



Full wwPDB X-ray Structure Validation Report ⓘ

May 5, 2016 – 06:04 PM EDT

PDB ID : 5AEX
Title : Crystal structure of *Saccharomyces cerevisiae* Mep2
Authors : Rutherford, J.C.; Chembath, A.; van den Berg, B.
Deposited on : 2015-01-12
Resolution : 3.20 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027457
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) : rb-20027457

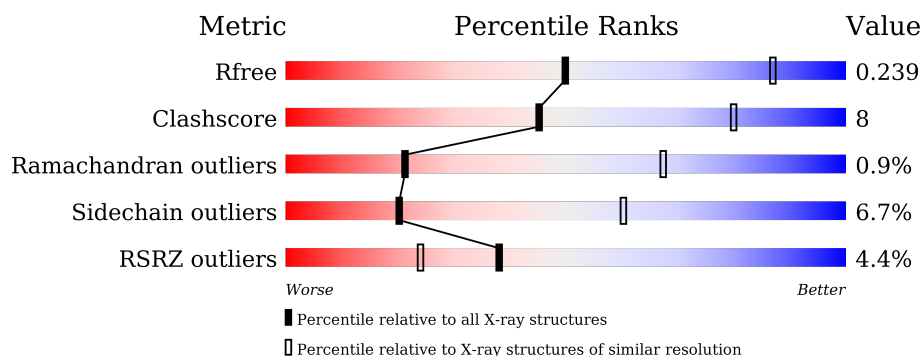
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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 $\leq 5\%$. 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	505	<div> <div>73%</div> <div>15%</div> <div>•</div> <div>10%</div> </div>
1	B	505	<div> <div>3%</div> <div>66%</div> <div>23%</div> <div>•</div> <div>10%</div> </div>
1	C	505	<div> <div>4%</div> <div>70%</div> <div>18%</div> <div>•</div> <div>10%</div> </div>
1	D	505	<div> <div>3%</div> <div>70%</div> <div>17%</div> <div>•</div> <div>10%</div> </div>
1	E	505	<div> <div>5%</div> <div>66%</div> <div>22%</div> <div>•</div> <div>10%</div> </div>
1	F	505	<div> <div>2%</div> <div>71%</div> <div>18%</div> <div>•</div> <div>10%</div> </div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	H	505	
1	I	505	
1	J	505	

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
2	PO4	D	1458	-	-	-	X
2	PO4	E	1458	-	-	-	X
2	PO4	I	1455	-	-	X	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 30576 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 AMMONIUM TRANSPORTER MEP2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	456	Total	C	N	O	S	41	0	0
			3437	2242	561	608	26			
1	B	456	Total	C	N	O	S	65	0	0
			3437	2242	561	608	26			
1	C	456	Total	C	N	O	S	69	0	0
			3437	2242	561	608	26			
1	D	455	Total	C	N	O	S	60	0	0
			3431	2239	560	606	26			
1	E	456	Total	C	N	O	S	95	0	0
			3437	2242	561	608	26			
1	F	456	Total	C	N	O	S	57	0	0
			3437	2242	561	608	26			
1	H	440	Total	C	N	O	S	117	0	0
			3320	2175	541	578	26			
1	I	447	Total	C	N	O	S	204	0	0
			3375	2206	549	594	26			
1	J	427	Total	C	N	O	S	258	0	0
			3220	2115	519	561	25			

There are 54 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	500	HIS	-	EXPRESSION TAG	UNP P41948
A	501	HIS	-	EXPRESSION TAG	UNP P41948
A	502	HIS	-	EXPRESSION TAG	UNP P41948
A	503	HIS	-	EXPRESSION TAG	UNP P41948
A	504	HIS	-	EXPRESSION TAG	UNP P41948
A	505	HIS	-	EXPRESSION TAG	UNP P41948
B	500	HIS	-	EXPRESSION TAG	UNP P41948
B	501	HIS	-	EXPRESSION TAG	UNP P41948
B	502	HIS	-	EXPRESSION TAG	UNP P41948
B	503	HIS	-	EXPRESSION TAG	UNP P41948
B	504	HIS	-	EXPRESSION TAG	UNP P41948

Continued on next page...

Continued from previous page...

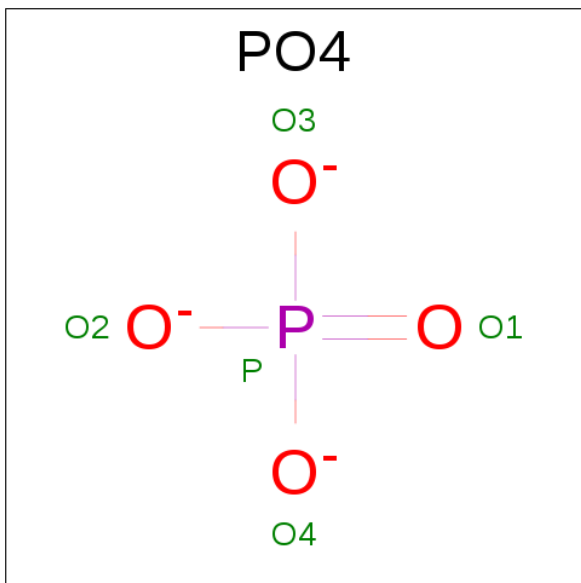
Chain	Residue	Modelled	Actual	Comment	Reference
B	505	HIS	-	EXPRESSION TAG	UNP P41948
C	500	HIS	-	EXPRESSION TAG	UNP P41948
C	501	HIS	-	EXPRESSION TAG	UNP P41948
C	502	HIS	-	EXPRESSION TAG	UNP P41948
C	503	HIS	-	EXPRESSION TAG	UNP P41948
C	504	HIS	-	EXPRESSION TAG	UNP P41948
C	505	HIS	-	EXPRESSION TAG	UNP P41948
D	500	HIS	-	EXPRESSION TAG	UNP P41948
D	501	HIS	-	EXPRESSION TAG	UNP P41948
D	502	HIS	-	EXPRESSION TAG	UNP P41948
D	503	HIS	-	EXPRESSION TAG	UNP P41948
D	504	HIS	-	EXPRESSION TAG	UNP P41948
D	505	HIS	-	EXPRESSION TAG	UNP P41948
E	500	HIS	-	EXPRESSION TAG	UNP P41948
E	501	HIS	-	EXPRESSION TAG	UNP P41948
E	502	HIS	-	EXPRESSION TAG	UNP P41948
E	503	HIS	-	EXPRESSION TAG	UNP P41948
E	504	HIS	-	EXPRESSION TAG	UNP P41948
E	505	HIS	-	EXPRESSION TAG	UNP P41948
F	500	HIS	-	EXPRESSION TAG	UNP P41948
F	501	HIS	-	EXPRESSION TAG	UNP P41948
F	502	HIS	-	EXPRESSION TAG	UNP P41948
F	503	HIS	-	EXPRESSION TAG	UNP P41948
F	504	HIS	-	EXPRESSION TAG	UNP P41948
F	505	HIS	-	EXPRESSION TAG	UNP P41948
H	500	HIS	-	EXPRESSION TAG	UNP P41948
H	501	HIS	-	EXPRESSION TAG	UNP P41948
H	502	HIS	-	EXPRESSION TAG	UNP P41948
H	503	HIS	-	EXPRESSION TAG	UNP P41948
H	504	HIS	-	EXPRESSION TAG	UNP P41948
H	505	HIS	-	EXPRESSION TAG	UNP P41948
I	500	HIS	-	EXPRESSION TAG	UNP P41948
I	501	HIS	-	EXPRESSION TAG	UNP P41948
I	502	HIS	-	EXPRESSION TAG	UNP P41948
I	503	HIS	-	EXPRESSION TAG	UNP P41948
I	504	HIS	-	EXPRESSION TAG	UNP P41948
I	505	HIS	-	EXPRESSION TAG	UNP P41948
J	500	HIS	-	EXPRESSION TAG	UNP P41948
J	501	HIS	-	EXPRESSION TAG	UNP P41948
J	502	HIS	-	EXPRESSION TAG	UNP P41948
J	503	HIS	-	EXPRESSION TAG	UNP P41948
J	504	HIS	-	EXPRESSION TAG	UNP P41948

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
J	505	HIS	-	EXPRESSION TAG	UNP P41948

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).

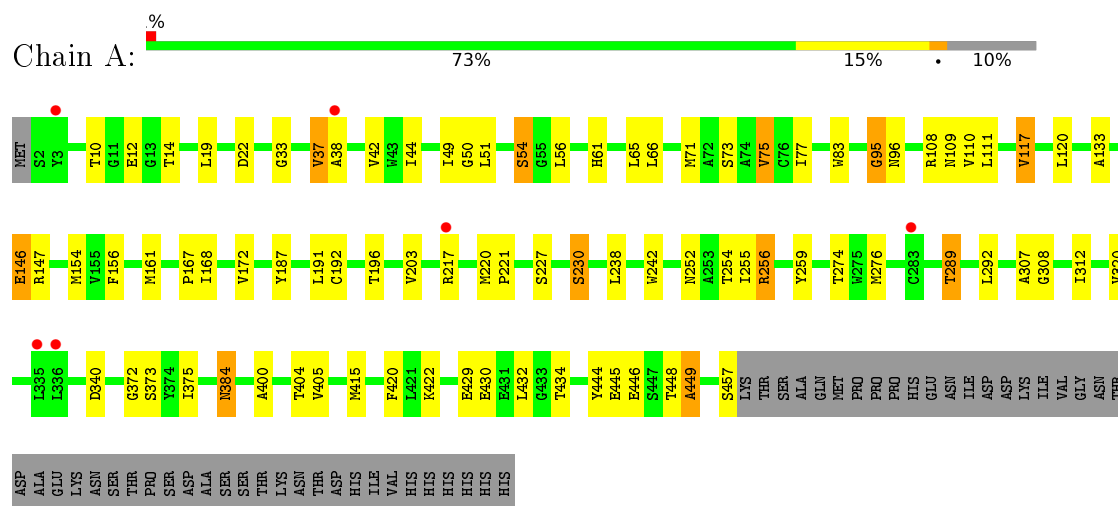


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	F	1	Total	O	P	0	0
			5	4	1		
2	H	1	Total	O	P	0	0
			5	4	1		
2	I	1	Total	O	P	0	0
			5	4	1		
2	J	1	Total	O	P	0	0
			5	4	1		

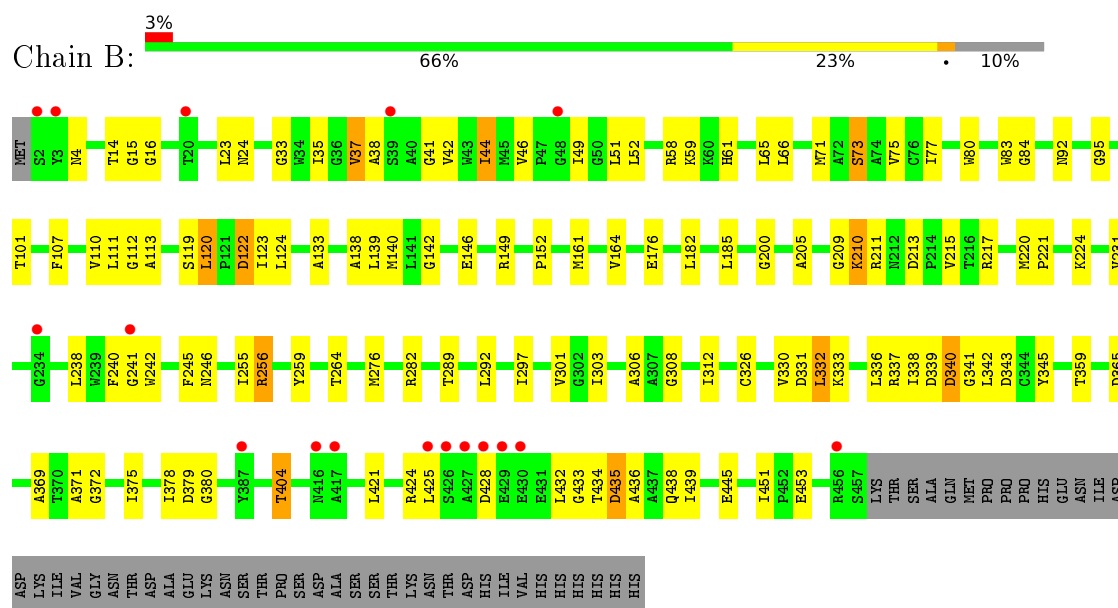
3 Residue-property plots

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: AMMONIUM TRANSPORTER MEP2

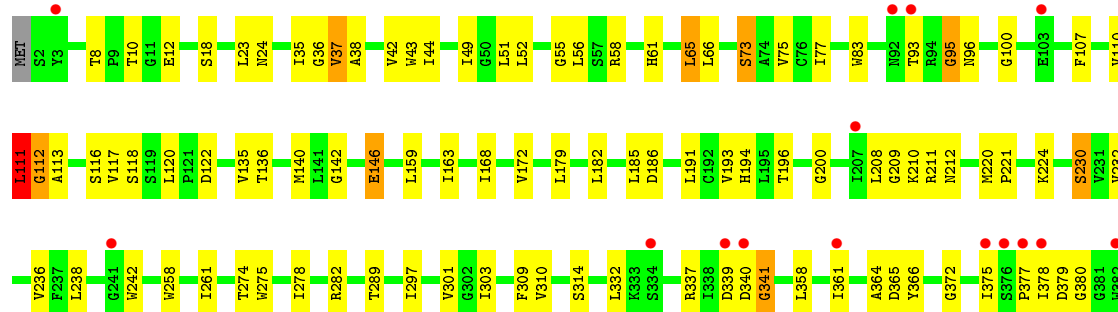


• Molecule 1: AMMONIUM TRANSPORTER MEP2

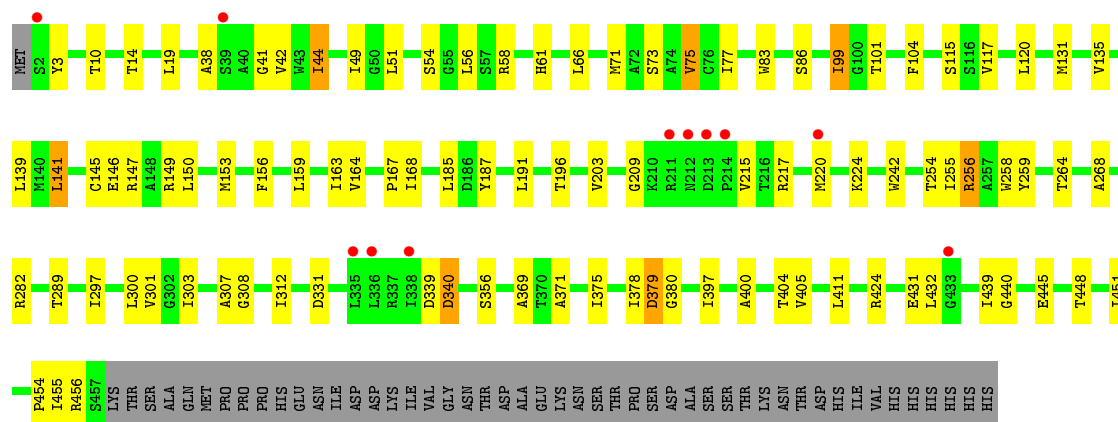


• Molecule 1: AMMONIUM TRANSPORTER MEP2

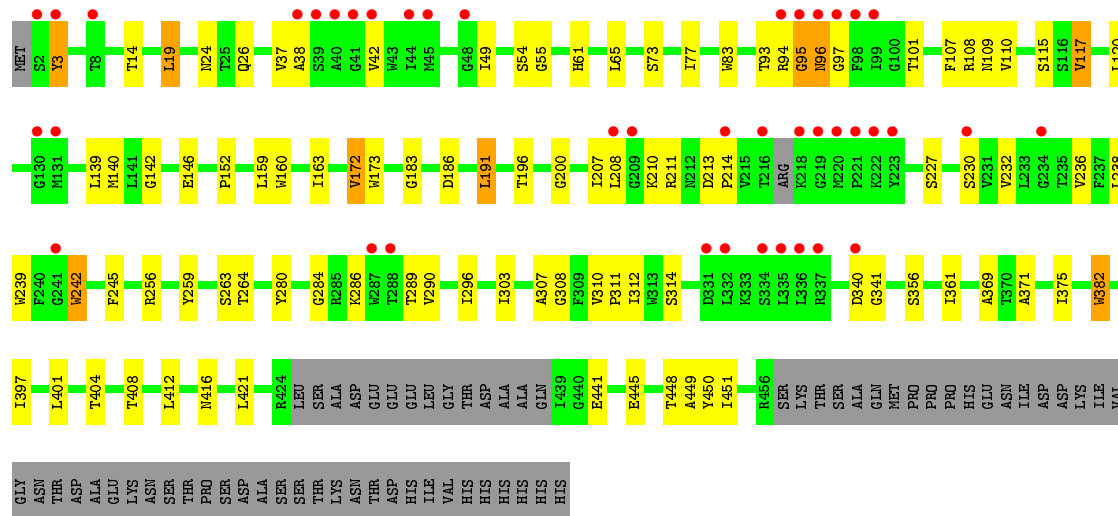




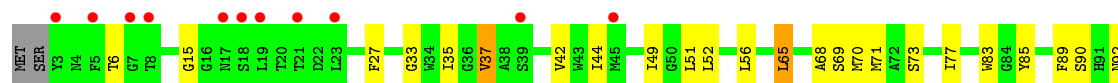
- Molecule 1: AMMONIUM TRANSPORTER MEP2

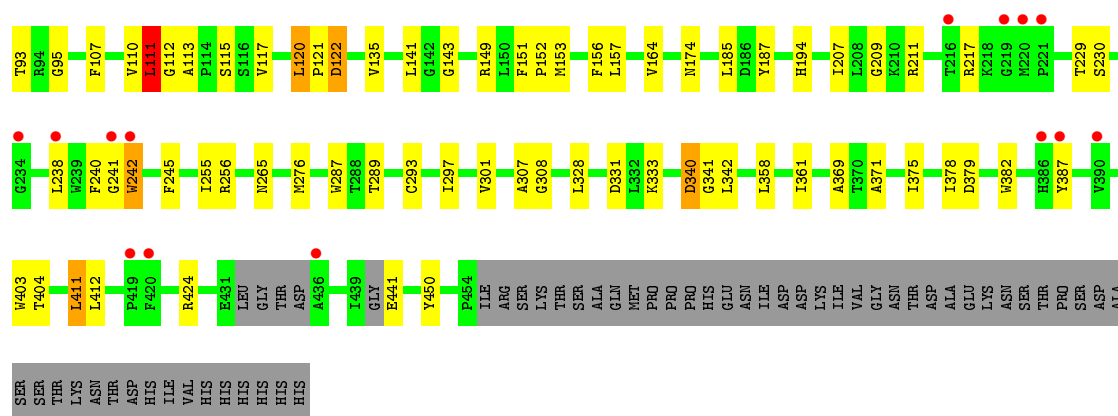


- Molecule 1: AMMONIUM TRANSPORTER MEP2

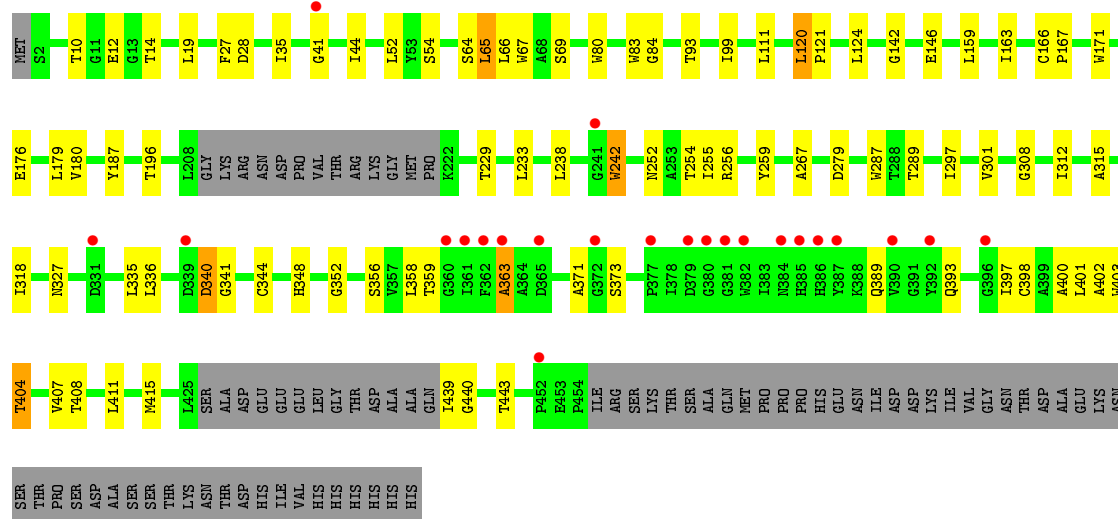


- Molecule 1: AMMONIUM TRANSPORTER MEP2





• Molecule 1: AMMONIUM TRANSPORTER MEP2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	103.69Å 232.35Å 279.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.28 – 3.20 49.28 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.28-3.20) 95.6 (49.28-3.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 3.19Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.189 , 0.241 0.185 , 0.239	Depositor DCC
R_{free} test set	2343 reflections (2.19%)	DCC
Wilson B-factor (Å ²)	86.5	Xtriage
Anisotropy	0.265	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 46.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	30576	wwPDB-VP
Average B, all atoms (Å ²)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.35% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: PO4

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	0.72	1/3534 (0.0%)	0.62	1/4818 (0.0%)
1	B	0.64	0/3534	0.61	2/4818 (0.0%)
1	C	0.54	0/3534	0.57	0/4818
1	D	0.65	0/3528	0.60	2/4810 (0.0%)
1	E	0.57	0/3534	0.59	1/4818 (0.0%)
1	F	0.69	0/3534	0.60	0/4818
1	H	0.53	0/3415	0.53	0/4654
1	I	0.41	0/3470	0.51	1/4729 (0.0%)
1	J	0.37	0/3313	0.48	0/4518
All	All	0.58	1/31396 (0.0%)	0.57	7/42801 (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
1	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	117	VAL	CB-CG1	-5.34	1.41	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	111	LEU	CA-CB-CG	6.78	130.89	115.30
1	A	120	LEU	CA-CB-CG	5.57	128.11	115.30
1	B	340	ASP	N-CA-C	5.55	125.98	111.00

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	336	LEU	CA-CB-CG	5.45	127.83	115.30
1	B	241	GLY	N-CA-C	-5.29	99.89	113.10
1	I	111	LEU	CA-CB-CG	5.25	127.38	115.30
1	D	120	LEU	CA-CB-CG	5.01	126.82	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	240	PHE	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	3437	0	3401	37	0
1	B	3437	0	3401	69	0
1	C	3437	0	3401	53	1
1	D	3431	0	3396	52	0
1	E	3437	0	3401	63	0
1	F	3437	0	3401	47	0
1	H	3320	0	3295	48	0
1	I	3375	0	3337	53	1
1	J	3220	0	3188	51	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
2	E	5	0	0	1	0
2	F	5	0	0	0	0
2	H	5	0	0	0	0
2	I	5	0	0	2	0
2	J	5	0	0	0	0
All	All	30576	0	30221	454	1

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 8.

All (454) 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:J:10:THR:HG22	1:J:12:GLU:H	1.37	0.88
1:F:256:ARG:NH1	1:F:308:GLY:O	2.10	0.84
1:A:10:THR:HG22	1:A:12:GLU:H	1.47	0.78
1:C:400:ALA:O	1:C:404:THR:OG1	2.01	0.78
1:H:49:ILE:HG13	1:H:238:LEU:HD21	1.65	0.78
1:B:205:ALA:HB2	1:B:341:GLY:HA2	1.67	0.77
1:B:378:ILE:HG13	1:B:380:GLY:H	1.49	0.76
1:I:256:ARG:NH1	1:I:308:GLY:O	2.19	0.76
1:D:342:LEU:O	1:D:346:SER:OG	2.02	0.76
1:H:95:GLY:HA3	1:H:97:GLY:N	2.02	0.75
1:E:220:MET:HG3	1:E:221:PRO:HD2	1.68	0.74
1:J:400:ALA:O	1:J:404:THR:OG1	2.04	0.74
1:D:110:VAL:HG23	1:D:122:ASP:HB3	1.70	0.72
1:D:447:SER:HB3	1:E:224:LYS:HG2	1.73	0.71
1:D:220:MET:HG3	1:D:221:PRO:HD2	1.73	0.70
1:B:333:LYS:HG2	1:B:342:LEU:HD12	1.74	0.70
1:H:448:THR:O	1:H:450:TYR:N	2.24	0.70
1:B:256:ARG:NH1	1:B:308:GLY:O	2.24	0.69
1:I:68:ALA:O	1:I:71:MET:N	2.25	0.69
1:F:86:SER:OG	1:F:101:THR:O	2.09	0.69
1:D:369:ALA:HA	1:D:375:ILE:HG21	1.74	0.69
1:E:49:ILE:HG13	1:E:238:LEU:HD21	1.74	0.69
1:C:220:MET:HG3	1:C:221:PRO:HD2	1.75	0.69
1:F:400:ALA:O	1:F:404:THR:HG23	1.93	0.69
1:C:367:VAL:O	1:C:370:THR:OG1	2.12	0.67
1:E:372:GLY:O	1:E:375:ILE:HG13	1.93	0.67
1:H:61:HIS:CD2	1:H:146:GLU:HG2	2.30	0.67
1:J:252:ASN:OD1	1:J:254:THR:OG1	2.11	0.67
1:E:10:THR:HG22	1:E:12:GLU:H	1.59	0.66
1:H:3:TYR:OH	1:H:19:LEU:O	2.14	0.65
1:B:340:ASP:N	1:B:341:GLY:HA3	2.11	0.65
1:I:340:ASP:OD1	1:I:340:ASP:N	2.30	0.65
1:B:220:MET:HG3	1:B:221:PRO:HD2	1.79	0.64
1:F:282:ARG:NH1	1:F:331:ASP:OD2	2.27	0.64
1:H:55:GLY:HA3	1:H:230:SER:HB3	1.80	0.63
1:A:252:ASN:OD1	1:A:254:THR:OG1	2.09	0.63
1:B:58:ARG:NH2	1:B:339:ASP:OD2	2.23	0.63
1:B:65:LEU:HD21	1:B:142:GLY:HA2	1.81	0.63
1:H:451:ILE:HG12	1:I:287:TRP:HB2	1.82	0.62
1:E:73:SER:O	1:E:77:ILE:HG12	1.99	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:256:ARG:NH1	1:D:308:GLY:O	2.33	0.62
1:F:378:ILE:HG13	1:F:380:GLY:H	1.63	0.62
1:B:110:VAL:HG23	1:B:122:ASP:HB3	1.81	0.62
1:I:209:GLY:HA2	1:I:424:ARG:NH1	2.14	0.62
1:A:340:ASP:OD1	1:A:340:ASP:N	2.32	0.62
1:B:133:ALA:HB1	1:B:161:MET:HE1	1.82	0.62
1:I:65:LEU:HA	1:I:68:ALA:HB3	1.80	0.62
1:D:340:ASP:N	1:D:340:ASP:OD1	2.33	0.61
1:B:438:GLN:O	1:B:445:GLU:N	2.29	0.61
1:F:297:ILE:O	1:F:301:VAL:HG23	2.01	0.61
1:C:49:ILE:HG13	1:C:238:LEU:HD21	1.82	0.61
2:I:1455:PO4:O4	1:J:229:THR:OG1	2.18	0.61
1:C:340:ASP:OD1	1:C:340:ASP:N	2.29	0.60
1:F:61:HIS:ND1	1:F:145:CYS:SG	2.74	0.60
1:B:49:ILE:HD13	1:B:139:LEU:HG	1.83	0.60
1:E:110:VAL:HG23	1:E:122:ASP:HB3	1.83	0.60
1:J:80:TRP:HA	1:J:84:GLY:HA3	1.83	0.60
1:E:65:LEU:HD21	1:E:142:GLY:HA2	1.83	0.60
1:F:163:ILE:HD11	1:F:411:LEU:HD13	1.83	0.60
1:D:371:ALA:HB3	1:D:375:ILE:HG22	1.84	0.59
1:A:71:MET:O	1:A:75:VAL:HG13	2.02	0.59
1:C:196:THR:HA	1:C:404:THR:HG21	1.84	0.59
1:I:369:ALA:HA	1:I:375:ILE:HG21	1.83	0.59
1:H:65:LEU:HD11	1:H:142:GLY:HA2	1.84	0.59
1:C:209:GLY:HA2	1:C:424:ARG:NH1	2.16	0.59
1:H:159:LEU:O	1:H:163:ILE:HG12	2.03	0.58
1:J:352:GLY:O	1:J:356:SER:OG	2.20	0.58
1:B:46:VAL:HG13	1:B:138:ALA:HB2	1.86	0.58
1:D:297:ILE:O	1:D:301:VAL:HG23	2.04	0.58
1:E:196:THR:HA	1:E:404:THR:HG21	1.84	0.58
1:C:149:ARG:HG2	1:C:454:PRO:HG2	1.86	0.57
1:F:71:MET:O	1:F:75:VAL:HG13	2.03	0.57
1:H:196:THR:HA	1:H:404:THR:HG21	1.85	0.57
1:A:196:THR:HG22	1:A:404:THR:HG21	1.85	0.57
1:C:174:ASN:OD1	1:C:175:ALA:N	2.37	0.57
1:H:289:THR:HG23	1:H:290:VAL:HG13	1.86	0.57
1:D:191:LEU:O	1:D:196:THR:HG23	2.05	0.57
1:F:149:ARG:HG2	1:F:454:PRO:HG2	1.87	0.57
1:B:111:LEU:O	1:B:113:ALA:N	2.37	0.57
1:F:369:ALA:HA	1:F:375:ILE:HG21	1.87	0.57
1:B:58:ARG:HG2	1:B:224:LYS:HB3	1.87	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:4:ASN:N	1:B:4:ASN:OD1	2.36	0.56
1:B:38:ALA:O	1:B:42:VAL:HG23	2.06	0.56
1:D:51:LEU:HD21	1:D:66:LEU:HD22	1.85	0.56
1:I:49:ILE:HG13	1:I:238:LEU:HD21	1.86	0.56
1:J:267:ALA:HB2	1:J:315:ALA:HB1	1.87	0.56
1:D:435:ASP:N	1:D:435:ASP:OD1	2.36	0.56
1:F:86:SER:HB2	1:F:104:PHE:HB2	1.87	0.56
1:C:282:ARG:NH2	1:C:331:ASP:OD2	2.34	0.56
1:I:194:HIS:CE1	1:I:242:TRP:HZ2	2.24	0.56
1:F:58:ARG:HG2	1:F:224:LYS:HB3	1.88	0.56
1:H:38:ALA:O	1:H:42:VAL:HG23	2.05	0.56
1:B:15:GLY:HA3	1:B:113:ALA:HB2	1.88	0.56
1:A:444:TYR:HB3	1:A:448:THR:HG21	1.86	0.55
1:E:415:MET:HG2	1:E:421:LEU:HB3	1.88	0.55
1:H:340:ASP:N	1:H:340:ASP:OD1	2.37	0.55
1:I:85:TYR:OH	1:I:122:ASP:OD1	2.21	0.55
1:J:259:TYR:CZ	1:J:312:ILE:HG12	2.42	0.55
1:E:310:VAL:HB	1:E:314:SER:OG	2.07	0.55
1:J:10:THR:HG22	1:J:12:GLU:N	2.15	0.55
1:D:65:LEU:HD21	1:D:142:GLY:HA2	1.87	0.55
1:D:49:ILE:HD13	1:D:139:LEU:HG	1.88	0.55
1:F:99:ILE:HD12	1:F:167:PRO:HG3	1.87	0.55
1:B:213:ASP:O	1:B:337:ARG:NH1	2.40	0.55
1:C:282:ARG:NH1	1:C:328:LEU:HD23	2.22	0.55
1:I:73:SER:O	1:I:77:ILE:HG12	2.07	0.55
1:I:141:LEU:HD11	1:I:157:LEU:HD11	1.89	0.55
1:E:55:GLY:HA3	1:E:230:SER:HB3	1.88	0.54
1:H:186:ASP:OD2	1:H:191:LEU:N	2.37	0.54
1:J:27:PHE:CE2	1:J:121:PRO:HB3	2.42	0.54
1:A:220:MET:HG3	1:A:221:PRO:HD2	1.89	0.54
1:C:453:GLU:HG2	1:C:454:PRO:HD2	1.89	0.54
1:H:160:TRP:HE1	1:H:408:THR:HG1	1.55	0.54
1:H:451:ILE:HG22	1:I:289:THR:HG22	1.90	0.54
1:I:149:ARG:HB2	1:I:152:PRO:HG2	1.89	0.54
1:J:411:LEU:O	1:J:415:MET:HB2	2.07	0.54
1:B:451:ILE:HG12	1:C:287:TRP:HB2	1.88	0.54
1:B:71:MET:O	1:B:75:VAL:HG13	2.07	0.54
1:H:24:ASN:OD1	1:I:255:ILE:HG12	2.07	0.54
1:C:86:SER:OG	1:C:101:THR:O	2.25	0.54
1:C:256:ARG:NH1	1:C:308:GLY:O	2.39	0.54
1:E:209:GLY:HA2	1:E:424:ARG:NH1	2.22	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:41:GLY:O	1:J:44:ILE:HG22	2.07	0.54
1:I:110:VAL:HG23	1:I:122:ASP:HB3	1.90	0.53
1:C:203:VAL:HG21	1:C:405:VAL:HA	1.90	0.53
1:C:431:GLU:HB2	1:C:455:ILE:HG13	1.90	0.53
1:A:372:GLY:O	1:A:375:ILE:HG12	2.08	0.53
1:F:49:ILE:HD13	1:F:139:LEU:HG	1.90	0.53
1:I:143:GLY:HA2	1:I:341:GLY:HA3	1.90	0.53
1:D:41:GLY:O	1:D:44:ILE:HG22	2.08	0.53
1:E:168:ILE:O	1:E:172:VAL:HG23	2.09	0.53
1:I:333:LYS:HG2	1:I:342:LEU:HD12	1.91	0.53
1:I:52:LEU:O	1:I:56:LEU:HB2	2.08	0.53
1:J:340:ASP:N	1:J:340:ASP:OD1	2.41	0.53
1:A:156:PHE:HA	1:A:415:MET:HE1	1.90	0.53
1:B:209:GLY:HA2	1:B:424:ARG:NH1	2.24	0.53
1:H:256:ARG:NH1	1:H:308:GLY:O	2.42	0.53
1:E:51:LEU:HD21	1:E:66:LEU:HD22	1.92	0.52
1:I:333:LYS:HD3	1:I:342:LEU:HB2	1.90	0.52
1:E:111:LEU:O	1:E:113:ALA:N	2.35	0.52
1:F:73:SER:O	1:F:77:ILE:HG12	2.08	0.52
1:E:275:TRP:O	1:E:275:TRP:HD1	1.93	0.52
1:H:191:LEU:O	1:H:196:THR:HG23	2.08	0.52
1:E:10:THR:HG22	1:E:12:GLU:N	2.23	0.52
1:I:89:PHE:HA	1:I:174:ASN:HD22	1.74	0.52
1:I:240:PHE:N	1:I:241:GLY:HA3	2.24	0.52
1:D:397:ILE:O	1:D:401:LEU:HB2	2.10	0.52
1:F:187:TYR:CD2	1:F:307:ALA:HB3	2.44	0.52
1:C:86:SER:HB2	1:C:104:PHE:HB2	1.92	0.52
1:I:361:ILE:O	1:I:382:TRP:HB3	2.10	0.52
1:B:33:GLY:O	1:B:37:VAL:HG12	2.10	0.51
1:E:38:ALA:O	1:E:42:VAL:HG23	2.10	0.51
1:H:42:VAL:HG13	1:H:245:PHE:HD1	1.76	0.51
1:J:65:LEU:HD21	1:J:142:GLY:HA2	1.93	0.51
1:J:393:GLN:O	1:J:397:ILE:HG13	2.10	0.51
1:J:66:LEU:O	1:J:69:SER:OG	2.24	0.51
1:C:14:THR:HG23	1:C:16:GLY:H	1.75	0.51
1:A:187:TYR:CD2	1:A:307:ALA:HB3	2.46	0.51
1:A:289:THR:HG21	1:C:448:THR:O	2.10	0.51
1:A:33:GLY:O	1:A:37:VAL:HG12	2.11	0.51
1:F:378:ILE:HD12	1:F:379:ASP:H	1.75	0.51
1:H:369:ALA:HA	1:H:375:ILE:HG21	1.92	0.51
1:H:361:ILE:HA	1:H:382:TRP:HB2	1.93	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:159:LEU:O	1:J:163:ILE:HG12	2.11	0.51
1:A:422:LYS:HE2	1:A:429:GLU:OE2	2.11	0.51
1:B:369:ALA:HA	1:B:375:ILE:HG21	1.93	0.51
1:C:187:TYR:CD2	1:C:307:ALA:HB3	2.45	0.51
1:I:49:ILE:HD11	1:I:135:VAL:HG13	1.93	0.51
1:H:397:ILE:O	1:H:401:LEU:HB2	2.11	0.50
1:B:326:CYS:O	1:B:330:VAL:HG23	2.12	0.50
1:F:115:SER:OG	1:F:117:VAL:O	2.26	0.50
1:I:111:LEU:O	1:I:113:ALA:N	2.44	0.50
1:J:318:ILE:HA	1:J:358:LEU:HD13	1.93	0.50
1:B:52:LEU:HD12	1:B:231:VAL:HG13	1.92	0.50
1:F:38:ALA:O	1:F:42:VAL:HG23	2.11	0.50
1:J:163:ILE:HD11	1:J:411:LEU:HD22	1.94	0.50
1:C:41:GLY:O	1:C:44:ILE:HG22	2.11	0.50
1:E:191:LEU:O	1:E:196:THR:HG23	2.12	0.50
1:J:256:ARG:HB3	1:J:371:ALA:HB2	1.93	0.50
1:E:384:ASN:O	1:E:386:HIS:ND1	2.44	0.49
1:E:258:TRP:HE3	1:E:261:ILE:HD12	1.78	0.49
1:E:36:GLY:HA3	1:F:258:TRP:CZ3	2.47	0.49
1:A:449:ALA:O	1:B:289:THR:N	2.45	0.49
1:C:35:ILE:HG21	1:C:120:LEU:HD21	1.94	0.49
1:D:62:ALA:HA	1:D:439:ILE:HD11	1.94	0.49
1:H:108:ARG:O	1:H:109:ASN:HB2	2.12	0.49
1:F:445:GLU:O	1:F:448:THR:OG1	2.24	0.49
1:B:59:LYS:NZ	1:B:213:ASP:OD2	2.30	0.49
1:B:80:TRP:HA	1:B:84:GLY:HA3	1.94	0.49
1:C:333:LYS:HE3	1:C:339:ASP:OD1	2.11	0.49
1:D:361:ILE:HA	1:D:382:TRP:HB2	1.94	0.49
1:J:93:THR:OG1	1:J:176:GLU:OE1	2.30	0.49
1:C:73:SER:O	1:C:77:ILE:HG12	2.13	0.49
1:D:187:TYR:CD2	1:D:307:ALA:HB3	2.48	0.49
1:J:54:SER:HB2	1:J:65:LEU:HB2	1.94	0.49
1:F:131:MET:O	1:F:135:VAL:HG23	2.12	0.49
1:C:95:GLY:C	1:C:97:GLY:H	2.15	0.49
1:A:73:SER:O	1:A:77:ILE:HG12	2.13	0.48
1:I:185:LEU:HD11	1:I:378:ILE:HG12	1.95	0.48
1:J:403:TRP:CE2	1:J:407:VAL:HG21	2.49	0.48
1:B:338:ILE:HG22	1:B:339:ASP:H	1.79	0.48
1:E:140:MET:HE3	1:E:200:GLY:HA3	1.95	0.48
1:B:49:ILE:HG13	1:B:238:LEU:HD21	1.95	0.48
1:B:336:LEU:HB3	1:B:338:ILE:HG13	1.95	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:136:THR:OG1	1:D:193:VAL:HG13	2.14	0.48
1:H:371:ALA:HB3	1:H:375:ILE:HG22	1.96	0.48
1:J:187:TYR:CE1	1:J:308:GLY:HA3	2.49	0.48
1:I:340:ASP:HA	1:I:341:GLY:HA2	1.59	0.48
1:F:141:LEU:HD12	1:F:153:MET:SD	2.53	0.48
1:A:256:ARG:NH1	1:A:308:GLY:O	2.47	0.48
1:B:73:SER:O	1:B:77:ILE:HG12	2.14	0.48
1:F:58:ARG:NH2	1:F:339:ASP:OD2	2.38	0.48
1:I:33:GLY:O	1:I:37:VAL:HG12	2.14	0.48
1:A:400:ALA:O	1:A:404:THR:HG23	2.13	0.48
1:I:90:SER:OG	1:I:92:ASN:O	2.32	0.48
1:D:95:GLY:HA2	1:D:96:ASN:HA	1.54	0.47
1:C:110:VAL:HG13	1:C:122:ASP:HB3	1.95	0.47
1:E:403:TRP:CE2	1:E:407:VAL:HG21	2.49	0.47
1:C:213:ASP:OD1	1:C:215:VAL:HG23	2.15	0.47
1:D:73:SER:O	1:D:77:ILE:HG12	2.15	0.47
1:E:93:THR:HA	1:E:100:GLY:HA2	1.95	0.47
1:B:164:VAL:CG1	1:B:404:THR:HG22	2.44	0.47
1:B:297:ILE:O	1:B:301:VAL:HG23	2.14	0.47
1:E:380:GLY:HA3	1:E:384:ASN:HD22	1.79	0.47
1:H:259:TYR:CZ	1:H:312:ILE:HG12	2.49	0.47
1:I:35:ILE:HG13	1:I:120:LEU:HD21	1.95	0.47
1:A:133:ALA:HB1	1:A:161:MET:HE1	1.97	0.47
1:C:202:LEU:HB2	1:C:345:TYR:CD1	2.49	0.47
1:C:49:ILE:HD13	1:C:139:LEU:HG	1.96	0.47
1:J:196:THR:HA	1:J:404:THR:HG21	1.97	0.47
1:B:342:LEU:HA	1:B:345:TYR:HB3	1.96	0.47
1:I:358:LEU:HD23	1:I:361:ILE:HD12	1.97	0.47
1:A:51:LEU:HD21	1:A:66:LEU:HD22	1.97	0.47
1:C:115:SER:OG	1:C:117:VAL:O	2.28	0.47
1:H:49:ILE:HD13	1:H:139:LEU:HG	1.95	0.47
1:B:42:VAL:HG13	1:B:245:PHE:HD1	1.80	0.47
1:E:24:ASN:OD1	1:F:255:ILE:HG13	2.15	0.47
1:B:35:ILE:HG21	1:B:120:LEU:HD21	1.97	0.47
1:B:16:GLY:HA3	1:B:23:LEU:HD21	1.97	0.47
1:D:49:ILE:HG13	1:D:238:LEU:HD21	1.95	0.47
1:E:358:LEU:HD23	1:E:361:ILE:HD12	1.97	0.47
1:E:51:LEU:HG	1:E:66:LEU:HD13	1.97	0.47
1:A:147:ARG:NH2	1:A:430:GLU:OE2	2.41	0.47
1:F:264:THR:HA	1:F:303:ILE:HG21	1.97	0.47
1:B:246:ASN:ND2	1:B:264:THR:OG1	2.25	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:365:ASP:N	1:B:365:ASP:OD1	2.48	0.46
1:C:434:THR:O	1:C:436:ALA:N	2.45	0.46
1:F:150:LEU:HD22	1:F:451:ILE:HD11	1.96	0.46
1:I:256:ARG:HD3	1:I:371:ALA:HB2	1.95	0.46
1:E:441:GLU:N	2:E:1458:PO4:O1	2.48	0.46
1:A:168:ILE:O	1:A:172:VAL:HG12	2.15	0.46
1:E:58:ARG:NH1	1:E:221:PRO:O	2.49	0.46
1:F:146:GLU:HG3	1:F:147:ARG:HG2	1.96	0.46
1:D:187:TYR:CE2	1:D:307:ALA:HB3	2.51	0.46
1:D:434:THR:O	1:D:436:ALA:N	2.48	0.46
1:H:94:ARG:HD3	1:H:101:THR:HG22	1.97	0.46
1:H:238:LEU:O	1:H:242:TRP:HB2	2.16	0.46
1:C:51:LEU:HD21	1:C:66:LEU:HD22	1.97	0.46
1:E:194:HIS:CG	1:E:301:VAL:HG13	2.50	0.46
1:F:164:VAL:O	1:F:168:ILE:HG13	2.15	0.46
1:J:279:ASP:OD1	1:J:327:ASN:ND2	2.36	0.46
1:A:259:TYR:CZ	1:A:312:ILE:HG12	2.50	0.46
1:A:255:ILE:HG13	1:C:24:ASN:OD1	2.16	0.46
1:D:61:HIS:CE1	1:D:146:GLU:HG2	2.51	0.46
1:F:209:GLY:HA2	1:F:424:ARG:NH1	2.31	0.46
1:H:340:ASP:HA	1:H:341:GLY:HA2	1.58	0.46
1:B:65:LEU:HA	1:B:65:LEU:HD12	1.71	0.46
1:J:146:GLU:HG2	1:J:340:ASP:OD2	2.16	0.46
1:B:333:LYS:NZ	1:B:343:ASP:OD1	2.41	0.46
1:D:166:CYS:HB2	1:D:167:PRO:HD3	1.98	0.46
1:E:232:VAL:O	1:E:236:VAL:HG23	2.16	0.46
1:J:64:SER:OG	1:J:443:THR:HA	2.15	0.46
1:A:61:HIS:ND1	1:A:146:GLU:HG2	2.30	0.46
1:B:276:MET:HE3	1:B:292:LEU:HD22	1.98	0.46
1:B:438:GLN:HA	1:B:445:GLU:HB2	1.98	0.46
1:D:174:ASN:OD1	1:D:175:ALA:N	2.49	0.46
1:H:232:VAL:O	1:H:236:VAL:HG23	2.16	0.46
1:H:77:ILE:HG21	1:I:265:ASN:O	2.16	0.46
1:J:52:LEU:HD21	1:J:297:ILE:HG13	1.98	0.46
1:A:61:HIS:CE1	1:A:146:GLU:HG2	2.52	0.45
1:A:49:ILE:HG13	1:A:238:LEU:HD21	1.98	0.45
1:B:282:ARG:NH2	1:B:331:ASP:OD2	2.41	0.45
1:D:217:ARG:HB2	1:D:337:ARG:HH22	1.82	0.45
1:D:164:VAL:CG1	1:D:404:THR:HG22	2.46	0.45
1:E:172:VAL:HG22	1:E:179:LEU:HD12	1.98	0.45
1:E:303:ILE:O	1:E:303:ILE:HG13	2.17	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:203:VAL:HG21	1:F:405:VAL:HA	1.98	0.45
1:J:238:LEU:O	1:J:242:TRP:HB2	2.16	0.45
1:A:227:SER:HB3	1:A:230:SER:HB2	1.98	0.45
1:B:24:ASN:OD1	1:C:255:ILE:HG12	2.16	0.45
1:D:211:ARG:HB3	1:D:211:ARG:HE	1.65	0.45
1:A:276:MET:HE3	1:A:292:LEU:HD22	1.99	0.45
1:H:140:MET:HE3	1:H:200:GLY:HA3	1.99	0.45
1:J:335:LEU:O	1:J:336:LEU:HD12	2.16	0.45
1:D:16:GLY:HA3	1:D:23:LEU:HD21	1.99	0.45
1:E:136:THR:OG1	1:E:193:VAL:HG13	2.16	0.45
1:B:51:LEU:HD21	1:B:66:LEU:HD22	1.98	0.45
1:B:92:ASN:O	1:B:101:THR:HG23	2.17	0.45
1:D:282:ARG:NH2	1:D:331:ASP:OD2	2.34	0.45
1:H:412:LEU:O	1:H:416:ASN:HB2	2.17	0.45
1:I:441:GLU:N	2:I:1455:PO4:O1	2.50	0.45
1:J:35:ILE:HG13	1:J:120:LEU:HD21	1.99	0.45
1:F:340:ASP:N	1:F:340:ASP:OD1	2.49	0.45
1:J:397:ILE:O	1:J:401:LEU:HB2	2.17	0.45
1:E:61:HIS:CE1	1:E:146:GLU:HG2	2.52	0.44
1:F:159:LEU:O	1:F:163:ILE:HG12	2.17	0.44
1:I:238:LEU:O	1:I:242:TRP:N	2.51	0.44
1:D:174:ASN:O	1:D:180:VAL:HG21	2.18	0.44
1:J:340:ASP:HA	1:J:341:GLY:HA2	1.62	0.44
1:C:191:LEU:O	1:C:196:THR:HG23	2.17	0.44
1:D:228:VAL:O	1:D:232:VAL:HG23	2.17	0.44
1:C:210:LYS:O	1:C:338:ILE:HG23	2.18	0.44
1:E:35:ILE:HG21	1:E:120:LEU:HD21	2.00	0.44
1:H:115:SER:OG	1:H:117:VAL:O	2.32	0.44
1:H:95:GLY:HA3	1:H:96:ASN:C	2.38	0.44
1:F:185:LEU:HD11	1:F:378:ILE:HG12	1.99	0.44
1:F:196:THR:HG22	1:F:404:THR:HG21	2.00	0.44
1:I:293:CYS:O	1:I:297:ILE:HG12	2.18	0.44
1:J:389:GLN:O	1:J:393:GLN:HG2	2.18	0.44
1:C:340:ASP:HA	1:C:341:GLY:HA2	1.78	0.44
1:C:163:ILE:HD11	1:C:411:LEU:HD13	2.00	0.44
1:F:371:ALA:HB3	1:F:375:ILE:CG2	2.48	0.44
1:A:95:GLY:HA2	1:A:96:ASN:HA	1.47	0.44
1:B:371:ALA:HB3	1:B:375:ILE:HG22	2.00	0.43
1:C:150:LEU:HD22	1:C:451:ILE:HD11	2.00	0.43
1:I:297:ILE:O	1:I:301:VAL:HG23	2.18	0.43
1:I:187:TYR:CD2	1:I:307:ALA:HB3	2.53	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:LYS:N	1:B:340:ASP:OD2	2.51	0.43
1:B:433:GLY:HA3	1:B:434:THR:HA	1.74	0.43
1:C:439:ILE:HA	1:C:440:GLY:HA3	1.76	0.43
1:D:326:CYS:O	1:D:330:VAL:HG23	2.17	0.43
1:D:37:VAL:HG21	1:E:37:VAL:HG21	1.99	0.43
1:J:439:ILE:HA	1:J:440:GLY:HA3	1.64	0.43
1:C:136:THR:OG1	1:C:193:VAL:HG13	2.19	0.43
1:D:188:ALA:HB3	1:D:304:THR:HG23	2.00	0.43
1:E:274:THR:O	1:E:278:ILE:HG13	2.19	0.43
1:E:411:LEU:O	1:E:415:MET:HB2	2.17	0.43
1:F:51:LEU:HD21	1:F:66:LEU:HD22	2.00	0.43
1:I:164:VAL:HA	1:I:403:TRP:HE1	1.82	0.43
1:B:185:LEU:HD11	1:B:378:ILE:HG12	2.01	0.43
1:I:68:ALA:O	1:I:70:MET:N	2.51	0.43
1:B:140:MET:CE	1:B:200:GLY:HA3	2.47	0.43
1:I:194:HIS:CE1	1:I:301:VAL:HG22	2.53	0.43
1:D:65:LEU:HA	1:D:65:LEU:HD12	1.65	0.43
1:C:238:LEU:O	1:C:242:TRP:HB2	2.18	0.43
1:I:141:LEU:HD12	1:I:153:MET:HE2	2.00	0.43
1:J:233:LEU:HA	1:J:233:LEU:HD23	1.76	0.43
1:B:152:PRO:HB3	1:B:421:LEU:O	2.19	0.43
1:C:185:LEU:N	1:C:393:GLN:OE1	2.45	0.43
1:E:401:LEU:O	1:E:405:VAL:HG23	2.19	0.43
1:B:107:PHE:CD2	1:B:110:VAL:HG11	2.54	0.43
1:I:56:LEU:HA	1:I:56:LEU:HD12	1.82	0.43
1:D:400:ALA:O	1:D:404:THR:OG1	2.20	0.43
1:E:278:ILE:O	1:E:282:ARG:HB2	2.18	0.43
1:E:365:ASP:H	1:E:384:ASN:HD21	1.67	0.43
1:H:172:VAL:HG22	1:H:173:TRP:CD1	2.54	0.43
1:B:123:ILE:HG12	1:C:259:TYR:HA	2.00	0.42
1:C:213:ASP:HA	1:C:214:PRO:HD2	1.80	0.42
1:C:35:ILE:HG13	1:C:120:LEU:HD21	2.00	0.42
1:E:212:ASN:HA	1:E:337:ARG:HB3	2.02	0.42
1:I:194:HIS:CG	1:I:301:VAL:HG13	2.54	0.42
1:A:38:ALA:O	1:A:42:VAL:HG23	2.19	0.42
1:I:151:PHE:HB2	1:J:287:TRP:HZ2	1.84	0.42
1:J:166:CYS:HB2	1:J:167:PRO:HD3	2.02	0.42
1:D:340:ASP:HA	1:D:341:GLY:HA2	1.60	0.42
1:J:171:TRP:O	1:J:180:VAL:HG23	2.20	0.42
1:J:179:LEU:HA	1:J:179:LEU:HD23	1.83	0.42
1:D:87:LEU:HD22	1:D:99:ILE:HD11	2.01	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:339:ASP:O	1:E:341:GLY:N	2.52	0.42
1:A:203:VAL:HG21	1:A:405:VAL:HA	2.02	0.42
1:B:306:ALA:HB2	1:B:359:THR:OG1	2.20	0.42
1:C:365:ASP:H	1:C:384:ASN:HD21	1.68	0.42
1:D:50:GLY:O	1:D:54:SER:HB3	2.20	0.42
1:I:412:LEU:HD23	1:I:412:LEU:HA	1.84	0.42
1:A:384:ASN:HD22	1:A:384:ASN:HA	1.70	0.42
1:B:61:HIS:O	1:B:439:ILE:HD13	2.20	0.42
1:D:140:MET:HE3	1:D:200:GLY:HA3	2.02	0.42
1:E:95:GLY:HA2	1:E:96:ASN:HA	1.53	0.42
1:B:259:TYR:CZ	1:B:312:ILE:HG12	2.55	0.42
1:C:43:TRP:HB2	1:C:131:MET:SD	2.59	0.42
1:C:202:LEU:HD22	1:C:345:TYR:CE1	2.55	0.42
1:E:49:ILE:HD11	1:E:135:VAL:HG13	2.02	0.42
1:F:56:LEU:HD12	1:F:56:LEU:HA	1.83	0.41
1:H:310:VAL:HB	1:H:314:SER:OG	2.20	0.41
1:I:15:GLY:HA3	1:I:113:ALA:HB2	2.02	0.41
1:A:274:THR:OG1	1:A:320:VAL:HG22	2.21	0.41
1:B:149:ARG:HD2	1:B:425:LEU:HG	2.02	0.41
1:E:186:ASP:HB2	1:E:191:LEU:HD13	2.02	0.41
1:F:439:ILE:HA	1:F:440:GLY:HA3	1.83	0.41
1:B:436:ALA:C	1:B:438:GLN:H	2.23	0.41
1:E:185:LEU:HD22	1:E:309:PHE:CZ	2.56	0.41
1:H:280:TYR:HA	1:H:286:LYS:O	2.20	0.41
1:J:120:LEU:HG	1:J:124:LEU:HD23	2.02	0.41
1:A:22:ASP:OD1	1:B:372:GLY:HA3	2.19	0.41
1:D:140:MET:HE2	1:D:140:MET:HB2	1.99	0.41
1:E:23:LEU:HD13	1:E:112:GLY:HA3	2.03	0.41
1:E:140:MET:HE2	1:E:140:MET:HB2	1.93	0.41
1:F:41:GLY:O	1:F:44:ILE:HG22	2.21	0.41
1:H:264:THR:HA	1:H:303:ILE:HG21	2.02	0.41
1:I:328:LEU:O	1:I:331:ASP:N	2.51	0.41
1:J:359:THR:O	1:J:363:ALA:HB2	2.19	0.41
1:A:50:GLY:O	1:A:54:SER:HB3	2.20	0.41
1:D:57:SER:OG	1:D:59:LYS:O	2.36	0.41
1:E:297:ILE:O	1:E:301:VAL:HG23	2.19	0.41
1:F:153:MET:CE	1:F:156:PHE:HD2	2.32	0.41
1:F:259:TYR:CZ	1:F:312:ILE:HG12	2.55	0.41
1:F:371:ALA:HB3	1:F:375:ILE:HG22	2.03	0.41
1:I:27:PHE:CE2	1:I:121:PRO:HB3	2.55	0.41
1:J:65:LEU:HD12	1:J:65:LEU:HA	1.78	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:56:LEU:HA	1:D:56:LEU:HD12	1.88	0.41
1:E:208:LEU:HD13	1:E:424:ARG:HB3	2.02	0.41
1:H:152:PRO:HB3	1:H:421:LEU:O	2.19	0.41
1:J:398:CYS:O	1:J:402:ALA:N	2.46	0.41
1:E:159:LEU:O	1:E:163:ILE:HG12	2.20	0.41
1:E:443:THR:OG1	1:E:444:TYR:N	2.51	0.41
1:H:289:THR:O	1:J:67:TRP:NE1	2.38	0.41
1:B:332:LEU:HD23	1:B:342:LEU:HD13	2.03	0.41
1:H:263:SER:OG	1:H:312:ILE:HD13	2.20	0.41
1:I:42:VAL:HG13	1:I:245:PHE:HD1	1.86	0.41
1:A:71:MET:SD	1:A:154:MET:HG2	2.61	0.41
1:D:289:THR:HG21	1:F:448:THR:O	2.21	0.41
1:D:425:LEU:HB2	1:D:430:GLU:HG2	2.02	0.41
1:D:328:LEU:O	1:D:330:VAL:N	2.54	0.41
1:D:453:GLU:HG2	1:D:454:PRO:HD2	2.02	0.41
1:D:71:MET:O	1:D:75:VAL:HG13	2.21	0.41
1:E:375:ILE:HG22	1:E:377:PRO:HD3	2.03	0.41
1:J:99:ILE:HG13	1:J:167:PRO:HA	2.03	0.41
1:B:264:THR:HA	1:B:303:ILE:HG21	2.02	0.41
1:I:361:ILE:HA	1:I:382:TRP:HB2	2.02	0.41
1:I:51:LEU:O	1:I:230:SER:OG	2.32	0.41
1:C:188:ALA:HB3	1:C:304:THR:HG23	2.02	0.40
1:E:52:LEU:O	1:E:56:LEU:HB2	2.21	0.40
1:E:43:TRP:NE1	1:E:73:SER:OG	2.40	0.40
1:F:356:SER:HB2	1:F:397:ILE:HD11	2.04	0.40
1:H:356:SER:HB2	1:H:397:ILE:HD11	2.03	0.40
1:I:156:PHE:CD1	1:I:411:LEU:HD11	2.56	0.40
1:J:35:ILE:HB	1:J:124:LEU:HD21	2.02	0.40
1:B:220:MET:CG	1:B:221:PRO:HD2	2.49	0.40
1:E:378:ILE:HA	1:E:378:ILE:HD12	1.91	0.40
1:H:239:TRP:CZ3	1:H:296:ILE:HD13	2.56	0.40
1:B:41:GLY:O	1:B:44:ILE:HG22	2.22	0.40
1:B:124:LEU:HD13	1:C:258:TRP:CG	2.56	0.40
1:E:435:ASP:HB3	1:E:438:GLN:OE1	2.21	0.40
1:H:310:VAL:HA	1:H:311:PRO:HD3	1.89	0.40
1:J:344:CYS:O	1:J:348:HIS:HB2	2.21	0.40
1:C:64:SER:OG	1:C:443:THR:HA	2.21	0.40
1:J:404:THR:O	1:J:408:THR:OG1	2.29	0.40
1:A:108:ARG:O	1:A:109:ASN:HB2	2.21	0.40
1:B:107:PHE:O	1:B:110:VAL:HG12	2.21	0.40
1:E:364:ALA:HB1	1:E:366:TYR:CE1	2.57	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:268:ALA:HB2	1:F:300:LEU:HG	2.03	0.40
1:H:207:ILE:O	1:H:208:LEU:HD23	2.20	0.40
1:H:227:SER:HB2	1:H:441:GLU:HG2	2.04	0.40
1:J:297:ILE:O	1:J:301:VAL:HG23	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:12:GLU:OE2	1:I:387:TYR:OH[4_545]	2.19	0.01

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	454/505 (90%)	429 (94%)	21 (5%)	4 (1%)	21	67
1	B	454/505 (90%)	422 (93%)	29 (6%)	3 (1%)	26	72
1	C	454/505 (90%)	421 (93%)	29 (6%)	4 (1%)	21	67
1	D	453/505 (90%)	419 (92%)	29 (6%)	5 (1%)	17	62
1	E	454/505 (90%)	421 (93%)	28 (6%)	5 (1%)	17	62
1	F	454/505 (90%)	429 (94%)	24 (5%)	1 (0%)	52	88
1	H	434/505 (86%)	406 (94%)	20 (5%)	8 (2%)	11	51
1	I	441/505 (87%)	402 (91%)	34 (8%)	5 (1%)	17	62
1	J	421/505 (83%)	387 (92%)	32 (8%)	2 (0%)	34	78
All	All	4019/4545 (88%)	3736 (93%)	246 (6%)	37 (1%)	21	67

All (37) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	112	GLY
1	D	435	ASP
1	I	112	GLY
1	A	95	GLY
1	B	95	GLY
1	C	372	GLY
1	C	433	GLY
1	D	95	GLY
1	E	95	GLY
1	E	112	GLY
1	E	340	ASP
1	H	445	GLU
1	H	449	ALA
1	I	69	SER
1	J	363	ALA
1	A	111	LEU
1	A	449	ALA
1	C	435	ASP
1	D	329	ALA
1	C	447	SER
1	D	445	GLU
1	H	307	ALA
1	I	115	SER
1	J	111	LEU
1	A	446	GLU
1	B	435	ASP
1	D	111	LEU
1	E	341	GLY
1	E	432	LEU
1	F	254	THR
1	H	95	GLY
1	H	96	ASN
1	H	284	GLY
1	H	214	PRO
1	H	183	GLY
1	I	95	GLY
1	I	207	ILE

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	353/398 (89%)	326 (92%)	27 (8%)	16	55
1	B	353/398 (89%)	328 (93%)	25 (7%)	18	57
1	C	353/398 (89%)	329 (93%)	24 (7%)	20	59
1	D	352/398 (88%)	322 (92%)	30 (8%)	13	47
1	E	353/398 (89%)	325 (92%)	28 (8%)	15	53
1	F	353/398 (89%)	329 (93%)	24 (7%)	20	59
1	H	341/398 (86%)	321 (94%)	20 (6%)	24	65
1	I	346/398 (87%)	325 (94%)	21 (6%)	23	64
1	J	330/398 (83%)	318 (96%)	12 (4%)	42	79
All	All	3134/3582 (88%)	2923 (93%)	211 (7%)	20	60

All (211) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	THR
1	A	19	LEU
1	A	37	VAL
1	A	44	ILE
1	A	54	SER
1	A	56	LEU
1	A	65	LEU
1	A	75	VAL
1	A	83	TRP
1	A	110	VAL
1	A	117	VAL
1	A	146	GLU
1	A	167	PRO
1	A	191	LEU
1	A	192	CYS
1	A	217	ARG
1	A	230	SER
1	A	242	TRP
1	A	256	ARG
1	A	289	THR
1	A	373	SER
1	A	384	ASN
1	A	420	PHE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	432	LEU
1	A	434	THR
1	A	445	GLU
1	A	457	SER
1	B	14	THR
1	B	37	VAL
1	B	44	ILE
1	B	73	SER
1	B	83	TRP
1	B	119	SER
1	B	120	LEU
1	B	122	ASP
1	B	146	GLU
1	B	176	GLU
1	B	182	LEU
1	B	210	LYS
1	B	211	ARG
1	B	215	VAL
1	B	217	ARG
1	B	242	TRP
1	B	255	ILE
1	B	256	ARG
1	B	332	LEU
1	B	379	ASP
1	B	404	THR
1	B	428	ASP
1	B	432	LEU
1	B	435	ASP
1	B	453	GLU
1	C	12	GLU
1	C	19	LEU
1	C	44	ILE
1	C	54	SER
1	C	73	SER
1	C	75	VAL
1	C	83	TRP
1	C	90	SER
1	C	182	LEU
1	C	191	LEU
1	C	242	TRP
1	C	256	ARG
1	C	276	MET

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	288	THR
1	C	289	THR
1	C	340	ASP
1	C	343	ASP
1	C	373	SER
1	C	404	THR
1	C	431	GLU
1	C	434	THR
1	C	435	ASP
1	C	447	SER
1	C	453	GLU
1	D	6	THR
1	D	14	THR
1	D	19	LEU
1	D	54	SER
1	D	65	LEU
1	D	73	SER
1	D	75	VAL
1	D	83	TRP
1	D	90	SER
1	D	107	PHE
1	D	111	LEU
1	D	146	GLU
1	D	172	VAL
1	D	211	ARG
1	D	215	VAL
1	D	217	ARG
1	D	242	TRP
1	D	256	ARG
1	D	286	LYS
1	D	289	THR
1	D	332	LEU
1	D	346	SER
1	D	382	TRP
1	D	401	LEU
1	D	404	THR
1	D	420	PHE
1	D	435	ASP
1	D	450	TYR
1	D	453	GLU
1	D	455	ILE
1	E	8	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	E	18	SER
1	E	37	VAL
1	E	44	ILE
1	E	65	LEU
1	E	73	SER
1	E	75	VAL
1	E	83	TRP
1	E	107	PHE
1	E	111	LEU
1	E	116	SER
1	E	117	VAL
1	E	118	SER
1	E	146	GLU
1	E	182	LEU
1	E	210	LYS
1	E	211	ARG
1	E	230	SER
1	E	242	TRP
1	E	289	THR
1	E	332	LEU
1	E	379	ASP
1	E	404	THR
1	E	426	SER
1	E	428	ASP
1	E	430	GLU
1	E	432	LEU
1	E	457	SER
1	F	3	TYR
1	F	10	THR
1	F	14	THR
1	F	19	LEU
1	F	44	ILE
1	F	54	SER
1	F	75	VAL
1	F	83	TRP
1	F	99	ILE
1	F	120	LEU
1	F	141	LEU
1	F	191	LEU
1	F	215	VAL
1	F	217	ARG
1	F	220	MET

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	F	242	TRP
1	F	256	ARG
1	F	289	THR
1	F	340	ASP
1	F	379	ASP
1	F	431	GLU
1	F	432	LEU
1	F	455	ILE
1	F	456	ARG
1	H	3	TYR
1	H	14	THR
1	H	19	LEU
1	H	26	GLN
1	H	37	VAL
1	H	54	SER
1	H	73	SER
1	H	83	TRP
1	H	93	THR
1	H	107	PHE
1	H	110	VAL
1	H	117	VAL
1	H	120	LEU
1	H	172	VAL
1	H	191	LEU
1	H	210	LYS
1	H	211	ARG
1	H	213	ASP
1	H	242	TRP
1	H	382	TRP
1	I	6	THR
1	I	37	VAL
1	I	44	ILE
1	I	65	LEU
1	I	83	TRP
1	I	93	THR
1	I	107	PHE
1	I	111	LEU
1	I	117	VAL
1	I	120	LEU
1	I	122	ASP
1	I	211	ARG
1	I	217	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	I	229	THR
1	I	242	TRP
1	I	276	MET
1	I	340	ASP
1	I	379	ASP
1	I	404	THR
1	I	411	LEU
1	I	450	TYR
1	J	14	THR
1	J	19	LEU
1	J	28	ASP
1	J	65	LEU
1	J	83	TRP
1	J	120	LEU
1	J	242	TRP
1	J	255	ILE
1	J	289	THR
1	J	340	ASP
1	J	373	SER
1	J	404	THR

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

Mol	Chain	Res	Type
1	E	368	ASN
1	E	384	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

9 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PO4	A	1458	-	4,4,4	0.99	0	6,6,6	0.24	0
2	PO4	B	1458	-	4,4,4	0.80	0	6,6,6	0.23	0
2	PO4	C	1458	-	4,4,4	0.86	0	6,6,6	0.23	0
2	PO4	D	1458	-	4,4,4	0.83	0	6,6,6	0.23	0
2	PO4	E	1458	-	4,4,4	0.90	0	6,6,6	0.23	0
2	PO4	F	1458	-	4,4,4	0.95	0	6,6,6	0.24	0
2	PO4	H	1457	-	4,4,4	0.71	0	6,6,6	0.23	0
2	PO4	I	1455	-	4,4,4	0.71	0	6,6,6	0.23	0
2	PO4	J	1455	-	4,4,4	0.72	0	6,6,6	0.23	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	PO4	A	1458	-	-	0/0/0/0	0/0/0/0
2	PO4	B	1458	-	-	0/0/0/0	0/0/0/0
2	PO4	C	1458	-	-	0/0/0/0	0/0/0/0
2	PO4	D	1458	-	-	0/0/0/0	0/0/0/0
2	PO4	E	1458	-	-	0/0/0/0	0/0/0/0
2	PO4	F	1458	-	-	0/0/0/0	0/0/0/0
2	PO4	H	1457	-	-	0/0/0/0	0/0/0/0
2	PO4	I	1455	-	-	0/0/0/0	0/0/0/0
2	PO4	J	1455	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	1458	PO4	1	0
2	I	1455	PO4	2	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	456/505 (90%)	-0.21	6 (1%) 79 67	45, 62, 92, 173	11 (2%)
1	B	456/505 (90%)	-0.02	17 (3%) 45 30	44, 71, 109, 192	18 (3%)
1	C	456/505 (90%)	-0.06	19 (4%) 40 26	49, 80, 122, 155	19 (4%)
1	D	455/505 (90%)	-0.12	13 (2%) 55 41	44, 69, 107, 158	16 (3%)
1	E	456/505 (90%)	0.02	23 (5%) 32 19	46, 76, 115, 211	25 (5%)
1	F	456/505 (90%)	-0.18	11 (2%) 62 47	42, 64, 99, 164	15 (3%)
1	H	440/505 (87%)	0.19	41 (9%) 11 6	57, 83, 138, 160	31 (7%)
1	I	447/505 (88%)	0.09	25 (5%) 28 16	64, 104, 151, 196	55 (12%)
1	J	426/505 (84%)	0.10	23 (5%) 29 17	68, 119, 158, 189	71 (16%)
All	All	4048/4545 (89%)	-0.02	178 (4%) 38 24	42, 77, 139, 211	261 (6%)

All (178) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	382	TRP	6.1
1	C	379	ASP	6.0
1	C	381	GLY	5.8
1	H	216	THR	5.5
1	C	219	GLY	5.5
1	C	378	ILE	5.3
1	E	377	PRO	5.3
1	H	334	SER	5.2
1	H	219	GLY	4.6
1	C	377	PRO	4.6
1	I	241	GLY	4.5
1	E	378	ILE	4.1
1	F	212	ASN	4.1
1	B	426	SER	4.1
1	H	234	GLY	4.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	H	241	GLY	4.0
1	B	3	TYR	4.0
1	E	457	SER	3.8
1	H	39	SER	3.8
1	E	382	TRP	3.7
1	I	3	TYR	3.6
1	I	17	ASN	3.6
1	I	45	MET	3.6
1	F	335	LEU	3.5
1	H	218	LYS	3.5
1	H	332	LEU	3.5
1	F	338	ILE	3.4
1	D	3	TYR	3.4
1	H	214	PRO	3.4
1	H	41	GLY	3.3
1	J	386	HIS	3.3
1	E	456	ARG	3.3
1	H	288	THR	3.3
1	J	385	HIS	3.3
1	D	219	GLY	3.2
1	H	340	ASP	3.2
1	F	336	LEU	3.2
1	J	381	GLY	3.2
1	J	452	PRO	3.2
1	H	48	GLY	3.2
1	E	241	GLY	3.2
1	B	430	GLU	3.2
1	B	241	GLY	3.2
1	J	361	ILE	3.2
1	H	45	MET	3.1
1	J	390	VAL	3.1
1	C	380	GLY	3.1
1	J	365	ASP	3.1
1	C	387	TYR	3.1
1	E	376	SER	3.1
1	B	425	LEU	3.1
1	D	455	ILE	3.0
1	H	42	VAL	3.0
1	H	94	ARG	3.0
1	E	334	SER	3.0
1	C	389	GLN	3.0
1	D	428	ASP	2.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	385	HIS	2.9
1	C	386	HIS	2.9
1	H	221	PRO	2.9
1	E	426	SER	2.9
1	H	223	TYR	2.9
1	J	331	ASP	2.9
1	H	208	LEU	2.9
1	J	384	ASN	2.8
1	C	241	GLY	2.8
1	H	98	PHE	2.8
1	B	387	TYR	2.8
1	B	427	ALA	2.8
1	J	360	GLY	2.8
1	H	38	ALA	2.8
1	E	103	GLU	2.8
1	F	39	SER	2.8
1	E	430	GLU	2.7
1	I	219	GLY	2.7
1	B	456	ARG	2.7
1	I	21	THR	2.7
1	I	18	SER	2.7
1	D	457	SER	2.7
1	H	287	TRP	2.7
1	H	337	ARG	2.7
1	I	234	GLY	2.7
1	J	380	GLY	2.7
1	J	387	TYR	2.6
1	D	241	GLY	2.6
1	E	3	TYR	2.6
1	H	335	LEU	2.6
1	I	386	HIS	2.6
1	D	427	ALA	2.6
1	J	372	GLY	2.6
1	E	425	LEU	2.6
1	A	3	TYR	2.5
1	E	92	ASN	2.5
1	A	217	ARG	2.5
1	B	2	SER	2.5
1	E	433	GLY	2.5
1	H	97	GLY	2.5
1	J	362	PHE	2.5
1	B	234	GLY	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	I	220	MET	2.5
1	J	377	PRO	2.5
1	D	429	GLU	2.5
1	E	429	GLU	2.5
1	I	39	SER	2.5
1	C	384	ASN	2.5
1	E	375	ILE	2.5
1	D	48	GLY	2.5
1	E	424	ARG	2.5
1	H	8	THR	2.5
1	J	241	GLY	2.5
1	F	214	PRO	2.5
1	H	2	SER	2.4
1	F	433	GLY	2.4
1	C	427	ALA	2.4
1	D	418	ILE	2.4
1	I	242	TRP	2.4
1	H	131	MET	2.4
1	H	130	GLY	2.4
1	J	396	GLY	2.4
1	J	392	TYR	2.4
1	I	8	THR	2.4
1	I	23	LEU	2.4
1	C	428	ASP	2.4
1	H	96	ASN	2.4
1	H	3	TYR	2.4
1	J	382	TRP	2.3
1	H	40	ALA	2.3
1	C	425	LEU	2.3
1	I	19	LEU	2.3
1	H	331	ASP	2.3
1	H	222	LYS	2.3
1	B	429	GLU	2.3
1	I	221	PRO	2.3
1	B	48	GLY	2.3
1	F	213	ASP	2.3
1	A	38	ALA	2.3
1	A	283	CYS	2.3
1	I	387	TYR	2.3
1	B	20	THR	2.3
1	J	363	ALA	2.3
1	J	339	ASP	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	I	5	PHE	2.3
1	J	41	GLY	2.2
1	H	220	MET	2.2
1	B	428	ASP	2.2
1	I	216	THR	2.2
1	F	2	SER	2.2
1	H	95	GLY	2.2
1	A	335	LEU	2.2
1	A	336	LEU	2.2
1	H	336	LEU	2.2
1	H	99	ILE	2.2
1	J	379	ASP	2.2
1	D	9	PRO	2.1
1	D	234	GLY	2.1
1	D	419	PRO	2.1
1	B	416	ASN	2.1
1	F	211	ARG	2.1
1	F	220	MET	2.1
1	I	238	LEU	2.1
1	E	93	THR	2.1
1	H	209	GLY	2.1
1	B	39	SER	2.1
1	E	207	ILE	2.1
1	H	230	SER	2.1
1	E	340	ASP	2.1
1	I	7	GLY	2.1
1	C	390	VAL	2.1
1	E	339	ASP	2.1
1	I	420	PHE	2.0
1	B	417	ALA	2.0
1	I	436	ALA	2.0
1	E	361	ILE	2.0
1	C	220	MET	2.0
1	I	390	VAL	2.0
1	H	44	ILE	2.0
1	C	364	ALA	2.0
1	I	419	PRO	2.0

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

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

6.3 Carbohydrates [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. 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	PO4	D	1458	5/5	0.96	0.32	4.10	142,143,144,147	0
2	PO4	E	1458	5/5	0.96	0.25	3.31	113,113,114,114	0
2	PO4	F	1458	5/5	0.95	0.21	1.94	97,97,100,104	0
2	PO4	J	1455	5/5	0.91	0.25	1.78	163,163,163,163	0
2	PO4	B	1458	5/5	0.95	0.21	1.40	99,100,101,102	0
2	PO4	A	1458	5/5	0.94	0.19	1.02	98,101,104,105	0
2	PO4	H	1457	5/5	0.81	0.23	0.73	200,202,202,204	0
2	PO4	C	1458	5/5	0.97	0.18	0.71	110,110,111,112	0
2	PO4	I	1455	5/5	0.67	0.20	0.53	178,181,181,182	0

6.5 Other polymers [i](#)

There are no such residues in this entry.