Edwards, A.M.; Bountra, C.; Weigelt, J.; Bochkarev, A.; Park, H.; Structural Genomics Consortium (SGC)
Deposited on : 2010-06-09
Resolution : 2.26 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

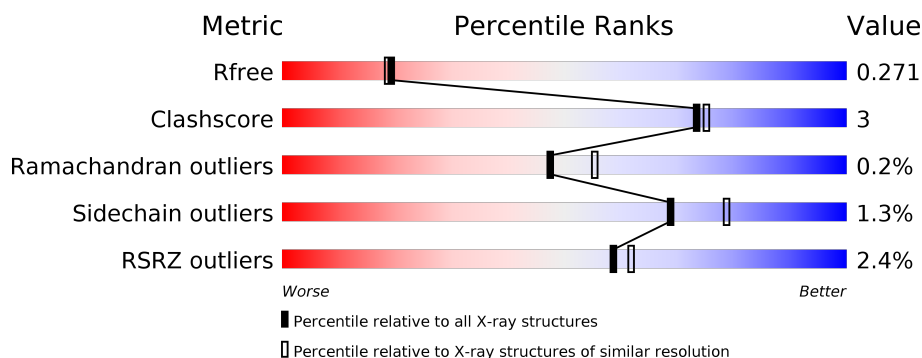
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.26 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}	130704	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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 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	197	<div> <div>2%</div> <div> <div></div> <div>85%</div> <div>6%</div> <div>9%</div> </div> </div>
1	B	197	<div> <div>4%</div> <div> <div></div> <div>85%</div> <div>6%</div> <div>9%</div> </div> </div>
1	C	197	<div> <div>2%</div> <div> <div></div> <div>84%</div> <div>7%</div> <div>9%</div> </div> </div>
1	D	197	<div> <div>2%</div> <div> <div></div> <div>82%</div> <div>10%</div> <div>9%</div> </div> </div>
1	E	197	<div> <div>3%</div> <div> <div></div> <div>86%</div> <div>6%</div> <div>9%</div> </div> </div>
1	F	197	<div> <div>%</div> <div> <div></div> <div>80%</div> <div>8%</div> <div>11%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	UNX	A	12	-	-	-	X
3	UNX	A	13	-	-	-	X
3	UNX	A	15	-	-	-	X
3	UNX	A	21	-	-	-	X
3	UNX	A	25	-	-	-	X
3	UNX	A	7	-	-	-	X
3	UNX	A	8	-	-	-	X
3	UNX	B	10	-	-	-	X
3	UNX	B	11	-	-	-	X
3	UNX	B	22	-	-	-	X
3	UNX	B	26	-	-	-	X
3	UNX	B	4	-	-	-	X
3	UNX	B	9	-	-	-	X
3	UNX	C	14	-	-	-	X
3	UNX	C	23	-	-	-	X
3	UNX	C	24	-	-	-	X
3	UNX	C	30	-	-	-	X
3	UNX	C	32	-	-	-	X
3	UNX	D	27	-	-	-	X
3	UNX	D	33	-	-	-	X
3	UNX	D	34	-	-	-	X
3	UNX	D	35	-	-	-	X
3	UNX	D	36	-	-	-	X
3	UNX	E	20	-	-	-	X
3	UNX	E	28	-	-	-	X
3	UNX	E	31	-	-	-	X
3	UNX	E	39	-	-	-	X
3	UNX	F	29	-	-	-	X
3	UNX	F	37	-	-	-	X
3	UNX	F	38	-	-	-	X
3	UNX	F	6	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8713 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 55 kDa erythrocyte membrane protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	180	Total	C	N	O	S	0	1	0
			1403	892	241	267	3			
1	B	180	Total	C	N	O	S	0	2	0
			1404	895	237	269	3			
1	C	180	Total	C	N	O	S	0	3	0
			1414	900	242	269	3			
1	D	180	Total	C	N	O	S	0	2	0
			1420	902	240	275	3			
1	F	175	Total	C	N	O	S	0	2	0
			1367	872	234	258	3			
1	E	180	Total	C	N	O	S	0	2	0
			1413	898	240	272	3			

There are 108 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	264	MET	-	EXPRESSION TAG	UNP Q00013
A	265	HIS	-	EXPRESSION TAG	UNP Q00013
A	266	HIS	-	EXPRESSION TAG	UNP Q00013
A	267	HIS	-	EXPRESSION TAG	UNP Q00013
A	268	HIS	-	EXPRESSION TAG	UNP Q00013
A	269	HIS	-	EXPRESSION TAG	UNP Q00013
A	270	HIS	-	EXPRESSION TAG	UNP Q00013
A	271	SER	-	EXPRESSION TAG	UNP Q00013
A	272	SER	-	EXPRESSION TAG	UNP Q00013
A	273	GLY	-	EXPRESSION TAG	UNP Q00013
A	274	ARG	-	EXPRESSION TAG	UNP Q00013
A	275	GLU	-	EXPRESSION TAG	UNP Q00013
A	276	ASN	-	EXPRESSION TAG	UNP Q00013
A	277	LEU	-	EXPRESSION TAG	UNP Q00013
A	278	TYR	-	EXPRESSION TAG	UNP Q00013
A	279	PHE	-	EXPRESSION TAG	UNP Q00013
A	280	GLN	-	EXPRESSION TAG	UNP Q00013

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Chain	Residue	Modelled	Actual	Comment	Reference
A	281	GLY	-	EXPRESSION TAG	UNP Q00013
B	264	MET	-	EXPRESSION TAG	UNP Q00013
B	265	HIS	-	EXPRESSION TAG	UNP Q00013
B	266	HIS	-	EXPRESSION TAG	UNP Q00013
B	267	HIS	-	EXPRESSION TAG	UNP Q00013
B	268	HIS	-	EXPRESSION TAG	UNP Q00013
B	269	HIS	-	EXPRESSION TAG	UNP Q00013
B	270	HIS	-	EXPRESSION TAG	UNP Q00013
B	271	SER	-	EXPRESSION TAG	UNP Q00013
B	272	SER	-	EXPRESSION TAG	UNP Q00013
B	273	GLY	-	EXPRESSION TAG	UNP Q00013
B	274	ARG	-	EXPRESSION TAG	UNP Q00013
B	275	GLU	-	EXPRESSION TAG	UNP Q00013
B	276	ASN	-	EXPRESSION TAG	UNP Q00013
B	277	LEU	-	EXPRESSION TAG	UNP Q00013
B	278	TYR	-	EXPRESSION TAG	UNP Q00013
B	279	PHE	-	EXPRESSION TAG	UNP Q00013
B	280	GLN	-	EXPRESSION TAG	UNP Q00013
B	281	GLY	-	EXPRESSION TAG	UNP Q00013
C	264	MET	-	EXPRESSION TAG	UNP Q00013
C	265	HIS	-	EXPRESSION TAG	UNP Q00013
C	266	HIS	-	EXPRESSION TAG	UNP Q00013
C	267	HIS	-	EXPRESSION TAG	UNP Q00013
C	268	HIS	-	EXPRESSION TAG	UNP Q00013
C	269	HIS	-	EXPRESSION TAG	UNP Q00013
C	270	HIS	-	EXPRESSION TAG	UNP Q00013
C	271	SER	-	EXPRESSION TAG	UNP Q00013
C	272	SER	-	EXPRESSION TAG	UNP Q00013
C	273	GLY	-	EXPRESSION TAG	UNP Q00013
C	274	ARG	-	EXPRESSION TAG	UNP Q00013
C	275	GLU	-	EXPRESSION TAG	UNP Q00013
C	276	ASN	-	EXPRESSION TAG	UNP Q00013
C	277	LEU	-	EXPRESSION TAG	UNP Q00013
C	278	TYR	-	EXPRESSION TAG	UNP Q00013
C	279	PHE	-	EXPRESSION TAG	UNP Q00013
C	280	GLN	-	EXPRESSION TAG	UNP Q00013
C	281	GLY	-	EXPRESSION TAG	UNP Q00013
D	264	MET	-	EXPRESSION TAG	UNP Q00013
D	265	HIS	-	EXPRESSION TAG	UNP Q00013
D	266	HIS	-	EXPRESSION TAG	UNP Q00013
D	267	HIS	-	EXPRESSION TAG	UNP Q00013
D	268	HIS	-	EXPRESSION TAG	UNP Q00013

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Chain	Residue	Modelled	Actual	Comment	Reference
D	269	HIS	-	EXPRESSION TAG	UNP Q00013
D	270	HIS	-	EXPRESSION TAG	UNP Q00013
D	271	SER	-	EXPRESSION TAG	UNP Q00013
D	272	SER	-	EXPRESSION TAG	UNP Q00013
D	273	GLY	-	EXPRESSION TAG	UNP Q00013
D	274	ARG	-	EXPRESSION TAG	UNP Q00013
D	275	GLU	-	EXPRESSION TAG	UNP Q00013
D	276	ASN	-	EXPRESSION TAG	UNP Q00013
D	277	LEU	-	EXPRESSION TAG	UNP Q00013
D	278	TYR	-	EXPRESSION TAG	UNP Q00013
D	279	PHE	-	EXPRESSION TAG	UNP Q00013
D	280	GLN	-	EXPRESSION TAG	UNP Q00013
D	281	GLY	-	EXPRESSION TAG	UNP Q00013
F	264	MET	-	EXPRESSION TAG	UNP Q00013
F	265	HIS	-	EXPRESSION TAG	UNP Q00013
F	266	HIS	-	EXPRESSION TAG	UNP Q00013
F	267	HIS	-	EXPRESSION TAG	UNP Q00013
F	268	HIS	-	EXPRESSION TAG	UNP Q00013
F	269	HIS	-	EXPRESSION TAG	UNP Q00013
F	270	HIS	-	EXPRESSION TAG	UNP Q00013
F	271	SER	-	EXPRESSION TAG	UNP Q00013
F	272	SER	-	EXPRESSION TAG	UNP Q00013
F	273	GLY	-	EXPRESSION TAG	UNP Q00013
F	274	ARG	-	EXPRESSION TAG	UNP Q00013
F	275	GLU	-	EXPRESSION TAG	UNP Q00013
F	276	ASN	-	EXPRESSION TAG	UNP Q00013
F	277	LEU	-	EXPRESSION TAG	UNP Q00013
F	278	TYR	-	EXPRESSION TAG	UNP Q00013
F	279	PHE	-	EXPRESSION TAG	UNP Q00013
F	280	GLN	-	EXPRESSION TAG	UNP Q00013
F	281	GLY	-	EXPRESSION TAG	UNP Q00013
E	264	MET	-	EXPRESSION TAG	UNP Q00013
E	265	HIS	-	EXPRESSION TAG	UNP Q00013
E	266	HIS	-	EXPRESSION TAG	UNP Q00013
E	267	HIS	-	EXPRESSION TAG	UNP Q00013
E	268	HIS	-	EXPRESSION TAG	UNP Q00013
E	269	HIS	-	EXPRESSION TAG	UNP Q00013
E	270	HIS	-	EXPRESSION TAG	UNP Q00013
E	271	SER	-	EXPRESSION TAG	UNP Q00013
E	272	SER	-	EXPRESSION TAG	UNP Q00013
E	273	GLY	-	EXPRESSION TAG	UNP Q00013
E	274	ARG	-	EXPRESSION TAG	UNP Q00013

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Chain	Residue	Modelled	Actual	Comment	Reference
E	275	GLU	-	EXPRESSION TAG	UNP Q00013
E	276	ASN	-	EXPRESSION TAG	UNP Q00013
E	277	LEU	-	EXPRESSION TAG	UNP Q00013
E	278	TYR	-	EXPRESSION TAG	UNP Q00013
E	279	PHE	-	EXPRESSION TAG	UNP Q00013
E	280	GLN	-	EXPRESSION TAG	UNP Q00013
E	281	GLY	-	EXPRESSION TAG	UNP Q00013

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is UNKNOWN ATOM OR ION (three-letter code: UNX) (formula: X).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	5	Total	X	0	0
			5	5		
3	E	4	Total	X	0	0
			4	4		
3	B	6	Total	X	0	0
			6	6		
3	C	5	Total	X	0	0
			5	5		
3	A	7	Total	X	0	0
			7	7		
3	F	4	Total	X	0	0
			4	4		

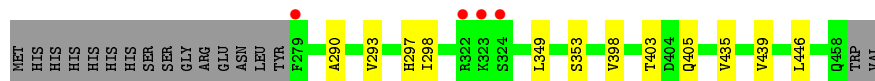
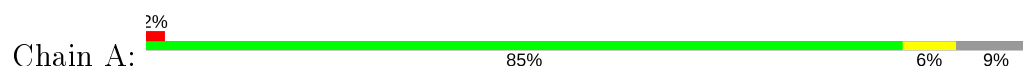
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	29	Total	O	0	0
			29	29		
4	B	45	Total	O	0	0
			45	45		
4	C	36	Total	O	0	0
			36	36		
4	D	32	Total	O	0	0
			32	32		
4	F	37	Total	O	0	0
			37	37		
4	E	22	Total	O	0	0
			22	22		

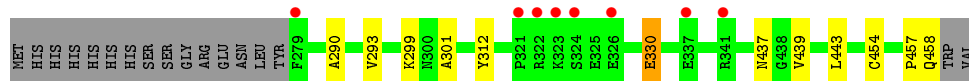
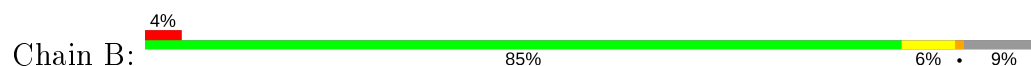
3 Residue-property plots [i](#)

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

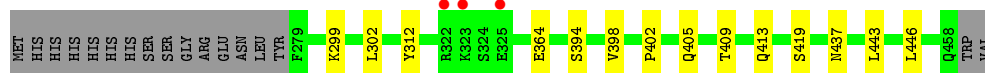
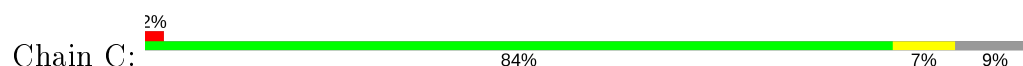
- Molecule 1: 55 kDa erythrocyte membrane protein



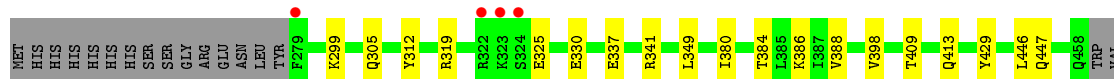
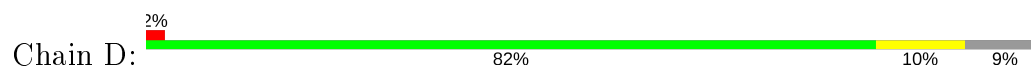
- Molecule 1: 55 kDa erythrocyte membrane protein



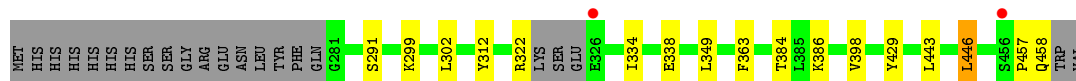
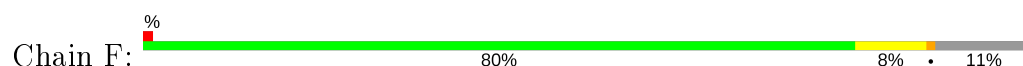
- Molecule 1: 55 kDa erythrocyte membrane protein



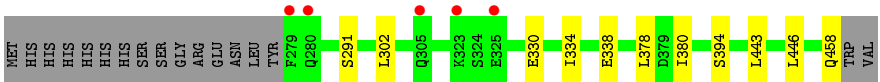
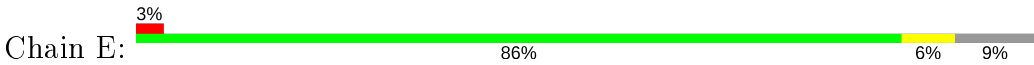
- Molecule 1: 55 kDa erythrocyte membrane protein



- Molecule 1: 55 kDa erythrocyte membrane protein



- Molecule 1: 55 kDa erythrocyte membrane protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.09Å 131.56Å 237.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.26 29.99 – 2.26	Depositor EDS
% Data completeness (in resolution range)	99.3 (30.00-2.26) 99.3 (29.99-2.26)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.26Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.235 , 0.271 0.234 , 0.271	Depositor DCC
R_{free} test set	2121 reflections (2.01%)	wwPDB-VP
Wilson B-factor (Å ²)	33.2	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 45.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8713	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 21.81 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.4930e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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

5.1 Standard geometry

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.78	0/1437	0.74	1/1951 (0.1%)
1	B	0.79	1/1441 (0.1%)	0.72	0/1957
1	C	0.79	1/1454 (0.1%)	0.73	0/1972
1	D	0.80	0/1456	0.75	0/1974
1	E	0.77	0/1449	0.74	0/1966
1	F	0.77	0/1401	0.73	0/1899
All	All	0.78	2/8638 (0.0%)	0.74	1/11719 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	454	CYS	CB-SG	5.91	1.92	1.82
1	C	364	GLU	CG-CD	5.40	1.60	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	349	LEU	CA-CB-CG	5.16	127.18	115.30

There are no chirality outliers.

There are no planarity outliers.

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	1403	0	1340	9	0
1	B	1404	0	1334	6	0
1	C	1414	0	1363	8	0
1	D	1420	0	1364	13	0
1	E	1413	0	1353	5	0
1	F	1367	0	1323	14	0
2	A	10	0	0	0	0
2	B	20	0	0	0	0
2	C	10	0	0	0	0
2	D	10	0	0	0	0
2	E	5	0	0	0	0
2	F	5	0	0	0	0
3	A	7	0	0	0	0
3	B	6	0	0	0	0
3	C	5	0	0	0	0
3	D	5	0	0	0	0
3	E	4	0	0	0	0
3	F	4	0	0	0	0
4	A	29	0	0	0	0
4	B	45	0	0	0	0
4	C	36	0	0	0	0
4	D	32	0	0	0	0
4	E	22	0	0	0	0
4	F	37	0	0	0	0
All	All	8713	0	8077	54	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 3.

All (54) 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:A:297[B]:HIS:ND1	1:A:439:VAL:HG21	1.47	1.27
1:F:302:LEU:CD2	1:F:443:LEU:HD11	1.99	0.92
1:F:302:LEU:HD21	1:F:443:LEU:HD11	1.66	0.77
1:F:302:LEU:HD23	1:F:443:LEU:HD11	1.76	0.66
1:C:398:VAL:HG21	1:C:446:LEU:HD11	1.79	0.64
1:D:319:ARG:NH2	1:D:325:GLU:OE2	2.34	0.61
1:B:330:GLU:H	1:B:330:GLU:CD	2.05	0.60
1:A:398:VAL:HG21	1:A:446:LEU:HD11	1.83	0.60
1:F:446:LEU:O	1:F:446:LEU:HD12	2.02	0.60
1:D:349:LEU:HD11	1:D:388:VAL:HG12	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:299:LYS:HD2	1:D:312:TYR:CD1	2.38	0.58
1:F:386:LYS:HD3	1:F:429:TYR:OH	2.06	0.56
1:F:302:LEU:HD21	1:F:443:LEU:CD1	2.35	0.55
1:D:386:LYS:HD3	1:D:429:TYR:OH	2.08	0.53
1:B:457:PRO:O	1:B:458:GLN:HB2	2.09	0.53
1:D:386:LYS:CD	1:D:429:TYR:OH	2.56	0.53
1:E:302:LEU:CD2	1:E:443:LEU:HD11	2.39	0.53
1:C:299:LYS:HD2	1:C:312:TYR:CG	2.45	0.52
1:E:446:LEU:HD12	1:E:446:LEU:O	2.09	0.52
1:A:405:GLN:NE2	1:C:405:GLN:OE1	2.43	0.51
1:F:299:LYS:HD2	1:F:312:TYR:CD1	2.46	0.51
1:B:301:ALA:HB1	1:B:443:LEU:HD11	1.93	0.51
1:E:334:ILE:HD12	1:E:338:GLU:HG2	1.93	0.50
1:C:402:PRO:HD3	1:C:419[B]:SER:OG	2.12	0.50
1:D:409:THR:O	1:D:413:GLN:HG3	2.13	0.49
1:B:299:LYS:HD2	1:B:312:TYR:CG	2.48	0.49
1:F:398:VAL:HG21	1:F:446:LEU:HD13	1.96	0.47
1:F:457:PRO:O	1:F:458:GLN:HB2	2.13	0.47
1:F:302:LEU:CD2	1:F:443:LEU:CD1	2.83	0.47
1:B:439:VAL:O	1:B:443:LEU:HG	2.14	0.47
1:D:349:LEU:HD11	1:D:388:VAL:CG1	2.45	0.47
1:A:297[B]:HIS:ND1	1:A:439:VAL:CG2	2.43	0.46
1:A:403:THR:HA	1:A:435:VAL:CG1	2.45	0.46
1:A:297[B]:HIS:CE1	1:A:439:VAL:HG21	2.38	0.45
1:C:409:THR:O	1:C:413:GLN:HG2	2.16	0.45
1:A:290:ALA:O	1:A:293:VAL:HG22	2.16	0.45
1:A:298:ILE:HD13	1:A:439:VAL:HG22	1.98	0.44
1:C:299:LYS:HD2	1:C:312:TYR:CD1	2.52	0.44
1:D:305:GLN:NE2	1:D:447:GLN:OE1	2.47	0.43
1:B:290:ALA:O	1:B:293:VAL:HG22	2.19	0.43
1:D:386:LYS:HD2	1:D:429:TYR:OH	2.19	0.43
1:F:446:LEU:C	1:F:446:LEU:HD12	2.39	0.43
1:D:380:ILE:HD12	1:D:384:THR:HB	2.00	0.42
1:D:398:VAL:HG21	1:D:446:LEU:CD1	2.50	0.42
1:E:446:LEU:HD12	1:E:446:LEU:C	2.40	0.42
1:F:384:THR:HG22	1:F:384:THR:O	2.20	0.42
1:E:378:LEU:HB3	1:E:380:ILE:HG23	2.02	0.42
1:C:302:LEU:HG	1:C:443:LEU:HD22	2.00	0.42
1:F:349:LEU:HA	1:F:363:PHE:CZ	2.54	0.41
1:D:337:GLU:O	1:D:341:ARG:HG3	2.20	0.41
1:D:398:VAL:HG21	1:D:446:LEU:HD13	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:ILE:CD1	1:A:439:VAL:HG22	2.50	0.41
1:F:334:ILE:HD12	1:F:338:GLU:HG2	2.03	0.41
1:C:398:VAL:HG21	1:C:446:LEU:CD1	2.48	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	179/197 (91%)	175 (98%)	4 (2%)	0	100	100
1	B	180/197 (91%)	173 (96%)	6 (3%)	1 (1%)	25	25
1	C	181/197 (92%)	176 (97%)	4 (2%)	1 (1%)	25	25
1	D	180/197 (91%)	177 (98%)	3 (2%)	0	100	100
1	E	180/197 (91%)	174 (97%)	6 (3%)	0	100	100
1	F	173/197 (88%)	171 (99%)	2 (1%)	0	100	100
All	All	1073/1182 (91%)	1046 (98%)	25 (2%)	2 (0%)	47	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	437	ASN
1	C	437	ASN

5.3.2 Protein sidechains [i](#)

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	147/175 (84%)	146 (99%)	1 (1%)	84	90
1	B	146/175 (83%)	145 (99%)	1 (1%)	84	90
1	C	150/175 (86%)	149 (99%)	1 (1%)	84	90
1	D	151/175 (86%)	149 (99%)	2 (1%)	69	79
1	E	149/175 (85%)	145 (97%)	4 (3%)	44	54
1	F	143/175 (82%)	140 (98%)	3 (2%)	53	62
All	All	886/1050 (84%)	874 (99%)	12 (1%)	69	76

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	353	SER
1	B	330	GLU
1	C	394	SER
1	D	330[A]	GLU
1	D	330[B]	GLU
1	F	291	SER
1	F	322	ARG
1	F	446	LEU
1	E	291	SER
1	E	330	GLU
1	E	394	SER
1	E	458	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	405	GLN
1	C	405	GLN

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

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 43 ligands modelled in this entry, 31 are unknown - leaving 12 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	$\# Z > 2$
2	SO4	B	7	-	4,4,4	0.22	0	6,6,6	1.14	1 (16%)
2	SO4	E	12	-	4,4,4	0.13	0	6,6,6	0.27	0
2	SO4	A	3	-	4,4,4	0.26	0	6,6,6	0.59	0
2	SO4	D	10	-	4,4,4	0.19	0	6,6,6	0.25	0
2	SO4	C	9	-	4,4,4	0.22	0	6,6,6	0.25	0
2	SO4	F	11	-	4,4,4	0.15	0	6,6,6	0.26	0
2	SO4	C	1	-	4,4,4	0.22	0	6,6,6	0.77	0
2	SO4	B	6	-	4,4,4	0.22	0	6,6,6	0.29	0
2	SO4	D	2	-	4,4,4	0.20	0	6,6,6	0.35	0
2	SO4	A	4	-	4,4,4	0.16	0	6,6,6	0.20	0
2	SO4	B	5	-	4,4,4	0.15	0	6,6,6	0.37	0
2	SO4	B	8	-	4,4,4	0.12	0	6,6,6	0.35	0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	7	SO4	O4-S-O3	-2.32	99.17	109.06

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	180/197 (91%)	-0.06	4 (2%) 62 65	17, 28, 53, 61	0
1	B	180/197 (91%)	-0.10	8 (4%) 34 37	15, 27, 52, 64	0
1	C	180/197 (91%)	-0.15	3 (1%) 70 73	17, 26, 51, 64	0
1	D	180/197 (91%)	-0.05	4 (2%) 62 65	16, 29, 50, 62	0
1	E	180/197 (91%)	-0.13	5 (2%) 53 55	18, 27, 55, 65	0
1	F	175/197 (88%)	-0.14	2 (1%) 80 82	17, 28, 47, 69	0
All	All	1075/1182 (90%)	-0.11	26 (2%) 59 62	15, 27, 52, 69	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	325	GLU	3.9
1	D	323	LYS	3.8
1	E	280	GLN	3.6
1	A	324	SER	3.6
1	E	279	PHE	3.4
1	B	324	SER	3.2
1	F	456	SER	3.0
1	B	323	LYS	3.0
1	D	279	PHE	2.8
1	C	323	LYS	2.8
1	C	322	ARG	2.8
1	D	324	SER	2.8
1	E	325	GLU	2.7
1	F	326	GLU	2.6
1	B	326	GLU	2.6
1	D	322	ARG	2.2
1	B	341	ARG	2.2
1	A	323	LYS	2.2
1	B	322	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	279	PHE	2.2
1	E	305	GLN	2.2
1	A	279	PHE	2.1
1	E	323	LYS	2.1
1	B	337	GLU	2.0
1	A	322	ARG	2.0
1	B	321	PRO	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 carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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(\AA^2)	Q<0.9
3	UNX	B	11	1/1	-0.27	3.09	2,2,2,2	1
3	UNX	A	7	1/1	-0.08	2.67	2,2,2,2	1
3	UNX	E	20	1/1	0.01	3.38	2,2,2,2	1
3	UNX	C	30	1/1	0.01	4.22	2,2,2,2	1
3	UNX	A	15	1/1	0.01	4.13	2,2,2,2	1
3	UNX	E	28	1/1	0.04	0.84	2,2,2,2	1
3	UNX	B	4	1/1	0.05	1.58	2,2,2,2	1
3	UNX	E	39	1/1	0.09	1.33	2,2,2,2	1
3	UNX	F	37	1/1	0.12	3.87	2,2,2,2	1
3	UNX	A	8	1/1	0.20	1.43	2,2,2,2	1
3	UNX	A	21	1/1	0.21	2.53	2,2,2,2	1
3	UNX	A	25	1/1	0.26	1.62	2,2,2,2	1
3	UNX	D	34	1/1	0.28	1.67	2,2,2,2	1
3	UNX	C	24	1/1	0.31	1.98	2,2,2,2	1
3	UNX	B	22	1/1	0.33	1.76	2,2,2,2	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	UNX	C	23	1/1	0.34	1.61	2,2,2,2	1
3	UNX	D	27	1/1	0.38	1.57	2,2,2,2	1
3	UNX	B	9	1/1	0.38	2.61	2,2,2,2	1
3	UNX	B	26	1/1	0.45	0.89	2,2,2,2	1
3	UNX	D	33	1/1	0.47	1.59	2,2,2,2	1
3	UNX	D	36	1/1	0.50	2.91	2,2,2,2	1
3	UNX	B	10	1/1	0.50	1.25	2,2,2,2	1
3	UNX	D	35	1/1	0.52	2.75	2,2,2,2	1
3	UNX	C	32	1/1	0.53	1.72	2,2,2,2	1
3	UNX	C	14	1/1	0.54	2.27	2,2,2,2	1
3	UNX	F	29	1/1	0.56	0.96	2,2,2,2	1
3	UNX	F	38	1/1	0.62	2.12	2,2,2,2	1
3	UNX	F	6	1/1	0.64	1.93	2,2,2,2	1
3	UNX	E	31	1/1	0.69	1.69	2,2,2,2	1
3	UNX	A	13	1/1	0.70	3.46	2,2,2,2	1
3	UNX	A	12	1/1	0.75	1.53	2,2,2,2	1
2	SO4	F	11	5/5	0.85	0.36	87,87,87,88	0
2	SO4	E	12	5/5	0.90	0.25	92,93,93,93	0
2	SO4	A	4	5/5	0.93	0.33	82,83,84,84	0
2	SO4	B	5	5/5	0.94	0.35	81,82,82,83	0
2	SO4	B	8	5/5	0.95	0.14	59,60,60,61	0
2	SO4	C	9	5/5	0.95	0.17	66,67,68,68	0
2	SO4	D	10	5/5	0.96	0.35	83,83,84,84	0
2	SO4	B	6	5/5	0.97	0.30	63,63,64,65	0
2	SO4	A	3	5/5	0.98	0.11	25,25,26,26	0
2	SO4	C	1	5/5	0.99	0.13	23,24,27,28	0
2	SO4	D	2	5/5	0.99	0.09	20,20,22,24	0
2	SO4	B	7	5/5	0.99	0.11	27,28,30,30	0

6.5 Other polymers

There are no such residues in this entry.