



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:18 AM GMT

PDB ID : 3OLR
Title : PTPN22 in complex with consensus phospho-tyrosine peptide 1
Authors : Yu, X.; Sun, J.-P.; Zhang, S.; Zhang, Z.-Y.
Deposited on : 2010-08-26
Resolution : 2.50 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

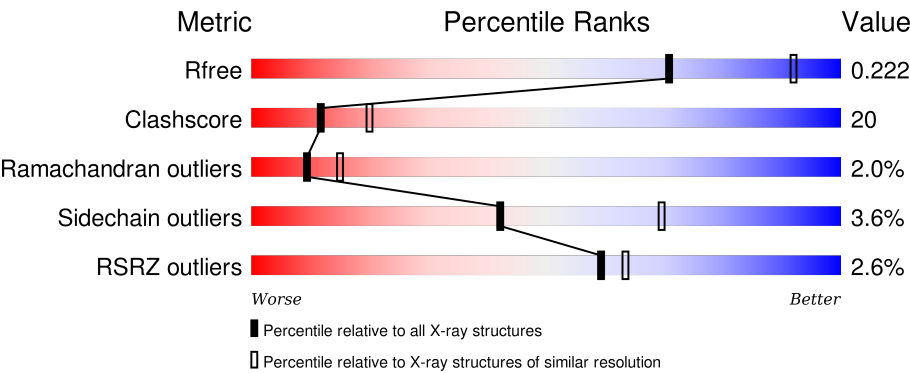
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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. 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 <=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	313	<div><div>2%</div><div>55%38%5%</div></div>
1	B	313	<div><div>2%</div><div>57%35%5%</div></div>
1	C	313	<div><div>2%</div><div>59%33%5%</div></div>
1	D	313	<div><div>2%</div><div>61%32%5%</div></div>
2	E	9	<div><div>11%11%</div><div>33%33%22%</div></div>

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Mol	Chain	Length	Quality of chain
2	F	9	<div><div><div></div><div></div><div></div><div></div></div><div>22%33%11%44%11%</div></div>
2	G	9	<div><div><div></div><div></div><div></div><div></div></div><div>33%11%56%33%</div></div>
2	H	9	<div><div><div></div><div></div><div></div><div></div></div><div>22%22%11%56%11%</div></div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 10534 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 Tyrosine-protein phosphatase non-receptor type 22.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	S	0	0	0
			2451	1573	402	459	17			
1	B	297	Total	C	N	O	S	0	0	0
			2451	1573	402	459	17			
1	C	297	Total	C	N	O	S	0	0	0
			2451	1573	402	459	17			
1	D	297	Total	C	N	O	S	0	0	0
			2451	1573	402	459	17			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
A	-18	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
A	-17	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
A	-16	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
A	-15	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
A	-14	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
A	-13	SER	-	EXPRESSION TAG	UNP Q9Y2R2
A	-12	SER	-	EXPRESSION TAG	UNP Q9Y2R2
A	-11	GLY	-	EXPRESSION TAG	UNP Q9Y2R2
A	-10	LEU	-	EXPRESSION TAG	UNP Q9Y2R2
A	-9	VAL	-	EXPRESSION TAG	UNP Q9Y2R2
A	-8	PRO	-	EXPRESSION TAG	UNP Q9Y2R2
A	-7	ARG	-	EXPRESSION TAG	UNP Q9Y2R2
A	-6	GLY	-	EXPRESSION TAG	UNP Q9Y2R2
A	-5	SER	-	EXPRESSION TAG	UNP Q9Y2R2
A	-4	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
A	-3	MET	-	EXPRESSION TAG	UNP Q9Y2R2
A	-2	ALA	-	EXPRESSION TAG	UNP Q9Y2R2
A	-1	SER	-	EXPRESSION TAG	UNP Q9Y2R2
A	227	SER	CYS	ENGINEERED MUTATION	UNP Q9Y2R2
B	-19	HIS	-	EXPRESSION TAG	UNP Q9Y2R2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
B	-17	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
B	-16	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
B	-15	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
B	-14	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
B	-13	SER	-	EXPRESSION TAG	UNP Q9Y2R2
B	-12	SER	-	EXPRESSION TAG	UNP Q9Y2R2
B	-11	GLY	-	EXPRESSION TAG	UNP Q9Y2R2
B	-10	LEU	-	EXPRESSION TAG	UNP Q9Y2R2
B	-9	VAL	-	EXPRESSION TAG	UNP Q9Y2R2
B	-8	PRO	-	EXPRESSION TAG	UNP Q9Y2R2
B	-7	ARG	-	EXPRESSION TAG	UNP Q9Y2R2
B	-6	GLY	-	EXPRESSION TAG	UNP Q9Y2R2
B	-5	SER	-	EXPRESSION TAG	UNP Q9Y2R2
B	-4	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
B	-3	MET	-	EXPRESSION TAG	UNP Q9Y2R2
B	-2	ALA	-	EXPRESSION TAG	UNP Q9Y2R2
B	-1	SER	-	EXPRESSION TAG	UNP Q9Y2R2
B	227	SER	CYS	ENGINEERED MUTATION	UNP Q9Y2R2
C	-19	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
C	-18	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
C	-17	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
C	-16	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
C	-15	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
C	-14	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
C	-13	SER	-	EXPRESSION TAG	UNP Q9Y2R2
C	-12	SER	-	EXPRESSION TAG	UNP Q9Y2R2
C	-11	GLY	-	EXPRESSION TAG	UNP Q9Y2R2
C	-10	LEU	-	EXPRESSION TAG	UNP Q9Y2R2
C	-9	VAL	-	EXPRESSION TAG	UNP Q9Y2R2
C	-8	PRO	-	EXPRESSION TAG	UNP Q9Y2R2
C	-7	ARG	-	EXPRESSION TAG	UNP Q9Y2R2
C	-6	GLY	-	EXPRESSION TAG	UNP Q9Y2R2
C	-5	SER	-	EXPRESSION TAG	UNP Q9Y2R2
C	-4	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
C	-3	MET	-	EXPRESSION TAG	UNP Q9Y2R2
C	-2	ALA	-	EXPRESSION TAG	UNP Q9Y2R2
C	-1	SER	-	EXPRESSION TAG	UNP Q9Y2R2
C	227	SER	CYS	ENGINEERED MUTATION	UNP Q9Y2R2
D	-19	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
D	-18	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
D	-17	HIS	-	EXPRESSION TAG	UNP Q9Y2R2

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
D	-15	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
D	-14	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
D	-13	SER	-	EXPRESSION TAG	UNP Q9Y2R2
D	-12	SER	-	EXPRESSION TAG	UNP Q9Y2R2
D	-11	GLY	-	EXPRESSION TAG	UNP Q9Y2R2
D	-10	LEU	-	EXPRESSION TAG	UNP Q9Y2R2
D	-9	VAL	-	EXPRESSION TAG	UNP Q9Y2R2
D	-8	PRO	-	EXPRESSION TAG	UNP Q9Y2R2
D	-7	ARG	-	EXPRESSION TAG	UNP Q9Y2R2
D	-6	GLY	-	EXPRESSION TAG	UNP Q9Y2R2
D	-5	SER	-	EXPRESSION TAG	UNP Q9Y2R2
D	-4	HIS	-	EXPRESSION TAG	UNP Q9Y2R2
D	-3	MET	-	EXPRESSION TAG	UNP Q9Y2R2
D	-2	ALA	-	EXPRESSION TAG	UNP Q9Y2R2
D	-1	SER	-	EXPRESSION TAG	UNP Q9Y2R2
D	227	SER	CYS	ENGINEERED MUTATION	UNP Q9Y2R2

- Molecule 2 is a protein called SKAP2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	9	Total	C	N	O	P	0	0	0
			87	53	9	24	1			
2	F	8	Total	C	N	O	P	0	0	0
			74	44	8	21	1			
2	G	9	Total	C	N	O	P	0	0	0
			87	53	9	24	1			
2	H	9	Total	C	N	O	P	0	0	0
			87	53	9	24	1			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	84	Total	O	0	0
			84	84		
3	B	95	Total	O	0	0
			95	95		
3	C	108	Total	O	0	0
			108	108		
3	D	96	Total	O	0	0
			96	96		
3	E	4	Total	O	0	0
			4	4		

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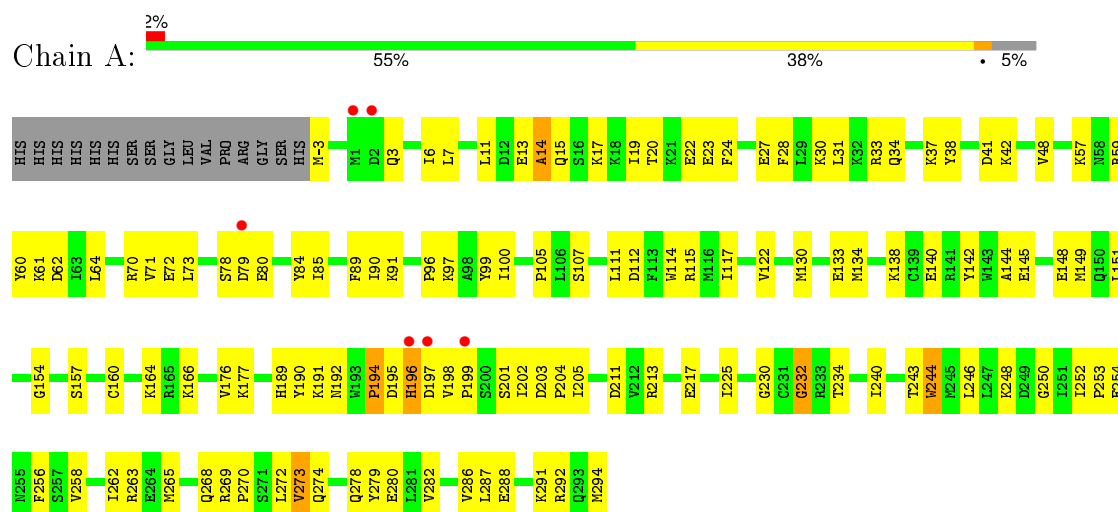
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	F	2	Total	O	0	0
			2	2		
3	G	4	Total	O	0	0
			4	4		
3	H	2	Total	O	0	0
			2	2		

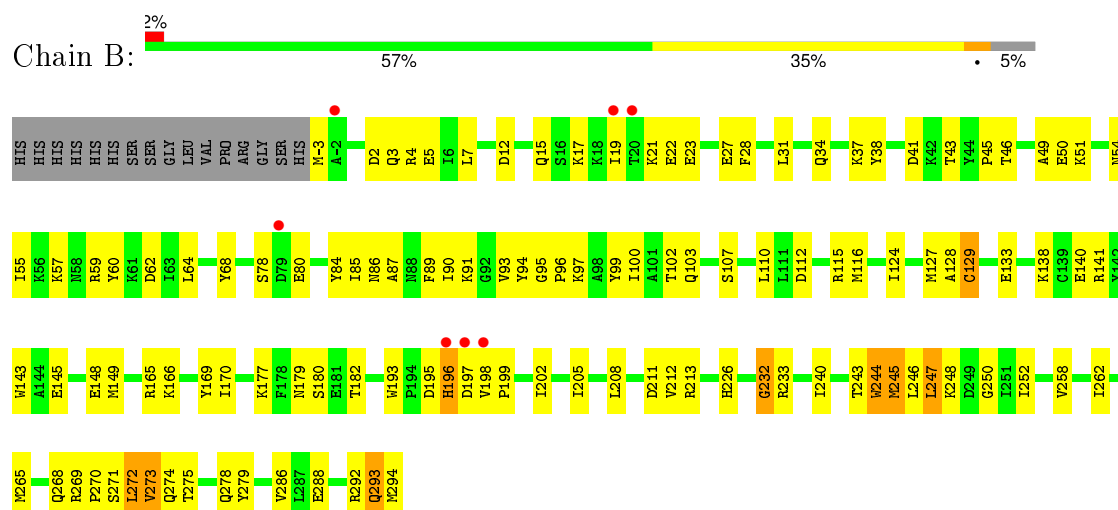
3 Residue-property plots

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: Tyrosine-protein phosphatase non-receptor type 22

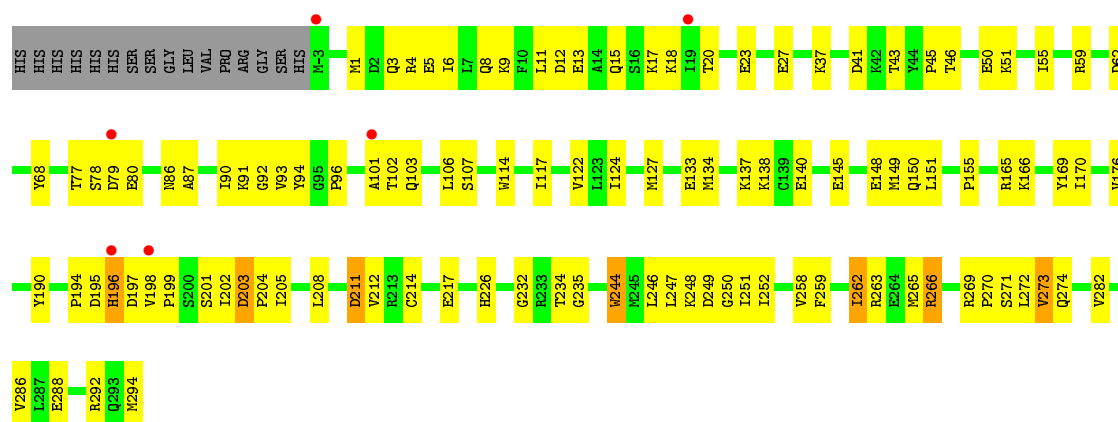


- Molecule 1: Tyrosine-protein phosphatase non-receptor type 22



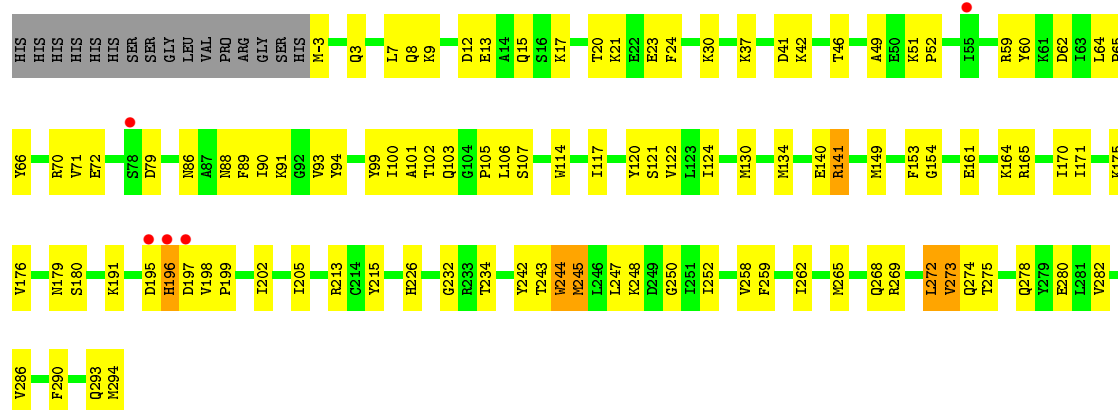
- Molecule 1: Tyrosine-protein phosphatase non-receptor type 22





- Molecule 1: Tyrosine-protein phosphatase non-receptor type 22

Chain D: 2% 61% 32% 5%



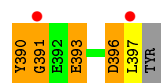
- Molecule 2: SKAP2

Chain E: 11% 33% 33% 22%



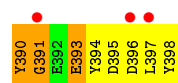
- Molecule 2: SKAP2

Chain F: 22% 33% 11% 44% 11%



- Molecule 2: SKAP2

Chain G: 33% 11% 56% 33%



● Molecule 2: SKAP2

Chain H: 22% 22% 11% 56% 11%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	45.90 Å 62.77 Å 117.46 Å 99.06° 96.53° 105.06°	Depositor
Resolution (Å)	50.00 – 2.50 31.38 – 2.06	Depositor EDS
% Data completeness (in resolution range)	89.1 (50.00-2.50) 89.5 (31.38-2.06)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 2.06 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.162 , 0.210 0.180 , 0.222	Depositor DCC
R_{free} test set	3826 reflections (9.17%)	DCC
Wilson B-factor (Å ²)	29.3	Xtriage
Anisotropy	0.075	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 50.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 71590 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10534	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.64% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: PTR

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	0.33	0/2508	0.57	0/3386
1	B	0.35	0/2508	0.59	0/3386
1	C	0.36	1/2508 (0.0%)	0.60	0/3386
1	D	0.35	0/2508	0.58	0/3386
2	E	0.49	0/71	0.84	1/92 (1.1%)
2	F	0.49	0/57	0.62	0/74
2	G	0.49	0/71	0.75	0/92
2	H	0.54	0/71	1.12	1/92 (1.1%)
All	All	0.35	1/10302 (0.0%)	0.59	2/13894 (0.0%)

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

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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	150	GLN	C-N	-5.22	1.22	1.34

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	397	LEU	N-CA-C	-5.39	96.44	111.00
2	H	397	LEU	N-CA-C	5.30	125.32	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	398	TYR	Sidechain

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	2451	0	2441	107	0
1	B	2451	0	2441	99	0
1	C	2451	0	2441	90	0
1	D	2451	0	2441	82	0
2	E	87	0	58	19	0
2	F	74	0	49	10	0
2	G	87	0	58	13	0
2	H	87	0	58	17	0
3	A	84	0	0	6	0
3	B	95	0	0	2	0
3	C	108	0	0	3	0
3	D	96	0	0	5	0
3	E	4	0	0	0	0
3	F	2	0	0	0	0
3	G	4	0	0	0	0
3	H	2	0	0	0	0
All	All	10534	0	9987	407	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 20.

All (407) 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:130:MET:HG2	1:A:191:LYS:HD2	1.42	0.99
1:A:265:MET:HB3	1:A:272:LEU:HD23	1.48	0.96
1:C:3:GLN:HE22	1:C:252:ILE:H	1.13	0.90
1:B:265:MET:HB3	1:B:272:LEU:HD23	1.54	0.89
2:F:393:GLU:HB3	2:F:396:ASP:HB2	1.55	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:393:GLU:HB3	2:E:396:ASP:HB2	1.57	0.87
1:A:134:MET:SD	2:E:398:TYR:HB2	2.17	0.85
1:A:3:GLN:HE22	1:A:252:ILE:H	1.25	0.84
2:H:395:ASP:O	2:H:397:LEU:N	2.09	0.84
1:C:265:MET:HB3	1:C:272:LEU:HD23	1.60	0.82
1:D:107:SER:N	1:D:140:GLU:HG3	1.94	0.81
1:A:197:ASP:O	1:A:199:PRO:HD3	1.79	0.80
2:F:393:GLU:CB	2:F:396:ASP:HB2	2.10	0.80
1:B:3:GLN:HE22	1:B:252:ILE:H	1.29	0.80
1:A:157:SER:HB2	1:A:177:LYS:HB2	1.65	0.79
1:D:105:PRO:O	1:D:140:GLU:HB2	1.82	0.79
1:A:17:LYS:HG3	1:A:23:GLU:HG2	1.65	0.78
1:B:202:ILE:O	1:B:205:ILE:HG22	1.85	0.76
1:B:107:SER:N	1:B:140:GLU:HG3	2.01	0.76
2:E:393:GLU:CB	2:E:396:ASP:HB2	2.17	0.74
1:B:262:ILE:HD12	1:B:273:VAL:HG21	1.69	0.74
1:B:12:ASP:HA	1:B:15:GLN:HE21	1.53	0.73
1:C:107:SER:N	1:C:140:GLU:HG3	2.03	0.73
2:H:393:GLU:HG2	2:H:396:ASP:HB2	1.71	0.73
1:D:3:GLN:HE21	1:D:250:GLY:HA2	1.53	0.72
1:D:265:MET:HB3	1:D:272:LEU:HD23	1.70	0.71
1:D:282:VAL:O	1:D:286:VAL:HG23	1.91	0.71
1:A:199:PRO:O	1:A:202:ILE:HG13	1.91	0.71
1:A:130:MET:HG2	1:A:191:LYS:CD	2.19	0.70
1:C:37:LYS:HE2	1:C:41:ASP:OD2	1.92	0.70
1:B:62:ASP:OD1	2:F:393:GLU:HA	1.92	0.69
1:D:93:VAL:HG21	1:D:245:MET:HG2	1.74	0.68
1:C:202:ILE:O	1:C:205:ILE:HG22	1.94	0.68
1:C:258:VAL:O	1:C:262:ILE:HG12	1.95	0.67
1:A:91:LYS:HZ2	1:A:91:LYS:HB2	1.59	0.67
1:D:49:ALA:HB1	1:D:64:LEU:HD22	1.75	0.67
1:D:197:ASP:O	1:D:199:PRO:HD3	1.93	0.67
1:D:17:LYS:HG3	1:D:23:GLU:HG2	1.76	0.67
2:H:393:GLU:HB3	2:H:396:ASP:H	1.59	0.66
1:A:27:GLU:HG2	1:A:279:TYR:OH	1.94	0.66
1:A:288:GLU:O	1:A:292:ARG:HG3	1.95	0.66
1:A:202:ILE:O	1:A:205:ILE:HG22	1.95	0.66
1:C:288:GLU:O	1:C:292:ARG:HG3	1.96	0.66
1:A:265:MET:HB3	1:A:272:LEU:CD2	2.24	0.66
1:B:17:LYS:HG3	1:B:23:GLU:HG2	1.78	0.65
1:A:91:LYS:HG2	1:A:96:PRO:HA	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:12:ASP:HA	1:D:15:GLN:HE21	1.61	0.65
1:C:197:ASP:O	1:C:199:PRO:HD3	1.97	0.65
1:D:244:TRP:HE3	1:D:244:TRP:O	1.80	0.65
1:A:112:ASP:HA	1:A:115:ARG:HD3	1.78	0.64
1:C:199:PRO:O	1:C:202:ILE:HG12	1.96	0.64
1:A:64:LEU:O	1:A:270:PRO:HG3	1.98	0.64
1:A:7:LEU:HD23	1:A:294:MET:HE1	1.79	0.64
1:A:195:ASP:OD2	1:A:198:VAL:HG23	1.98	0.64
1:A:189:HIS:CE1	1:A:191:LYS:HE2	2.33	0.64
1:B:12:ASP:HA	1:B:15:GLN:NE2	2.12	0.64
1:C:195:ASP:HB3	1:C:198:VAL:HB	1.80	0.63
1:C:1:MET:HE2	1:C:5:GLU:HG2	1.80	0.63
1:A:287:LEU:HG	1:A:291:LYS:HE3	1.78	0.63
1:C:117:ILE:HG23	1:C:122:VAL:HB	1.80	0.63
1:B:165:ARG:HG2	1:B:170:ILE:HG12	1.79	0.63
1:B:17:LYS:CG	1:B:23:GLU:HG2	2.28	0.62
1:D:247:LEU:HD11	1:D:290:PHE:HD2	1.63	0.62
1:D:202:ILE:O	1:D:205:ILE:HG22	1.99	0.62
2:E:395:ASP:O	2:E:397:LEU:N	2.32	0.62
1:A:7:LEU:O	1:A:11:LEU:HG	2.00	0.62
1:B:51:LYS:HB2	1:B:54:ASN:HD22	1.65	0.62
1:A:164:LYS:HZ1	1:C:8:GLN:HB3	1.64	0.62
2:H:395:ASP:C	2:H:397:LEU:H	2.01	0.61
1:C:217:GLU:HB2	3:C:298:HOH:O	2.00	0.61
1:A:262:ILE:HD12	1:A:273:VAL:HG21	1.82	0.61
1:A:90:ILE:HB	1:A:99:TYR:HB2	1.82	0.61
1:A:59:ARG:NH1	2:E:390:TYR:CE2	2.68	0.61
2:G:394:PTR:O	2:G:398:TYR:HD2	1.84	0.61
1:D:59:ARG:NH1	2:H:390:TYR:CE2	2.69	0.60
1:D:247:LEU:HD11	1:D:290:PHE:CD2	2.37	0.60
1:B:49:ALA:HB1	1:B:64:LEU:HD22	1.82	0.60
1:A:157:SER:O	1:A:176:VAL:HA	2.02	0.60
2:E:396:ASP:O	2:E:397:LEU:C	2.40	0.60
1:D:51:LYS:HB3	1:D:52:PRO:HD2	1.83	0.60
1:C:133:GLU:HB2	1:C:138:LYS:HG3	1.83	0.59
1:D:3:GLN:HE22	1:D:252:ILE:H	1.48	0.59
1:C:272:LEU:O	1:C:273:VAL:HB	2.03	0.58
1:B:127:MET:HG3	1:B:226:HIS:CE1	2.38	0.58
1:A:57:LYS:HE2	3:A:355:HOH:O	2.03	0.58
1:A:17:LYS:CG	1:A:23:GLU:HG2	2.31	0.58
1:D:62:ASP:OD1	2:H:393:GLU:HA	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:ILE:HG22	1:B:100:ILE:O	2.02	0.58
1:B:127:MET:HE2	1:B:129:CYS:O	2.03	0.58
1:A:145:GLU:O	1:A:148:GLU:HB3	2.04	0.57
1:C:90:ILE:HD11	1:C:269:ARG:HB2	1.86	0.57
1:A:269:ARG:HD2	1:A:270:PRO:HD2	1.85	0.57
1:D:272:LEU:O	1:D:273:VAL:HB	2.04	0.57
1:C:202:ILE:HD11	1:C:288:GLU:HG2	1.87	0.57
1:B:165:ARG:NH1	1:B:170:ILE:HD11	2.20	0.57
2:H:395:ASP:C	2:H:397:LEU:N	2.58	0.57
1:B:195:ASP:HB3	1:B:198:VAL:HB	1.87	0.57
1:A:217:GLU:HB2	3:A:421:HOH:O	2.04	0.57
1:A:3:GLN:NE2	1:A:252:ILE:H	2.00	0.57
1:B:258:VAL:O	1:B:262:ILE:HG12	2.05	0.57
1:A:195:ASP:HB3	1:A:198:VAL:HB	1.87	0.57
1:B:84:TYR:CG	1:B:85:ILE:N	2.73	0.56
1:A:258:VAL:O	1:A:262:ILE:HG12	2.05	0.56
1:C:4:ARG:NH1	1:C:294:MET:HE2	2.21	0.56
1:B:27:GLU:HG2	1:B:279:TYR:OH	2.06	0.56
1:B:19:ILE:HG22	1:B:19:ILE:O	2.05	0.56
1:C:195:ASP:OD2	1:C:198:VAL:HG23	2.05	0.56
1:C:1:MET:HG2	1:C:5:GLU:OE2	2.06	0.55
1:B:166:LYS:HD3	1:B:169:TYR:CE1	2.41	0.55
1:D:258:VAL:O	1:D:262:ILE:HG12	2.07	0.55
1:D:165:ARG:HG2	1:D:170:ILE:HG12	1.87	0.55
1:D:46:THR:HG22	1:D:66:TYR:HA	1.88	0.55
1:D:134:MET:SD	2:H:398:TYR:HB2	2.46	0.55
2:G:394:PTR:O	2:G:398:TYR:CD2	2.59	0.55
1:B:196:HIS:CD2	1:B:196:HIS:C	2.81	0.55
1:C:101:ALA:HB1	1:C:234:THR:HG21	1.88	0.55
1:C:208:LEU:O	1:C:212:VAL:HG23	2.06	0.55
2:E:390:TYR:O	2:E:391:GLY:C	2.45	0.55
1:A:263:ARG:NH2	1:B:96:PRO:HD2	2.22	0.55
1:C:15:GLN:O	1:C:18:LYS:HB3	2.07	0.54
1:B:87:ALA:HB2	1:B:102:THR:OG1	2.07	0.54
2:G:393:GLU:HG2	2:G:396:ASP:HB2	1.89	0.54
1:B:89:PHE:CZ	1:B:97:LYS:HE3	2.42	0.54
1:C:78:SER:C	1:C:80:GLU:H	2.10	0.54
1:C:269:ARG:HG2	1:C:272:LEU:HD13	1.88	0.54
1:B:102:THR:O	1:B:226:HIS:HB2	2.06	0.54
1:C:165:ARG:NH1	1:C:170:ILE:HD11	2.23	0.54
1:A:73:LEU:HD11	1:A:85:ILE:HB	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:259:PHE:HE2	1:C:263:ARG:HH21	1.55	0.54
1:B:269:ARG:HG2	1:B:272:LEU:HD13	1.89	0.54
1:C:59:ARG:NH1	2:G:390:TYR:CE2	2.76	0.53
2:F:397:LEU:N	2:F:397:LEU:HD22	2.23	0.53
1:B:265:MET:HB3	1:B:272:LEU:CD2	2.33	0.53
2:E:394:PTR:O	2:E:398:TYR:HD1	1.91	0.53
1:D:42:LYS:HE2	3:D:314:HOH:O	2.09	0.53
1:B:128:ALA:O	1:B:233:ARG:HD2	2.09	0.53
1:A:202:ILE:HD12	1:A:203:ASP:N	2.23	0.53
1:B:243:THR:HG21	1:B:286:VAL:HG22	1.89	0.53
1:A:30:LYS:HA	1:A:33:ARG:HD2	1.90	0.53
1:B:91:LYS:HG3	1:B:268:GLN:OE1	2.07	0.53
1:C:265:MET:HB3	1:C:272:LEU:CD2	2.35	0.53
1:C:51:LYS:O	1:C:55:ILE:HG13	2.09	0.53
1:A:148:GLU:O	1:A:148:GLU:HG2	2.09	0.53
1:B:28:PHE:CD1	1:B:31:LEU:HD12	2.44	0.53
1:D:13:GLU:HG2	1:D:17:LYS:HE2	1.91	0.53
1:B:51:LYS:HB2	1:B:54:ASN:ND2	2.24	0.53
1:A:62:ASP:HB3	2:E:393:GLU:OE2	2.09	0.52
1:A:269:ARG:HG2	1:A:272:LEU:HD13	1.91	0.52
1:A:201:SER:O	1:A:204:PRO:HD2	2.09	0.52
1:B:4:ARG:NH1	1:B:294:MET:HE2	2.24	0.52
2:E:397:LEU:N	2:E:397:LEU:HD23	2.24	0.52
1:B:3:GLN:HE21	1:B:250:GLY:HA2	1.73	0.52
1:D:244:TRP:HB3	3:D:308:HOH:O	2.08	0.52
1:B:86:ASN:HB3	1:B:103:GLN:HG2	1.92	0.52
1:C:17:LYS:HG3	1:C:23:GLU:HG2	1.92	0.52
1:C:12:ASP:HA	1:C:15:GLN:HE21	1.74	0.52
1:B:274:GLN:HB3	1:B:278:GLN:OE1	2.09	0.52
1:D:102:THR:O	1:D:226:HIS:HB2	2.09	0.52
1:D:198:VAL:HG21	3:D:340:HOH:O	2.08	0.52
1:A:84:TYR:CG	1:A:85:ILE:N	2.77	0.51
1:A:190:TYR:OH	1:A:194:PRO:HD3	2.11	0.51
1:C:271:SER:HB3	1:C:274:GLN:OE1	2.10	0.51
1:D:269:ARG:CG	1:D:272:LEU:HD13	2.40	0.51
1:A:117:ILE:HG23	1:A:122:VAL:HB	1.93	0.51
1:A:144:ALA:O	1:A:160:CYS:HB2	2.10	0.51
1:D:93:VAL:CG2	1:D:245:MET:HG2	2.39	0.51
1:C:102:THR:O	1:C:226:HIS:HB2	2.11	0.51
1:A:78:SER:C	1:A:80:GLU:H	2.14	0.51
1:D:60:TYR:CE1	2:H:390:TYR:HE1	2.29	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:390:TYR:C	2:H:390:TYR:CD1	2.84	0.51
1:C:62:ASP:OD1	2:G:393:GLU:HA	2.11	0.51
1:A:89:PHE:CD1	1:A:97:LYS:HG2	2.45	0.51
1:A:6:ILE:HD13	1:A:254:GLU:HA	1.93	0.51
1:C:92:GLY:O	1:D:30:LYS:HE3	2.11	0.51
1:A:22:GLU:HA	3:A:313:HOH:O	2.11	0.51
1:C:166:LYS:HE3	1:C:211:ASP:OD2	2.11	0.51
1:A:34:GLN:HG3	1:A:38:TYR:CE2	2.46	0.51
1:C:77:THR:CG2	1:C:155:PRO:HG3	2.41	0.51
1:C:46:THR:O	1:C:50:GLU:HG3	2.11	0.51
1:A:202:ILE:C	1:A:202:ILE:HD12	2.31	0.50
1:B:197:ASP:O	1:B:199:PRO:HD3	2.11	0.50
1:A:71:VAL:HG22	1:A:89:PHE:CZ	2.46	0.50
1:A:100:ILE:HG22	1:A:100:ILE:O	2.10	0.50
1:B:148:GLU:O	1:B:148:GLU:HG2	2.11	0.50
1:D:37:LYS:HE2	1:D:41:ASP:OD2	2.11	0.50
1:C:246:LEU:HB2	1:C:252:ILE:HD13	1.93	0.50
1:C:235:GLY:HA2	1:C:272:LEU:HG	1.91	0.50
1:C:140:GLU:HG2	3:C:389:HOH:O	2.11	0.50
1:A:246:LEU:HB3	1:A:252:ILE:HD13	1.94	0.50
1:D:7:LEU:HD21	1:D:247:LEU:HD21	1.93	0.50
1:B:100:ILE:HG21	1:B:116:MET:CE	2.41	0.50
1:C:87:ALA:HB2	1:C:102:THR:CB	2.42	0.50
1:B:59:ARG:NH1	2:F:390:TYR:CE2	2.79	0.50
1:B:37:LYS:HE2	1:B:41:ASP:OD2	2.11	0.50
1:C:86:ASN:HB3	1:C:103:GLN:HG2	1.93	0.50
1:B:199:PRO:O	1:B:202:ILE:HG12	2.12	0.50
1:B:240:ILE:O	1:B:244:TRP:HB2	2.12	0.50
1:A:48:VAL:HG21	1:A:70:ARG:HE	1.77	0.50
1:D:269:ARG:HG2	1:D:272:LEU:CD1	2.42	0.50
2:F:390:TYR:O	2:F:391:GLY:C	2.50	0.50
1:C:114:TRP:CZ3	1:C:176:VAL:HG21	2.47	0.49
1:C:203:ASP:HB2	1:C:204:PRO:HD3	1.94	0.49
1:A:90:ILE:HD11	1:A:269:ARG:HB2	1.94	0.49
1:C:1:MET:CE	1:C:5:GLU:HG2	2.40	0.49
1:B:3:GLN:O	1:B:7:LEU:HD13	2.12	0.49
1:B:4:ARG:HH12	1:B:294:MET:HA	1.77	0.49
1:B:140:GLU:HG2	3:B:326:HOH:O	2.11	0.49
1:B:93:VAL:HG21	1:B:245:MET:HG2	1.94	0.49
3:C:367:HOH:O	2:G:398:TYR:HB3	2.13	0.49
1:B:269:ARG:CG	1:B:272:LEU:HD13	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:275:THR:OG1	1:D:278:GLN:HG3	2.13	0.49
1:D:244:TRP:CZ3	1:D:248:LYS:HG3	2.48	0.48
1:B:294:MET:HA	1:B:294:MET:CE	2.44	0.48
1:D:20:THR:HB	1:D:23:GLU:HB2	1.95	0.48
1:A:134:MET:CE	2:E:398:TYR:HB2	2.43	0.48
1:D:93:VAL:HG23	1:D:94:TYR:CD2	2.48	0.48
2:H:397:LEU:O	2:H:398:TYR:OXT	2.32	0.48
1:B:55:ILE:HG22	1:B:55:ILE:O	2.13	0.48
1:B:232:GLY:HA2	1:B:274:GLN:H	1.79	0.48
1:D:195:ASP:HB3	1:D:198:VAL:HB	1.96	0.48
1:C:87:ALA:HB2	1:C:102:THR:OG1	2.14	0.48
1:D:130:MET:HG2	1:D:191:LYS:HG2	1.95	0.48
1:B:244:TRP:O	1:B:244:TRP:HE3	1.97	0.47
1:B:288:GLU:O	1:B:292:ARG:HG3	2.14	0.47
1:A:20:THR:HB	1:A:23:GLU:HB2	1.97	0.47
1:C:195:ASP:O	1:C:196:HIS:C	2.52	0.47
1:B:244:TRP:CZ3	1:B:248:LYS:HG3	2.49	0.47
1:A:105:PRO:O	1:A:140:GLU:HB2	2.14	0.47
1:A:282:VAL:O	1:A:286:VAL:HG23	2.15	0.47
1:C:148:GLU:O	1:C:148:GLU:HG2	2.13	0.47
1:D:86:ASN:HB3	1:D:103:GLN:HG2	1.96	0.47
1:B:46:THR:O	1:B:50:GLU:HG3	2.15	0.47
1:A:24:PHE:CG	1:A:280:GLU:HG2	2.49	0.47
2:E:397:LEU:O	2:E:398:TYR:CD1	2.67	0.47
2:H:393:GLU:CG	2:H:396:ASP:HB2	2.44	0.47
1:D:20:THR:HG22	1:D:21:LYS:N	2.29	0.47
2:H:390:TYR:O	2:H:391:GLY:C	2.52	0.47
1:B:89:PHE:CD1	1:B:97:LYS:HG2	2.50	0.47
1:B:193:TRP:CH2	1:B:233:ARG:HG2	2.50	0.47
1:A:278:GLN:O	1:A:282:VAL:HG23	2.13	0.47
1:A:232:GLY:HA2	1:A:274:GLN:H	1.78	0.47
1:A:60:TYR:CE1	2:E:390:TYR:HE1	2.33	0.47
1:C:13:GLU:HG2	1:C:17:LYS:HE2	1.97	0.47
1:B:57:LYS:HE2	3:B:369:HOH:O	2.14	0.47
1:A:272:LEU:O	1:A:273:VAL:HB	2.15	0.47
1:C:269:ARG:HD2	1:C:270:PRO:HD2	1.96	0.46
1:A:17:LYS:C	1:A:19:ILE:H	2.17	0.46
1:D:244:TRP:CE3	1:D:244:TRP:O	2.65	0.46
1:B:28:PHE:HD1	1:B:31:LEU:HD12	1.81	0.46
1:B:208:LEU:O	1:B:212:VAL:HG23	2.15	0.46
1:A:196:HIS:C	1:A:196:HIS:CD2	2.88	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:LEU:CB	1:A:252:ILE:HD13	2.45	0.46
1:D:274:GLN:HB3	1:D:278:GLN:OE1	2.16	0.46
1:C:145:GLU:O	1:C:148:GLU:HB3	2.15	0.46
1:D:179:ASN:O	1:D:180:SER:HB2	2.15	0.46
1:C:198:VAL:HG12	1:C:201:SER:H	1.81	0.46
1:A:263:ARG:HH21	1:B:95:GLY:HA3	1.78	0.46
1:A:28:PHE:HD1	1:A:31:LEU:HD12	1.79	0.46
1:A:37:LYS:HE2	1:A:41:ASP:OD2	2.16	0.46
2:F:393:GLU:HB3	2:F:396:ASP:H	1.80	0.46
2:E:394:PTR:O	2:E:398:TYR:CD1	2.69	0.46
1:C:43:THR:O	1:C:45:PRO:HD3	2.16	0.46
1:D:232:GLY:HA2	1:D:274:GLN:H	1.81	0.46
1:A:243:THR:HG21	1:A:286:VAL:HG13	1.98	0.46
1:A:225:ILE:HG22	1:A:234:THR:HG23	1.98	0.46
1:C:3:GLN:HE21	1:C:250:GLY:HA2	1.80	0.46
1:C:134:MET:SD	2:G:398:TYR:HB2	2.55	0.46
1:D:164:LYS:HB3	1:D:171:ILE:HB	1.97	0.45
1:C:11:LEU:O	1:C:15:GLN:HG3	2.16	0.45
1:B:2:ASP:OD1	1:B:5:GLU:N	2.43	0.45
1:A:61:LYS:NZ	2:E:391:GLY:O	2.49	0.45
1:A:269:ARG:CG	1:A:272:LEU:HD13	2.46	0.45
1:D:88:ASN:OD1	1:D:269:ARG:HD2	2.16	0.45
1:B:145:GLU:O	1:B:148:GLU:HB3	2.17	0.45
1:A:107:SER:N	1:A:140:GLU:HG3	2.31	0.45
1:B:112:ASP:HA	1:B:115:ARG:HD3	1.97	0.45
1:B:197:ASP:C	1:B:199:PRO:HD3	2.37	0.45
1:C:166:LYS:HD2	1:C:169:TYR:CE1	2.52	0.45
1:A:13:GLU:O	1:A:15:GLN:N	2.49	0.45
1:C:127:MET:HB2	1:C:226:HIS:NE2	2.31	0.45
1:A:78:SER:C	1:A:80:GLU:N	2.69	0.45
1:D:93:VAL:HG12	1:D:242:TYR:CE2	2.52	0.45
1:C:134:MET:CE	2:G:398:TYR:HB2	2.46	0.45
1:A:166:LYS:O	1:C:4:ARG:HD2	2.16	0.45
2:E:396:ASP:O	2:E:397:LEU:HG	2.17	0.45
1:A:17:LYS:C	1:A:19:ILE:N	2.70	0.45
1:D:265:MET:HB3	1:D:272:LEU:CD2	2.45	0.45
1:B:110:LEU:HD21	1:B:140:GLU:HB3	1.99	0.44
1:B:195:ASP:O	1:B:196:HIS:C	2.55	0.44
1:B:89:PHE:CE2	1:B:97:LYS:HE3	2.52	0.44
1:A:133:GLU:HB2	1:A:138:LYS:HG3	1.99	0.44
1:C:59:ARG:NH2	1:C:106:LEU:HD23	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:90:ILE:HB	1:D:99:TYR:HB2	1.98	0.44
1:D:161:GLU:OE2	1:D:175:LYS:HD2	2.18	0.44
1:C:91:LYS:HG2	1:C:96:PRO:HA	1.99	0.44
1:C:3:GLN:NE2	1:C:252:ILE:H	1.96	0.44
2:G:390:TYR:CD1	2:G:390:TYR:C	2.90	0.44
1:D:3:GLN:HE21	1:D:250:GLY:CA	2.28	0.44
1:A:59:ARG:O	2:E:390:TYR:CD1	2.71	0.44
1:C:124:ILE:HG21	1:C:212:VAL:HG13	2.00	0.44
1:D:101:ALA:HB1	1:D:234:THR:HG21	1.99	0.44
1:A:244:TRP:CE3	1:A:248:LYS:HG3	2.53	0.44
1:C:246:LEU:CB	1:C:252:ILE:HD13	2.48	0.44
1:D:124:ILE:HD11	1:D:215:TYR:HB3	2.00	0.44
1:A:3:GLN:HE21	1:A:250:GLY:HA2	1.83	0.43
1:B:244:TRP:CE3	1:B:248:LYS:HG3	2.53	0.43
1:D:243:THR:HG21	1:D:286:VAL:HG13	2.00	0.43
2:E:390:TYR:CD1	2:E:390:TYR:C	2.92	0.43
1:C:282:VAL:O	1:C:286:VAL:HG23	2.18	0.43
1:A:243:THR:HG23	1:A:256:PHE:CZ	2.53	0.43
2:H:395:ASP:O	2:H:397:LEU:HD12	2.19	0.43
1:D:117:ILE:HG23	1:D:122:VAL:HB	2.01	0.43
1:D:30:LYS:HB2	1:D:259:PHE:HE1	1.82	0.43
1:C:77:THR:HG21	1:C:155:PRO:HG3	2.00	0.43
1:C:6:ILE:O	1:C:9:LYS:HB3	2.18	0.43
2:H:393:GLU:HB3	2:H:396:ASP:N	2.29	0.43
1:A:230:GLY:HA2	1:A:234:THR:OG1	2.19	0.43
1:D:202:ILE:O	1:D:202:ILE:HD12	2.19	0.43
2:G:390:TYR:O	2:G:391:GLY:C	2.57	0.43
1:C:235:GLY:CA	1:C:272:LEU:HG	2.49	0.43
2:H:390:TYR:C	2:H:390:TYR:HD1	2.22	0.43
1:B:294:MET:HE2	1:B:294:MET:HA	2.01	0.43
1:C:93:VAL:HG23	1:C:94:TYR:CD2	2.53	0.43
1:A:240:ILE:O	1:A:244:TRP:HB2	2.19	0.42
1:D:141:ARG:HB2	3:D:356:HOH:O	2.19	0.42
1:C:249:ASP:O	1:C:251:ILE:HG23	2.18	0.42
1:B:90:ILE:HB	1:B:99:TYR:HD2	1.85	0.42
1:D:3:GLN:NE2	1:D:250:GLY:C	2.72	0.42
1:B:21:LYS:O	1:B:22:GLU:HB2	2.19	0.42
1:A:142:TYR:C	1:A:142:TYR:CD1	2.93	0.42
1:A:42:LYS:NZ	1:A:42:LYS:HB2	2.35	0.42
1:B:78:SER:C	1:B:80:GLU:H	2.22	0.42
1:D:59:ARG:O	2:H:391:GLY:HA2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:213:ARG:HA	1:D:213:ARG:HD3	1.88	0.42
1:A:164:LYS:NZ	1:C:12:ASP:OD1	2.49	0.42
1:C:211:ASP:O	1:C:214:CYS:HB2	2.19	0.42
1:B:59:ARG:O	2:F:390:TYR:CD1	2.73	0.42
1:B:195:ASP:OD2	1:B:198:VAL:HG23	2.19	0.42
1:A:213:ARG:HD3	1:A:213:ARG:HA	1.88	0.42
1:D:114:TRP:CZ3	1:D:176:VAL:HG21	2.54	0.42
1:D:195:ASP:O	1:D:196:HIS:C	2.57	0.42
1:B:177:LYS:HG2	1:B:182:THR:OG1	2.20	0.42
1:D:9:LYS:HE3	3:D:323:HOH:O	2.19	0.42
1:A:253:PRO:HG3	3:A:320:HOH:O	2.19	0.42
1:C:93:VAL:O	1:D:30:LYS:HD3	2.20	0.42
1:C:266:ARG:HH11	1:C:266:ARG:HG3	1.84	0.42
1:D:120:TYR:O	1:D:121:SER:C	2.59	0.42
1:C:269:ARG:CG	1:C:272:LEU:HD13	2.49	0.41
1:A:91:LYS:HG3	1:A:268:GLN:OE1	2.20	0.41
2:F:390:TYR:CD1	2:F:390:TYR:C	2.93	0.41
1:B:34:GLN:HG3	1:B:38:TYR:CE2	2.55	0.41
1:D:294:MET:HE2	1:D:294:MET:HA	2.01	0.41
1:D:272:LEU:HD12	1:D:272:LEU:HA	1.90	0.41
1:D:202:ILE:HD12	1:D:202:ILE:C	2.41	0.41
1:A:13:GLU:O	1:A:14:ALA:C	2.59	0.41
1:B:133:GLU:HB2	1:B:138:LYS:HG3	2.01	0.41
1:B:193:TRP:CZ3	1:B:233:ARG:HG2	2.55	0.41
1:A:244:TRP:CE3	1:A:244:TRP:O	2.74	0.41
1:B:246:LEU:HB2	1:B:252:ILE:HD13	2.02	0.41
1:D:64:LEU:HA	1:D:65:PRO:HD2	1.91	0.41
1:A:263:ARG:HD2	3:A:304:HOH:O	2.20	0.41
1:A:70:ARG:NH2	1:A:72:GLU:HG2	2.36	0.41
1:D:71:VAL:HG22	1:D:89:PHE:CZ	2.56	0.41
1:D:244:TRP:CE3	1:D:248:LYS:HG3	2.56	0.41
1:D:59:ARG:NH2	1:D:106:LEU:HD23	2.36	0.41
1:C:137:LYS:HG3	1:C:138:LYS:N	2.36	0.41
1:C:87:ALA:HB2	1:C:102:THR:HB	2.01	0.41
2:G:395:ASP:O	2:G:397:LEU:N	2.53	0.41
1:B:247:LEU:HD21	1:B:293:GLN:HG2	2.03	0.41
1:B:27:GLU:HG2	1:B:279:TYR:CZ	2.56	0.41
1:A:151:LEU:HD12	1:A:151:LEU:N	2.36	0.41
1:B:213:ARG:HA	1:B:213:ARG:HD3	1.83	0.41
2:E:396:ASP:C	2:E:397:LEU:HG	2.41	0.41
1:A:114:TRP:CZ3	1:A:176:VAL:HG21	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:51:LYS:O	1:B:55:ILE:HG13	2.21	0.41
1:B:270:PRO:O	1:B:271:SER:HB2	2.20	0.41
1:A:190:TYR:CZ	1:A:192:ASN:O	2.73	0.41
1:C:244:TRP:HE3	1:C:244:TRP:O	2.04	0.41
1:D:100:ILE:HG22	1:D:100:ILE:O	2.20	0.41
1:C:151:LEU:HD12	1:C:151:LEU:N	2.36	0.41
1:D:24:PHE:CG	1:D:280:GLU:HG2	2.55	0.41
1:A:203:ASP:HB2	1:A:204:PRO:HD3	2.03	0.41
1:B:275:THR:OG1	1:B:278:GLN:HG3	2.21	0.41
1:C:190:TYR:OH	1:C:194:PRO:HD3	2.21	0.41
1:D:153:PHE:O	1:D:154:GLY:C	2.58	0.41
1:A:111:LEU:HD13	1:A:154:GLY:HA3	2.03	0.41
1:C:59:ARG:O	2:G:390:TYR:CD1	2.74	0.40
1:B:93:VAL:CG2	1:B:245:MET:HG2	2.51	0.40
1:D:70:ARG:NH1	1:D:72:GLU:HG2	2.36	0.40
1:D:20:THR:HG22	1:D:21:LYS:H	1.85	0.40
1:B:17:LYS:HG2	1:B:23:GLU:HG2	2.03	0.40
1:A:195:ASP:O	1:A:196:HIS:C	2.59	0.40
1:C:20:THR:CG2	1:C:23:GLU:HB2	2.51	0.40
1:C:23:GLU:O	1:C:27:GLU:HB2	2.22	0.40
1:B:100:ILE:HG21	1:B:116:MET:HE3	2.04	0.40
1:A:166:LYS:C	1:C:4:ARG:HD2	2.42	0.40
1:B:124:ILE:HG21	1:B:212:VAL:HG13	2.03	0.40
1:D:91:LYS:HG3	1:D:268:GLN:OE1	2.20	0.40
1:B:43:THR:O	1:B:45:PRO:HD3	2.22	0.40
1:B:64:LEU:O	1:B:270:PRO:HG3	2.21	0.40
1:B:60:TYR:CE1	2:F:390:TYR:HE1	2.39	0.40
1:A:138:LYS:HD2	3:A:296:HOH:O	2.21	0.40
1:B:141:ARG:HA	1:B:143:TRP:CZ3	2.56	0.40
1:C:62:ASP:CG	2:G:393:GLU:HA	2.42	0.40
1:B:93:VAL:HG23	1:B:94:TYR:CD2	2.57	0.40
1:C:244:TRP:CE3	1:C:248:LYS:HG3	2.56	0.40
1:B:179:ASN:O	1:B:180:SER:HB2	2.21	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	295/313 (94%)	259 (88%)	31 (10%)	5 (2%)	11	19
1	B	295/313 (94%)	263 (89%)	29 (10%)	3 (1%)	19	34
1	C	295/313 (94%)	274 (93%)	17 (6%)	4 (1%)	14	24
1	D	295/313 (94%)	277 (94%)	16 (5%)	2 (1%)	26	46
2	E	6/9 (67%)	1 (17%)	2 (33%)	3 (50%)	0	0
2	F	5/9 (56%)	2 (40%)	1 (20%)	2 (40%)	0	0
2	G	6/9 (67%)	0	4 (67%)	2 (33%)	0	0
2	H	6/9 (67%)	2 (33%)	1 (17%)	3 (50%)	0	0
All	All	1203/1288 (93%)	1078 (90%)	101 (8%)	24 (2%)	9	15

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	196	HIS
1	D	196	HIS
2	E	393	GLU
2	E	396	ASP
2	F	393	GLU
2	G	393	GLU
2	H	393	GLU
1	A	14	ALA
1	B	196	HIS
1	C	196	HIS
2	E	391	GLY
2	F	391	GLY
2	G	391	GLY
2	H	391	GLY
2	H	396	ASP
1	A	232	GLY
1	A	273	VAL

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Mol	Chain	Res	Type
1	C	273	VAL
1	D	273	VAL
1	B	273	VAL
1	C	79	ASP
1	B	232	GLY
1	C	232	GLY
1	A	194	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	273/287 (95%)	268 (98%)	5 (2%)	66	88
1	B	273/287 (95%)	263 (96%)	10 (4%)	41	68
1	C	273/287 (95%)	265 (97%)	8 (3%)	50	77
1	D	273/287 (95%)	264 (97%)	9 (3%)	45	73
2	E	7/7 (100%)	4 (57%)	3 (43%)	0	0
2	F	6/7 (86%)	4 (67%)	2 (33%)	0	0
2	G	7/7 (100%)	6 (86%)	1 (14%)	4	7
2	H	7/7 (100%)	5 (71%)	2 (29%)	0	0
All	All	1119/1176 (95%)	1079 (96%)	40 (4%)	42	69

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-3	MET
1	A	79	ASP
1	A	149	MET
1	A	211	ASP
1	A	244	TRP
1	B	-3	MET
1	B	68	TYR
1	B	129	CYS

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Mol	Chain	Res	Type
1	B	149	MET
1	B	211	ASP
1	B	244	TRP
1	B	245	MET
1	B	247	LEU
1	B	272	LEU
1	B	293	GLN
1	C	68	TYR
1	C	149	MET
1	C	203	ASP
1	C	211	ASP
1	C	244	TRP
1	C	247	LEU
1	C	262	ILE
1	C	266	ARG
1	D	-3	MET
1	D	8	GLN
1	D	79	ASP
1	D	141	ARG
1	D	149	MET
1	D	244	TRP
1	D	245	MET
1	D	272	LEU
1	D	293	GLN
2	E	390	TYR
2	E	396	ASP
2	E	397	LEU
2	F	390	TYR
2	F	396	ASP
2	G	390	TYR
2	H	390	TYR
2	H	398	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	8	GLN
1	A	15	GLN
1	A	293	GLN
1	B	3	GLN
1	B	8	GLN

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Mol	Chain	Res	Type
1	B	15	GLN
1	B	54	ASN
1	B	196	HIS
1	B	293	GLN
1	C	3	GLN
1	C	8	GLN
1	C	15	GLN
1	C	293	GLN
1	D	3	GLN
1	D	8	GLN
1	D	15	GLN
1	D	293	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PTR	E	394	2	14,16,17	0.90	0	18,22,24	0.71	0
2	PTR	F	394	2	14,16,17	0.95	0	18,22,24	0.73	0
2	PTR	G	394	2	14,16,17	0.86	0	18,22,24	0.74	0
2	PTR	H	394	2	14,16,17	0.95	0	18,22,24	0.78	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
2	PTR	E	394	2	-	0/9/11/13	0/1/1/1
2	PTR	F	394	2	-	0/9/11/13	0/1/1/1
2	PTR	G	394	2	-	0/9/11/13	0/1/1/1
2	PTR	H	394	2	-	0/9/11/13	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	394	PTR	2	0
2	G	394	PTR	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	297/313 (94%)	-0.14	6 (2%) 68 72	17, 32, 53, 71	0
1	B	297/313 (94%)	-0.13	7 (2%) 62 66	14, 31, 58, 65	0
1	C	297/313 (94%)	-0.24	6 (2%) 68 72	12, 26, 51, 74	0
1	D	297/313 (94%)	-0.27	5 (1%) 73 76	11, 27, 50, 70	0
2	E	8/9 (88%)	1.32	1 (12%) 5 5	35, 38, 41, 42	0
2	F	7/9 (77%)	1.45	2 (28%) 1 0	35, 37, 40, 41	0
2	G	8/9 (88%)	1.36	3 (37%) 0 0	35, 38, 42, 43	0
2	H	8/9 (88%)	1.56	2 (25%) 1 1	33, 38, 41, 42	0
All	All	1219/1288 (94%)	-0.15	32 (2%) 59 63	11, 30, 53, 74	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	197	ASP	4.7
2	H	391	GLY	3.9
2	F	397	LEU	3.9
1	D	197	ASP	3.7
1	C	-3	MET	3.6
1	A	1	MET	3.5
2	E	397	LEU	3.4
1	B	198	VAL	3.4
1	B	196	HIS	3.2
1	B	20	THR	3.0
1	B	19	ILE	2.9
1	B	-2	ALA	2.9
2	F	391	GLY	2.9
1	C	79	ASP	2.8
1	D	196	HIS	2.7
1	C	196	HIS	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	198	VAL	2.6
2	H	390	TYR	2.6
1	C	101	ALA	2.5
1	A	196	HIS	2.5
1	A	197	ASP	2.4
1	D	55	ILE	2.3
2	G	397	LEU	2.3
1	A	79	ASP	2.3
1	D	195	ASP	2.2
2	G	396	ASP	2.2
2	G	391	GLY	2.2
1	A	2	ASP	2.1
1	C	19	ILE	2.1
1	B	79	ASP	2.1
1	D	78	SER	2.0
1	A	199	PRO	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	LLDF	B-factors(\AA^2)	Q<0.9
2	PTR	E	394	16/17	0.94	0.12	-	23,27,34,34	0
2	PTR	F	394	16/17	0.97	0.09	-	22,31,34,34	0
2	PTR	G	394	16/17	0.94	0.12	-	24,26,30,31	0
2	PTR	H	394	16/17	0.95	0.12	-	16,18,29,31	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.