



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:46 AM GMT

PDB ID : 3PUK  
Title : Re-refinement of the crystal structure of Munc18-3 and Syntaxin4 N-peptide complex  
Authors : Hu, S.-H.; Christie, M.P.; Saez, N.J.; Latham, C.F.; Jarrott, R.; Lua, L.H.L.; Collins, B.M.; Martin, J.L.  
Deposited on : 2010-12-05  
Resolution : 3.05 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](mailto: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



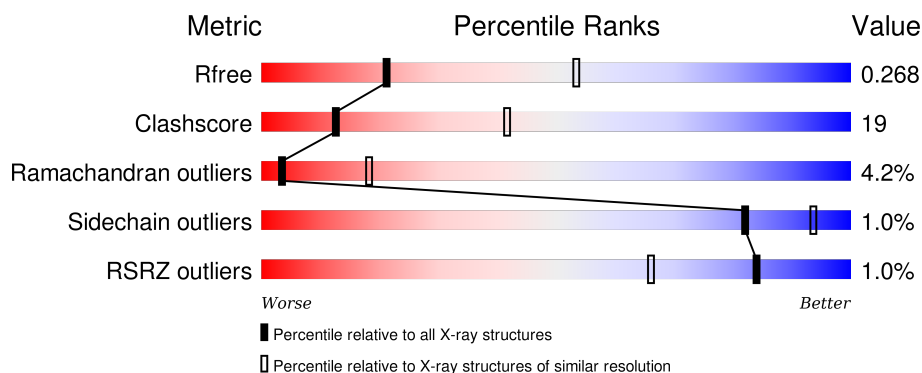
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.05 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	1191 (3.10-3.02)
Clashscore	102246	1303 (3.10-3.02)
Ramachandran outliers	100387	1254 (3.10-3.02)
Sidechain outliers	100360	1254 (3.10-3.02)
RSRZ outliers	91569	1197 (3.10-3.02)

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  $\geq 3$ , 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	592	<div> <div>10%</div> <div>59%</div> <div>31%</div> <div>7%</div> </div>
1	B	592	<div> <div>64%</div> <div>25%</div> <div>8%</div> </div>
2	C	10	<div> <div>10%</div> <div>80%</div> <div>10%</div> <div>10%</div> </div>
2	D	10	<div> <div>80%</div> <div>20%</div> </div>



## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8521 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 Syntaxin-binding protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	552	Total	C	N	O	S	200	0	0
			4213	2664	719	805	25			
1	B	545	Total	C	N	O	S	158	0	0
			4143	2622	705	788	28			

- Molecule 2 is a protein called Syntaxin-4 N-terminal peptide.

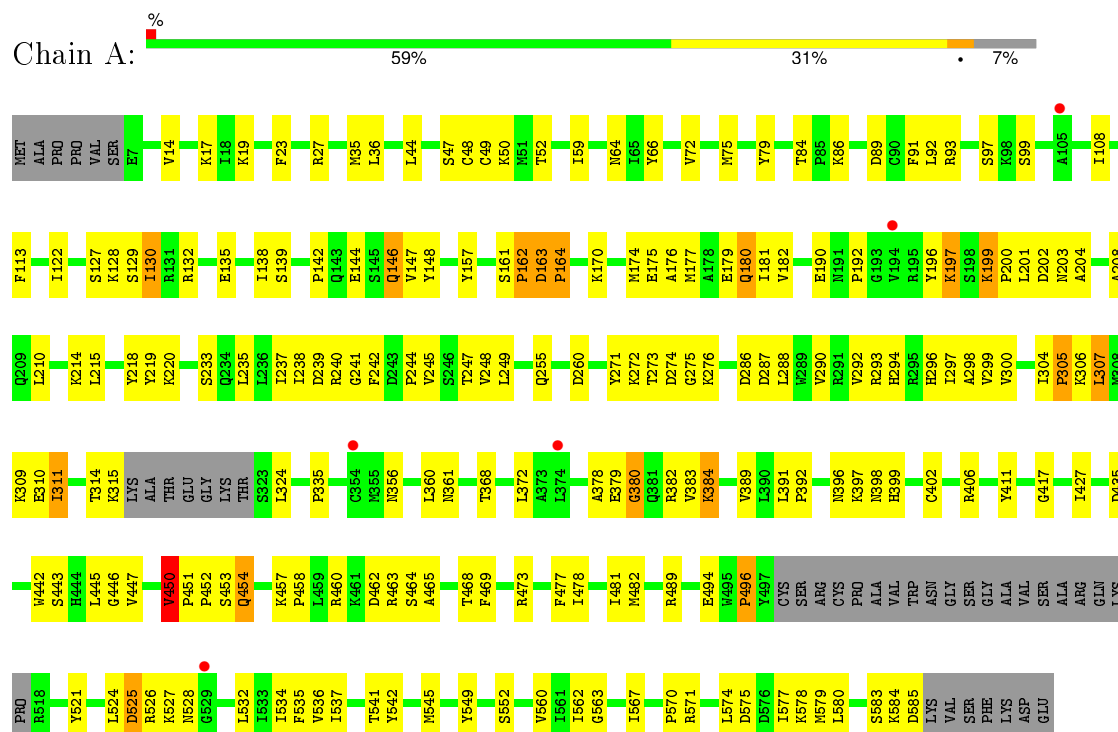
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	9	Total	C	N	O	0	0	0
			80	46	20	14			
2	D	10	Total	C	N	O	0	0	0
			85	49	21	15			



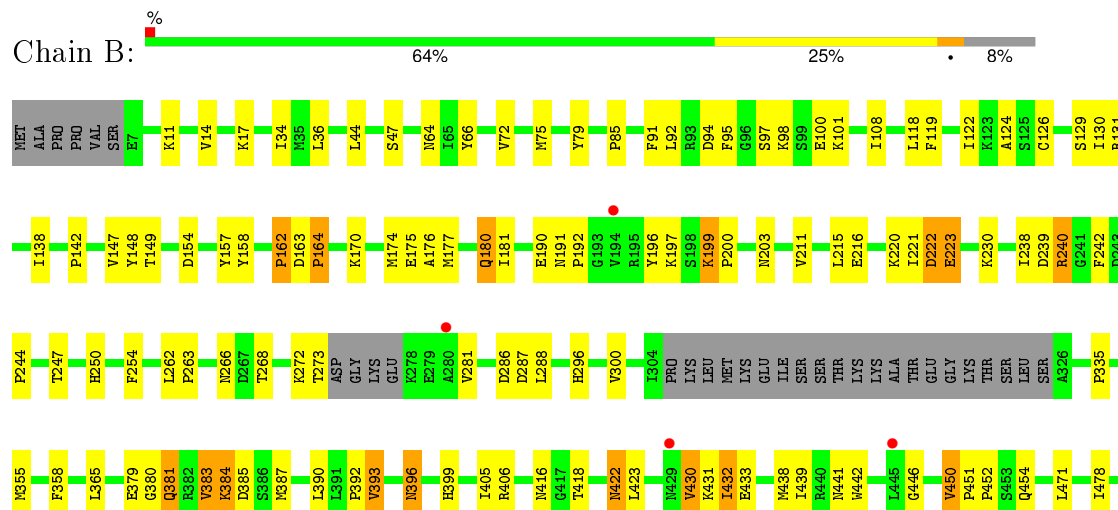
### 3 Residue-property plots

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.

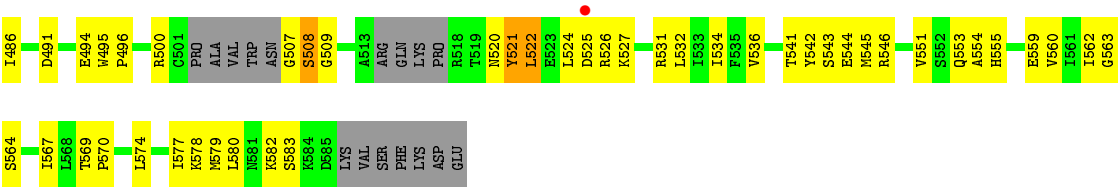
#### • Molecule 1: Syntaxin-binding protein 3



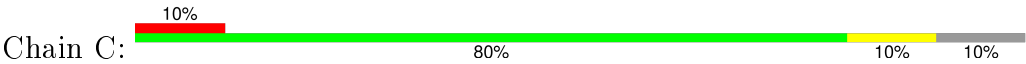
#### • Molecule 1: Syntaxin-binding protein 3



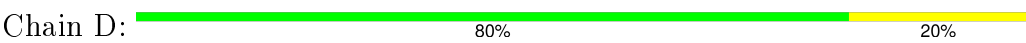




● Molecule 2: Syntaxis-4 N-terminal peptide



● Molecule 2: Syntaxis-4 N-terminal peptide





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	170.40 Å   170.40 Å   170.40 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	45.54 – 3.05 45.54 – 3.05	Depositor EDS
% Data completeness (in resolution range)	98.7 (45.54-3.05) 98.7 (45.54-3.05)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.01 (at 3.06 Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, $R_{free}$	0.239 , 0.279 0.226 , 0.268	Depositor DCC
$R_{free}$ test set	1572 reflections (5.05%)	DCC
Wilson B-factor (Å <sup>2</sup> )	99.8	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 148.4	EDS
Estimated twinning fraction	0.478 for l,-k,h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.53$ , $\langle L^2 \rangle = 0.37$	Xtriage
Outliers	3 of 31163 reflections (0.010%)	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	8521	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	148.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

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	0.22	0/4285	0.40	2/5807 (0.0%)
1	B	0.22	0/4214	0.39	1/5709 (0.0%)
2	C	0.19	0/80	0.37	0/105
2	D	0.19	0/85	0.38	0/112
All	All	0.22	0/8664	0.40	3/11733 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	335	PRO	N-CA-CB	6.05	110.56	103.30
1	A	305	PRO	N-CA-CB	6.04	110.55	103.30
1	A	335	PRO	N-CA-CB	5.90	110.38	103.30

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	4213	0	4003	169	0
1	B	4143	0	3944	132	0
2	C	80	0	78	1	0
2	D	85	0	80	3	0
All	All	8521	0	8105	304	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 19.

All (304) 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:199:LYS:HB3	1:A:200:PRO:HD3	1.38	1.03
1:B:450:VAL:HG13	1:B:451:PRO:HD3	1.40	1.02
1:A:163:ASP:H	1:A:164:PRO:HD3	1.30	0.96
2:D:9:ARG:HG3	2:D:10:GLN:H	1.29	0.95
1:A:526:ARG:HB3	1:A:527:LYS:HA	1.47	0.95
1:A:450:VAL:HG23	1:A:451:PRO:HD3	1.51	0.92
1:B:199:LYS:HB3	1:B:200:PRO:HD3	1.53	0.91
1:B:199:LYS:HB3	1:B:200:PRO:CD	2.01	0.89
1:A:127:SER:HB3	1:A:129:SER:N	1.90	0.87
1:B:526:ARG:H	1:B:527:LYS:HA	1.39	0.86
1:B:365:LEU:HD13	1:B:393:VAL:HG11	1.56	0.86
1:B:441:ASN:HD22	1:B:578:LYS:HA	1.42	0.84
1:A:271:TYR:HB3	1:A:272:LYS:HA	1.58	0.84
1:A:197:LYS:HG2	1:A:199:LYS:H	1.41	0.83
1:A:163:ASP:N	1:A:164:PRO:HD3	1.94	0.81
1:B:383:VAL:HG13	1:B:384:LYS:H	1.45	0.81
1:A:545:MET:HA	1:A:562:ILE:HD11	1.65	0.78
1:B:579:MET:HA	1:B:582:LYS:HE2	1.66	0.78
1:A:199:LYS:HB3	1:A:200:PRO:CD	2.13	0.77
1:B:100:GLU:CB	1:B:101:LYS:HA	2.14	0.77
1:A:463:ARG:HA	1:A:464:SER:OG	1.86	0.76
1:B:197:LYS:HB3	1:B:238:ILE:HG22	1.66	0.76
1:A:72:VAL:HG11	1:A:75:MET:HE3	1.68	0.75
1:A:524:LEU:HB2	1:A:525:ASP:CB	2.18	0.74
1:A:199:LYS:CB	1:A:200:PRO:HD3	2.15	0.73
1:A:526:ARG:HB3	1:A:527:LYS:CA	2.18	0.73
1:B:199:LYS:CB	1:B:200:PRO:CD	2.67	0.72
1:B:422:ASN:HD22	1:B:423:LEU:N	1.87	0.71
1:B:197:LYS:HE3	1:B:199:LYS:HB2	1.70	0.71
1:A:127:SER:HB2	1:A:130:ILE:HG13	1.72	0.71
1:A:526:ARG:CB	1:A:527:LYS:HA	2.13	0.70
1:B:162:PRO:HD3	1:B:580:LEU:O	1.92	0.69
1:A:245:VAL:HG12	1:A:249:LEU:HD21	1.73	0.69
1:A:199:LYS:CB	1:A:200:PRO:CD	2.69	0.69
1:A:14:VAL:HG21	1:A:142:PRO:HG3	1.75	0.69
1:B:222:ASP:O	1:B:223:GLU:HB2	1.91	0.69
1:A:272:LYS:CB	1:A:273:THR:HA	2.23	0.69

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:GLU:CB	1:B:101:LYS:HD3	2.23	0.68
1:B:430:VAL:HG13	1:B:431:LYS:H	1.58	0.68
1:A:176:ALA:O	1:A:180:GLN:HB2	1.92	0.68
1:B:240:ARG:O	1:B:240:ARG:HD3	1.94	0.67
1:B:268:THR:HG22	1:B:281:VAL:HG22	1.74	0.67
1:B:162:PRO:C	1:B:164:PRO:HD3	2.15	0.66
1:A:397:LYS:HA	1:A:398:ASN:C	2.15	0.66
1:B:240:ARG:HH12	1:B:242:PHE:HE1	1.45	0.65
1:B:545:MET:HE3	1:B:567:ILE:HG23	1.78	0.65
1:B:176:ALA:O	1:B:180:GLN:HB2	1.96	0.65
1:B:163:ASP:N	1:B:164:PRO:HD3	2.11	0.65
1:A:210:LEU:O	1:A:214:LYS:HB2	1.96	0.65
1:B:526:ARG:N	1:B:527:LYS:HA	2.03	0.64
1:B:380:GLY:HA2	1:B:381:GLN:CB	2.27	0.64
1:A:271:TYR:HB3	1:A:272:LYS:CA	2.26	0.64
1:A:442:TRP:CE2	1:A:577:ILE:HD13	2.34	0.63
1:B:524:LEU:O	1:B:525:ASP:HB2	1.99	0.62
1:A:113:PHE:HD2	1:A:135:GLU:OE2	1.82	0.62
1:A:545:MET:HE3	1:A:567:ILE:HG23	1.82	0.62
1:A:127:SER:HB3	1:A:130:ILE:H	1.63	0.62
1:B:247:THR:HA	1:B:541:THR:HG21	1.81	0.62
1:A:197:LYS:HD2	1:A:239:ASP:H	1.65	0.62
1:B:532:LEU:HD23	1:B:560:VAL:HG22	1.81	0.62
1:A:163:ASP:H	1:A:164:PRO:CD	2.09	0.61
1:B:72:VAL:HG11	1:B:75:MET:HE3	1.81	0.61
1:B:17:LYS:HD3	1:B:138:ILE:HD11	1.81	0.61
1:A:532:LEU:HD23	1:A:560:VAL:HG22	1.82	0.60
1:A:383:VAL:O	1:A:384:LYS:HB2	2.01	0.60
1:B:442:TRP:CE2	1:B:577:ILE:HD13	2.37	0.60
1:B:432:ILE:HA	1:B:433:GLU:CB	2.29	0.59
1:A:298:ALA:HB3	1:A:469:PHE:CE1	2.37	0.59
2:D:9:ARG:HG3	2:D:10:GLN:N	2.08	0.59
1:A:163:ASP:N	1:A:164:PRO:CD	2.64	0.59
1:A:417:GLY:HA3	1:A:452:PRO:HA	1.84	0.59
1:B:430:VAL:HG13	1:B:431:LYS:N	2.18	0.59
1:B:14:VAL:HG21	1:B:142:PRO:HG3	1.84	0.58
1:A:450:VAL:CG2	1:A:451:PRO:HD3	2.31	0.58
1:B:296:HIS:O	1:B:300:VAL:HG23	2.03	0.58
1:A:92:LEU:HD11	1:A:122:ILE:HG12	1.85	0.58
1:B:197:LYS:O	1:B:199:LYS:HG3	2.03	0.58
1:B:422:ASN:C	1:B:422:ASN:HD22	2.06	0.58

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:430:VAL:HG22	1:B:431:LYS:N	2.18	0.58
1:A:293:ARG:HG3	1:A:294:HIS:CD2	2.39	0.57
1:B:543:SER:O	1:B:546:ARG:HG2	2.03	0.57
1:B:441:ASN:ND2	1:B:578:LYS:HA	2.17	0.57
1:A:162:PRO:HB2	1:A:164:PRO:HD3	1.87	0.57
1:A:91:PHE:CZ	1:A:108:ILE:HD11	2.40	0.57
1:A:157:TYR:OH	1:A:583:SER:HA	2.03	0.57
1:A:397:LYS:CB	1:A:399:HIS:HB2	2.34	0.57
1:A:460:ARG:HD2	1:A:477:PHE:CD1	2.40	0.57
1:B:574:LEU:O	1:B:577:ILE:HG22	2.04	0.56
1:B:157:TYR:OH	1:B:583:SER:HA	2.05	0.56
1:A:460:ARG:HH11	1:A:477:PHE:HA	1.70	0.56
1:A:391:LEU:HB2	1:A:392:PRO:HD3	1.88	0.56
1:A:534:ILE:HB	1:A:562:ILE:HG22	1.87	0.55
1:B:34:ILE:HD11	1:B:75:MET:HE3	1.87	0.55
1:A:574:LEU:O	1:A:577:ILE:HG22	2.06	0.55
1:B:200:PRO:HB3	1:B:446:GLY:HA3	1.87	0.55
1:A:397:LYS:HA	1:A:399:HIS:N	2.21	0.55
1:A:247:THR:HA	1:A:541:THR:HG21	1.88	0.55
1:A:19:LYS:HA	1:A:23:PHE:HD1	1.72	0.55
1:A:453:SER:O	1:A:454:GLN:CB	2.55	0.55
1:A:356:ASN:O	1:A:360:LEU:HG	2.06	0.55
1:A:240:ARG:NH2	1:A:247:THR:HG22	2.22	0.55
1:A:240:ARG:NH2	1:A:541:THR:HG23	2.22	0.54
1:B:551:VAL:O	1:B:555:HIS:HB3	2.06	0.54
1:A:64:ASN:OD1	1:A:66:TYR:HB2	2.07	0.54
1:A:245:VAL:HG12	1:A:249:LEU:CD2	2.37	0.54
1:A:382:ARG:O	1:A:383:VAL:HB	2.08	0.54
1:B:405:ILE:HG23	1:B:439:ILE:HD11	1.88	0.54
1:B:520:ASN:O	1:B:521:TYR:CB	2.55	0.54
1:B:578:LYS:C	1:B:580:LEU:H	2.10	0.54
1:B:197:LYS:CE	1:B:199:LYS:HB2	2.36	0.53
1:B:430:VAL:HG22	1:B:431:LYS:H	1.73	0.53
1:B:64:ASN:OD1	1:B:66:TYR:HB2	2.07	0.53
1:B:36:LEU:HD22	1:B:79:TYR:CE2	2.43	0.53
1:A:578:LYS:C	1:A:580:LEU:H	2.11	0.53
1:A:463:ARG:HA	1:A:464:SER:CB	2.38	0.53
1:B:534:ILE:HB	1:B:562:ILE:HG22	1.90	0.53
1:B:418:THR:HB	1:B:422:ASN:HD21	1.74	0.53
1:B:272:LYS:O	1:B:273:THR:C	2.47	0.53
1:A:299:VAL:HG23	1:A:469:PHE:HZ	1.74	0.53

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:ASN:HB2	1:B:239:ASP:OD1	2.09	0.52
1:A:175:GLU:O	1:A:179:GLU:HG2	2.09	0.52
1:B:396:ASN:N	1:B:396:ASN:HD22	2.06	0.52
1:A:372:LEU:HD11	1:A:411:TYR:HA	1.90	0.52
1:A:218:TYR:O	1:A:219:TYR:HB2	2.09	0.52
1:A:139:SER:HB3	1:A:180:GLN:HG3	1.92	0.52
1:A:378:ALA:H	1:A:473:ARG:NH1	2.07	0.52
1:B:450:VAL:CG1	1:B:451:PRO:HD3	2.26	0.52
1:B:242:PHE:CZ	1:B:478:ILE:HG23	2.45	0.52
1:B:578:LYS:O	1:B:579:MET:HB2	2.08	0.52
1:B:542:TYR:CZ	1:B:567:ILE:HG21	2.45	0.52
2:D:9:ARG:CG	2:D:10:GLN:H	2.11	0.51
1:A:402:CYS:SG	1:A:571:ARG:HG2	2.51	0.51
1:B:230:LYS:HE3	1:B:500:ARG:CB	2.40	0.51
1:A:177:MET:O	1:A:181:ILE:HG13	2.10	0.51
1:B:190:GLU:HG3	1:B:531:ARG:HD2	1.92	0.51
1:A:146:GLN:HE22	1:A:552:SER:HB2	1.76	0.51
1:A:240:ARG:HD2	1:A:240:ARG:O	2.11	0.51
1:A:238:ILE:HD11	1:A:536:VAL:HG22	1.92	0.51
1:A:196:TYR:HD2	1:A:197:LYS:O	1.95	0.50
1:A:199:LYS:HE3	1:A:494:GLU:CG	2.42	0.50
1:A:298:ALA:HB3	1:A:469:PHE:HE1	1.75	0.50
1:A:196:TYR:CZ	1:A:496:PRO:HG2	2.47	0.50
1:B:521:TYR:O	1:B:522:LEU:CB	2.59	0.50
1:A:35:MET:HE2	1:A:59:ILE:HG21	1.93	0.50
1:A:127:SER:CB	1:A:130:ILE:H	2.25	0.50
1:A:240:ARG:HB2	1:A:537:ILE:O	2.12	0.50
1:A:36:LEU:HD22	1:A:79:TYR:CE2	2.47	0.50
1:B:197:LYS:NZ	1:B:200:PRO:HD2	2.26	0.50
1:A:368:THR:HG21	1:A:392:PRO:HG2	1.94	0.49
1:B:216:GLU:O	1:B:220:LYS:HG3	2.12	0.49
1:B:92:LEU:HD21	1:B:122:ILE:HG22	1.93	0.49
1:B:124:ALA:C	1:B:126:CYS:H	2.16	0.49
1:B:148:TYR:CE2	1:B:563:GLY:HA3	2.47	0.49
1:A:309:LYS:HA	1:A:311:ILE:N	2.27	0.49
1:B:438:MET:HG3	1:B:578:LYS:HE3	1.95	0.49
1:B:242:PHE:CE2	1:B:478:ILE:HG23	2.48	0.49
1:A:309:LYS:CB	1:A:310:GLU:HA	2.43	0.49
1:A:17:LYS:HD3	1:A:138:ILE:HD11	1.94	0.49
1:B:158:TYR:O	1:B:170:LYS:HG3	2.12	0.49
1:A:290:VAL:HA	1:A:293:ARG:HE	1.79	0.48

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:491:ASP:HB3	1:B:494:GLU:HB3	1.95	0.48
1:A:196:TYR:CE2	1:A:496:PRO:HG2	2.49	0.48
1:A:197:LYS:NZ	1:A:200:PRO:HD2	2.28	0.48
1:B:471:LEU:HD12	1:B:471:LEU:N	2.27	0.48
1:A:306:LYS:HA	1:A:307:LEU:CB	2.43	0.48
1:A:179:GLU:HB3	1:A:214:LYS:HE2	1.94	0.48
1:A:161:SER:OG	1:A:162:PRO:HD2	2.13	0.48
1:B:286:ASP:C	1:B:288:LEU:H	2.16	0.48
1:A:200:PRO:HG2	1:A:241:GLY:HA3	1.95	0.48
1:B:211:VAL:O	1:B:215:LEU:HG	2.14	0.48
1:A:89:ASP:O	1:A:93:ARG:HG3	2.14	0.48
1:A:97:SER:C	1:A:99:SER:H	2.17	0.48
1:B:119:PHE:O	1:B:122:ILE:HG12	2.14	0.47
1:A:48:CYS:O	1:A:49:CYS:HB3	2.14	0.47
1:A:199:LYS:HE3	1:A:494:GLU:HG2	1.96	0.47
1:A:235:LEU:HG	1:A:237:ILE:CD1	2.44	0.47
1:A:200:PRO:HB3	1:A:446:GLY:HA3	1.96	0.47
1:B:147:VAL:HG22	1:B:562:ILE:HD11	1.97	0.47
1:B:154:ASP:O	1:B:158:TYR:HD1	1.97	0.47
1:A:450:VAL:H	1:A:451:PRO:CD	2.28	0.47
1:B:416:ASN:O	1:B:452:PRO:HA	2.15	0.47
1:A:242:PHE:CZ	1:A:478:ILE:HG23	2.50	0.47
1:B:215:LEU:HD11	1:B:230:LYS:NZ	2.29	0.47
1:B:525:ASP:HA	1:B:526:ARG:HA	1.72	0.47
1:B:432:ILE:CA	1:B:433:GLU:CB	2.93	0.47
1:A:240:ARG:HH21	1:A:247:THR:HG22	1.80	0.46
1:A:255:GLN:HB2	1:A:293:ARG:HB2	1.97	0.46
1:A:296:HIS:O	1:A:300:VAL:HG23	2.15	0.46
1:A:286:ASP:C	1:A:288:LEU:H	2.19	0.46
1:B:238:ILE:HD11	1:B:536:VAL:HG22	1.97	0.46
1:B:221:ILE:O	1:B:223:GLU:N	2.49	0.46
1:A:477:PHE:O	1:A:481:ILE:HG13	2.16	0.46
1:B:221:ILE:O	1:B:222:ASP:C	2.54	0.46
1:A:175:GLU:HG3	1:A:176:ALA:N	2.30	0.46
1:B:175:GLU:HG3	1:B:176:ALA:N	2.31	0.46
1:A:292:VAL:HG13	1:A:300:VAL:HG22	1.97	0.46
1:A:584:LYS:HA	1:A:585:ASP:CB	2.44	0.46
1:A:460:ARG:NH1	1:A:477:PHE:HA	2.31	0.46
1:B:250:HIS:CE1	1:B:406:ARG:CZ	2.99	0.46
1:A:524:LEU:O	1:A:524:LEU:HD12	2.16	0.45
1:B:418:THR:HB	1:B:422:ASN:ND2	2.30	0.45

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:390:LEU:C	1:B:390:LEU:HD23	2.37	0.45
1:A:148:TYR:CZ	1:A:563:GLY:HA3	2.51	0.45
1:A:378:ALA:O	1:A:379:GLU:HB3	2.16	0.45
1:A:204:ALA:HB1	1:A:237:ILE:HG22	1.97	0.45
1:B:177:MET:O	1:B:181:ILE:HG13	2.16	0.45
1:B:525:ASP:HB3	1:B:526:ARG:CB	2.47	0.45
1:A:468:THR:HG22	1:A:469:PHE:N	2.32	0.45
1:B:191:ASN:OD1	1:B:215:LEU:HD13	2.17	0.45
1:B:85:PRO:HA	1:B:118:LEU:HD21	1.98	0.45
1:B:553:GLN:O	1:B:554:ALA:HB3	2.16	0.45
1:A:296:HIS:CE1	1:A:297:ILE:HG22	2.51	0.45
1:B:196:TYR:CZ	1:B:496:PRO:HB2	2.52	0.45
1:B:450:VAL:O	1:B:452:PRO:HD3	2.17	0.45
1:B:383:VAL:HG13	1:B:384:LYS:N	2.22	0.45
1:A:378:ALA:C	1:A:380:GLY:H	2.21	0.45
1:B:197:LYS:HE2	1:B:239:ASP:CB	2.47	0.44
1:A:91:PHE:CD2	1:A:92:LEU:HD12	2.52	0.44
1:B:478:ILE:HD11	1:B:544:GLU:HB3	1.98	0.44
1:A:147:VAL:HG11	1:A:545:MET:CE	2.47	0.44
1:A:435:ASP:HA	1:A:578:LYS:NZ	2.31	0.44
1:B:355:MET:O	1:B:358:PHE:HB3	2.18	0.44
1:B:197:LYS:NZ	1:B:242:PHE:HB3	2.32	0.44
1:A:478:ILE:O	1:A:482:MET:HG3	2.17	0.44
1:A:273:THR:HB	1:A:274:ASP:HB2	2.00	0.44
1:B:190:GLU:C	1:B:192:PRO:HD3	2.38	0.44
1:A:235:LEU:HG	1:A:237:ILE:HD11	1.99	0.44
1:A:545:MET:HA	1:A:562:ILE:CD1	2.44	0.44
1:B:197:LYS:HD2	1:B:199:LYS:HD2	1.99	0.43
1:B:431:LYS:O	1:B:432:ILE:C	2.56	0.43
1:A:575:ASP:O	1:A:578:LYS:HG3	2.18	0.43
1:A:389:VAL:O	1:A:392:PRO:HD2	2.18	0.43
1:A:23:PHE:O	1:A:27:ARG:HB2	2.19	0.43
1:B:191:ASN:O	1:B:230:LYS:HA	2.19	0.43
1:A:271:TYR:HA	1:A:272:LYS:O	2.18	0.43
1:A:91:PHE:HD2	1:A:92:LEU:HD12	1.82	0.43
1:A:182:VAL:HG21	1:A:214:LYS:HB3	1.99	0.43
1:B:545:MET:HG2	1:B:562:ILE:HD11	2.00	0.43
1:B:559:GLU:O	1:B:559:GLU:HG3	2.19	0.43
1:A:417:GLY:CA	1:A:452:PRO:HA	2.49	0.43
1:A:457:LYS:HA	1:A:458:PRO:HD3	1.89	0.43
1:A:240:ARG:HH11	1:A:242:PHE:HE2	1.66	0.42

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:TYR:C	1:A:220:LYS:H	2.22	0.42
1:A:383:VAL:O	1:A:384:LYS:CB	2.67	0.42
1:A:144:GLU:HB3	1:A:549:TYR:OH	2.18	0.42
1:A:215:LEU:HA	1:A:218:TYR:HB3	2.00	0.42
1:A:181:ILE:HD11	1:A:535:PHE:CZ	2.55	0.42
1:A:237:ILE:N	1:A:237:ILE:HD12	2.34	0.42
1:B:385:ASP:C	1:B:387:MET:H	2.22	0.42
1:A:170:LYS:O	1:A:174:MET:HG2	2.19	0.42
1:B:91:PHE:CZ	1:B:108:ILE:HD11	2.55	0.42
1:A:260:ASP:CG	1:A:570:PRO:HG2	2.39	0.42
1:A:482:MET:HG2	1:A:534:ILE:HD11	2.01	0.42
1:B:97:SER:HA	1:B:98:LYS:HA	1.63	0.42
1:A:378:ALA:HA	1:A:473:ARG:HH12	1.85	0.42
1:A:44:LEU:O	1:A:47:SER:HB3	2.18	0.42
1:A:578:LYS:O	1:A:579:MET:HB2	2.19	0.42
1:B:197:LYS:HE2	1:B:239:ASP:H	1.84	0.42
1:A:201:LEU:N	1:A:201:LEU:HD12	2.34	0.42
1:A:244:PRO:HB3	1:A:442:TRP:CZ3	2.55	0.42
1:A:275:GLY:HA2	1:A:276:LYS:HA	1.66	0.42
1:A:306:LYS:HA	1:A:307:LEU:HA	1.83	0.42
1:A:127:SER:HB3	1:A:129:SER:H	1.79	0.41
1:A:190:GLU:C	1:A:192:PRO:HD3	2.40	0.41
1:B:266:ASN:O	1:B:268:THR:HG23	2.19	0.41
1:B:262:LEU:HA	1:B:263:PRO:HD3	1.84	0.41
1:B:11:LYS:HB3	1:B:142:PRO:HB2	2.02	0.41
1:B:495:TRP:N	1:B:496:PRO:HD3	2.36	0.41
1:A:84:THR:HG22	1:A:86:LYS:H	1.85	0.41
1:B:507:GLY:O	1:B:508:SER:C	2.59	0.41
1:A:192:PRO:HA	1:A:233:SER:O	2.21	0.41
1:A:314:THR:HA	1:A:315:LYS:HA	1.56	0.41
1:A:443:SER:HA	1:A:447:VAL:O	2.21	0.41
1:A:132:ARG:HA	2:C:5:THR:HG23	2.03	0.41
1:A:201:LEU:HD11	1:A:445:LEU:C	2.41	0.41
1:B:244:PRO:HB3	1:B:442:TRP:CZ3	2.55	0.41
1:B:44:LEU:O	1:B:47:SER:HB3	2.21	0.41
1:B:129:SER:O	1:B:131:ARG:HG2	2.21	0.41
1:A:273:THR:HA	1:A:274:ASP:HA	1.86	0.41
1:B:94:ASP:HB2	1:B:95:PHE:CD1	2.56	0.41
1:A:196:TYR:HB3	1:A:208:ALA:HB2	2.03	0.41
1:B:199:LYS:CB	1:B:200:PRO:HD2	2.46	0.41
1:A:542:TYR:CE1	1:A:567:ILE:HG21	2.56	0.41

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:GLU:CB	1:B:101:LYS:CA	2.92	0.41
1:B:486:ILE:HG21	1:B:555:HIS:NE2	2.36	0.41
1:B:396:ASN:HB3	1:B:399:HIS:HD1	1.86	0.41
1:B:170:LYS:O	1:B:174:MET:HG2	2.21	0.41
1:A:427:ILE:HG22	1:A:427:ILE:O	2.21	0.41
1:B:569:THR:O	1:B:570:PRO:C	2.59	0.41
1:A:584:LYS:HA	1:A:585:ASP:C	2.41	0.41
1:A:450:VAL:O	1:A:452:PRO:HD3	2.22	0.40
1:A:147:VAL:HG11	1:A:545:MET:HE2	2.03	0.40
1:A:396:ASN:O	1:A:398:ASN:HB3	2.21	0.40
1:B:542:TYR:CE1	1:B:567:ILE:HG21	2.56	0.40
1:A:445:LEU:HD22	1:A:577:ILE:HD11	2.02	0.40
1:B:240:ARG:HA	1:B:240:ARG:HH11	1.87	0.40
1:A:396:ASN:O	1:A:399:HIS:ND1	2.54	0.40
1:A:50:LYS:C	1:A:52:THR:H	2.25	0.40
1:B:92:LEU:HD21	1:B:122:ILE:HA	2.03	0.40
1:A:248:VAL:HG13	1:A:406:ARG:HB3	2.03	0.40
1:A:201:LEU:HD22	1:A:203:ASN:ND2	2.36	0.40
1:A:391:LEU:O	1:A:392:PRO:C	2.60	0.40
1:A:146:GLN:HE22	1:A:552:SER:CB	2.34	0.40
1:B:149:THR:HA	1:B:564:SER:O	2.21	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	546/592 (92%)	426 (78%)	94 (17%)	26 (5%)	3	16
1	B	535/592 (90%)	421 (79%)	94 (18%)	20 (4%)	4	22
2	C	7/10 (70%)	7 (100%)	0	0	100	100
2	D	8/10 (80%)	6 (75%)	2 (25%)	0	100	100

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1096/1204 (91%)	860 (78%)	190 (17%)	46 (4%)	3	18

All (46) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	162	PRO
1	A	199	LYS
1	A	304	ILE
1	A	454	GLN
1	A	528	ASN
1	B	130	ILE
1	B	199	LYS
1	B	222	ASP
1	B	383	VAL
1	B	432	ILE
1	B	521	TYR
1	B	522	LEU
1	A	128	LYS
1	A	163	ASP
1	A	202	ASP
1	A	305	PRO
1	A	384	LYS
1	A	462	ASP
1	A	489	ARG
1	A	525	ASP
1	B	162	PRO
1	B	430	VAL
1	B	454	GLN
1	A	146	GLN
1	A	361	ASN
1	A	496	PRO
1	A	521	TYR
1	B	164	PRO
1	B	381	GLN
1	B	384	LYS
1	A	164	PRO
1	A	287	ASP
1	A	307	LEU
1	B	379	GLU
1	B	508	SER
1	A	465	ALA
1	B	287	ASP

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type
1	A	130	ILE
1	A	324	LEU
1	B	223	GLU
1	B	393	VAL
1	A	450	VAL
1	B	392	PRO
1	A	311	ILE
1	A	380	GLY
1	B	509	GLY

### 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	433/540 (80%)	430 (99%)	3 (1%)	88	95
1	B	428/540 (79%)	422 (99%)	6 (1%)	74	91
2	C	8/10 (80%)	8 (100%)	0	100	100
2	D	8/10 (80%)	8 (100%)	0	100	100
All	All	877/1100 (80%)	868 (99%)	9 (1%)	82	93

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	180	GLN
1	A	197	LYS
1	A	450	VAL
1	B	180	GLN
1	B	240	ARG
1	B	254	PHE
1	B	396	ASN
1	B	422	ASN
1	B	450	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:



Mol	Chain	Res	Type
1	A	146	GLN
1	A	166	ASN
1	B	146	GLN
1	B	250	HIS
1	B	349	ASN
1	B	396	ASN
1	B	422	ASN
1	B	470	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	552/592 (93%)	-0.17	5 (0%) 85 69	74, 150, 225, 327	55 (9%)
1	B	545/592 (92%)	-0.12	5 (0%) 85 69	77, 146, 221, 291	45 (8%)
2	C	9/10 (90%)	0.63	1 (11%) 7 2	115, 124, 164, 176	0
2	D	10/10 (100%)	-0.00	0 100 100	120, 147, 199, 219	0
All	All	1116/1204 (92%)	-0.14	11 (0%) 84 66	74, 148, 223, 327	100 (8%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	1	MET	6.9
1	A	105	ALA	3.2
1	B	194	VAL	3.2
1	A	194	VAL	3.1
1	B	429	ASN	2.8
1	B	280	ALA	2.6
1	A	529	GLY	2.5
1	B	445	LEU	2.3
1	A	374	LEU	2.2
1	A	354	CYS	2.1
1	B	525	ASP	2.0

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

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

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.



## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.