



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:05 PM GMT

PDB ID : 1NDF
Title : Carnitine Acetyltransferase in Complex with Carnitine
Authors : Jogl, G.; Tong, L.
Deposited on : 2002-12-09
Resolution : 1.90 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

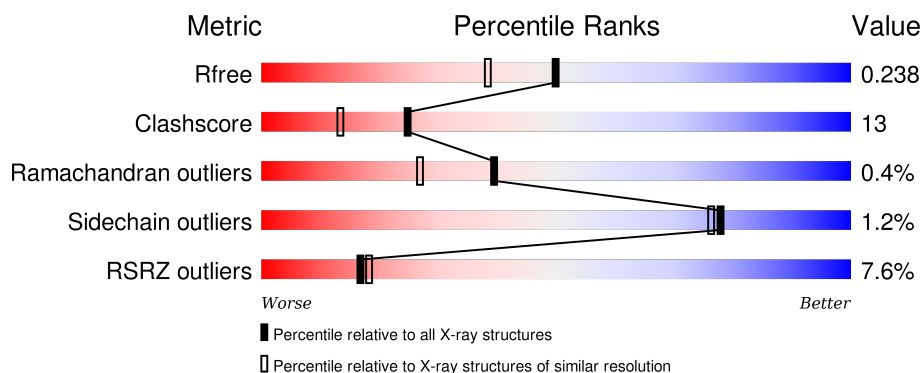
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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	596	<div> <div>12%</div> <div>74%</div> <div>26%</div> </div>
1	B	596	<div> <div>4%</div> <div>78%</div> <div>21%</div> </div>

2 Entry composition [i](#)

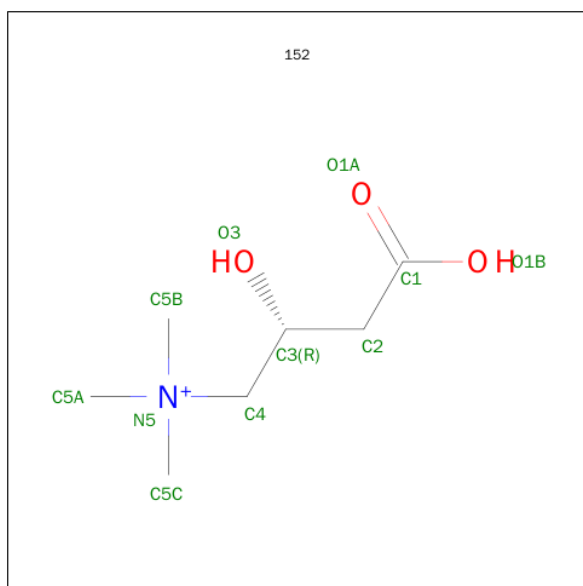
There are 3 unique types of molecules in this entry. The entry contains 10606 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Carnitine Acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	596	Total	C	N	O	S	0	0	0
			4757	3034	820	875	28			
1	B	596	Total	C	N	O	S	0	0	0
			4757	3034	820	875	28			

- Molecule 2 is CARNITINE (three-letter code: 152) (formula: $C_7H_{16}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			11	7	1	3		
2	B	1	Total	C	N	O	0	0
			11	7	1	3		

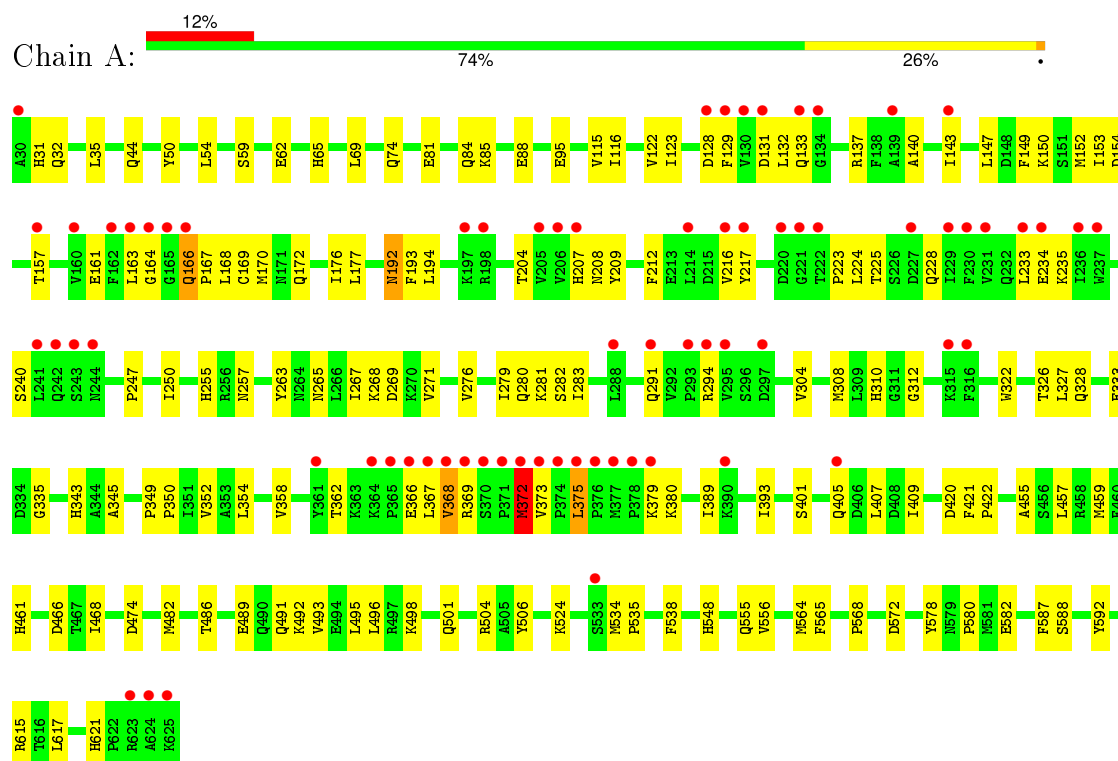
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	469	Total 469	O 469	0	0
3	B	601	Total 601	O 601	0	0

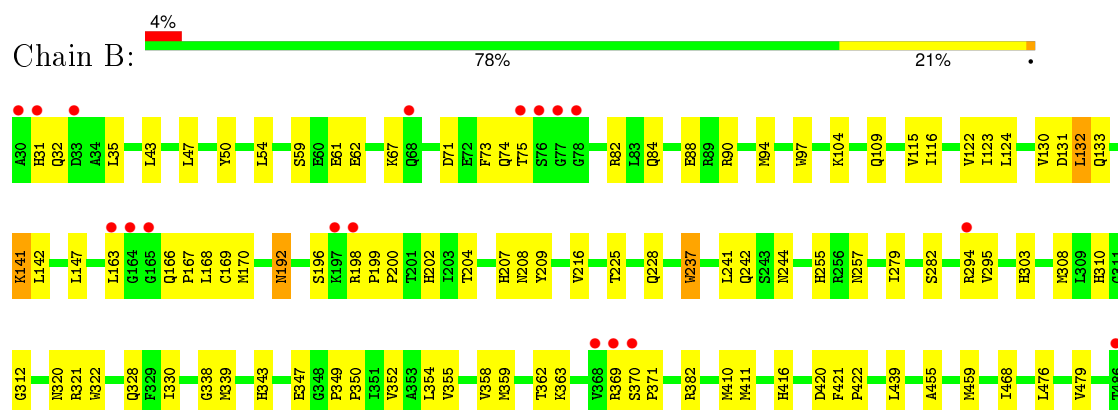
3 Residue-property plots

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: Carnitine Acetyltransferase



• Molecule 1: Carnitine Acetyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	160.68 Å 91.67 Å 122.61 Å 90.00° 128.84° 90.00°	Depositor
Resolution (Å)	19.95 – 1.90 19.94 – 1.89	Depositor EDS
% Data completeness (in resolution range)	96.7 (19.95-1.90) 96.3 (19.94-1.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.16 (at 1.89 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.200 , 0.241 0.198 , 0.238	Depositor DCC
R_{free} test set	7946 reflections (7.54%)	DCC
Wilson B-factor (Å ²)	17.9	Xtriage
Anisotropy	0.525	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 54.8	EDS
Estimated twinning fraction	0.019 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 108440 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10606	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.80% 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: 152

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	0.31	0/4872	0.55	0/6600
1	B	0.33	0/4872	0.58	0/6600
All	All	0.32	0/9744	0.56	0/13200

There are no bond length outliers.

There are no bond angle outliers.

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	4757	0	4733	134	0
1	B	4757	0	4733	108	0
2	A	11	0	15	2	0
2	B	11	0	15	1	0
3	A	469	0	0	8	0
3	B	601	0	0	10	0
All	All	10606	0	9496	244	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 13.

All (244) 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:161:GLU:H	1:A:170:MET:HE3	1.20	1.06
1:B:132:LEU:H	1:B:132:LEU:HD22	1.29	0.96
1:B:308:MET:HE3	1:B:330:ILE:HD12	1.48	0.96
1:B:94:MET:CE	1:B:97:TRP:HA	2.09	0.83
1:A:310:HIS:HD2	1:A:312:GLY:H	1.25	0.82
1:B:310:HIS:HD2	1:B:312:GLY:H	1.29	0.79
1:B:308:MET:HE1	1:B:330:ILE:HB	1.64	0.79
1:A:489:GLU:O	1:A:493:VAL:HG23	1.83	0.78
1:A:504:ARG:HB3	1:A:504:ARG:HH11	1.49	0.78
1:B:90:ARG:NE	1:B:94:MET:HE1	1.98	0.77
1:B:141:LYS:HD2	1:B:362:THR:HA	1.68	0.75
1:A:534:MET:CE	1:A:535:PRO:HD2	2.17	0.72
1:B:31:HIS:HE1	1:B:167:PRO:HB3	1.54	0.72
1:A:161:GLU:N	1:A:170:MET:HE3	2.02	0.72
1:B:493:VAL:HG22	1:B:617:LEU:HG	1.73	0.70
1:A:122:VAL:HG22	1:A:564:MET:HG2	1.74	0.70
1:B:32:GLN:HE22	1:B:170:MET:H	1.38	0.70
1:B:534:MET:HB3	1:B:539:MET:HE2	1.74	0.69
1:B:310:HIS:CD2	1:B:312:GLY:H	2.09	0.69
1:A:228:GLN:HB3	1:A:375:LEU:HD11	1.74	0.69
1:A:493:VAL:HG22	1:A:617:LEU:HG	1.75	0.68
1:B:349:PRO:O	1:B:352:VAL:HG22	1.94	0.68
1:B:555:GLN:HE21	1:B:580:PRO:HG2	1.59	0.68
1:A:457:LEU:HG	3:A:5100:HOH:O	1.94	0.68
1:A:265:ASN:HA	1:A:268:LYS:HE3	1.77	0.67
1:A:310:HIS:CD2	1:A:312:GLY:H	2.11	0.67
1:B:131:ASP:OD1	1:B:133:GLN:HG3	1.94	0.66
1:A:504:ARG:HB3	1:A:504:ARG:NH1	2.10	0.66
1:A:32:GLN:HE22	1:A:170:MET:H	1.44	0.66
1:A:482:MET:HE2	1:A:492:LYS:HD3	1.78	0.66
1:A:204:THR:HG23	1:A:282:SER:HB3	1.78	0.65
1:A:534:MET:HE2	1:A:535:PRO:HD2	1.77	0.65
1:A:255:HIS:HD2	1:A:257:ASN:H	1.44	0.65
1:B:497:ARG:O	1:B:501:GLN:HG2	1.97	0.65
1:B:94:MET:HE3	1:B:97:TRP:HA	1.78	0.64
1:B:308:MET:HE2	1:B:339:MET:N	2.13	0.64
1:B:47:LEU:HD12	1:B:74:GLN:HB3	1.78	0.64
1:A:482:MET:CE	1:A:492:LYS:HB3	2.28	0.63
1:B:308:MET:CE	1:B:330:ILE:HD12	2.26	0.63
1:A:349:PRO:HB2	1:A:350:PRO:HD3	1.80	0.62
1:B:534:MET:HB3	1:B:539:MET:CE	2.29	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:491:GLN:HE21	1:A:495:LEU:HG	1.65	0.62
1:A:407:LEU:HD21	1:A:409:ILE:HD11	1.79	0.62
1:A:247:PRO:O	1:A:250:ILE:HG22	1.99	0.62
1:A:31:HIS:HE1	1:A:167:PRO:HB3	1.64	0.61
1:A:421:PHE:HB3	1:A:422:PRO:HD3	1.82	0.61
1:A:482:MET:CE	1:A:482:MET:HA	2.31	0.61
1:B:132:LEU:CD2	1:B:132:LEU:H	2.06	0.60
1:B:90:ARG:HG2	1:B:94:MET:HE2	1.83	0.60
1:A:225:THR:H	1:A:228:GLN:HE21	1.50	0.60
1:B:370:SER:HB3	1:B:371:PRO:CD	2.31	0.59
1:A:482:MET:HE2	1:A:492:LYS:HB3	1.85	0.59
1:B:47:LEU:CD1	1:B:74:GLN:HB3	2.31	0.59
1:A:225:THR:OG1	1:A:228:GLN:HG3	2.03	0.59
1:A:276:VAL:O	1:A:280:GLN:HG3	2.03	0.59
1:A:216:VAL:C	1:A:224:LEU:HD13	2.23	0.59
1:A:493:VAL:HG22	1:A:617:LEU:CG	2.32	0.58
1:B:369:ARG:N	3:B:6100:HOH:O	2.36	0.58
1:A:84:GLN:O	1:A:88:GLU:HG3	2.04	0.57
1:B:512:ARG:HG3	1:B:512:ARG:HH11	1.68	0.57
1:B:322:TRP:H	1:B:328:GLN:NE2	2.02	0.57
1:B:421:PHE:HB3	1:B:422:PRO:HD3	1.86	0.57
1:A:524:LYS:HG3	1:A:534:MET:HE1	1.86	0.57
1:A:123:ILE:HG13	1:A:565:PHE:CE2	2.40	0.57
1:B:94:MET:HE2	1:B:97:TRP:HA	1.87	0.57
1:B:294:ARG:HG3	3:B:6232:HOH:O	2.04	0.57
1:B:255:HIS:HD2	1:B:257:ASN:H	1.51	0.57
2:A:5001:152:H5B3	2:A:5001:152:O3	2.05	0.56
1:B:59:SER:OG	1:B:62:GLU:HG3	2.05	0.56
1:A:163:LEU:HB3	1:A:166:GLN:OE1	2.05	0.56
1:A:131:ASP:OD1	1:A:133:GLN:HG3	2.05	0.56
1:A:163:LEU:O	1:A:163:LEU:HD13	2.06	0.55
1:B:504:ARG:HB3	1:B:504:ARG:HH11	1.71	0.55
1:A:493:VAL:HG22	1:A:617:LEU:CD1	2.36	0.55
1:A:482:MET:HE2	1:A:482:MET:HA	1.89	0.55
1:B:207:HIS:HE1	3:B:6426:HOH:O	1.90	0.55
1:B:568:PRO:HD2	1:B:592:TYR:CZ	2.42	0.55
1:B:294:ARG:HG3	1:B:294:ARG:HH11	1.72	0.55
1:B:382:ARG:NH1	1:B:382:ARG:HB2	2.21	0.55
1:A:534:MET:HE3	1:A:535:PRO:HD2	1.89	0.55
1:B:370:SER:HB3	1:B:371:PRO:HD2	1.88	0.55
1:A:498:LYS:HD3	3:A:5054:HOH:O	2.07	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:555:GLN:HE21	1:A:580:PRO:HG2	1.72	0.54
1:B:90:ARG:HG2	1:B:94:MET:CE	2.38	0.54
1:B:358:VAL:O	1:B:362:THR:HG23	2.07	0.54
1:A:352:VAL:HG11	1:A:556:VAL:HG12	1.90	0.54
1:A:132:LEU:HD22	1:A:132:LEU:H	1.73	0.54
1:A:50:TYR:CZ	1:A:54:LEU:HD11	2.43	0.54
1:A:228:GLN:HG2	1:A:372:MET:HB3	1.90	0.54
1:A:501:GLN:NE2	3:A:5287:HOH:O	2.42	0.53
1:B:347:GLU:O	1:B:350:PRO:HD2	2.09	0.53
1:A:129:PHE:CZ	1:A:335:GLY:HA2	2.43	0.53
1:A:294:ARG:HG3	3:A:5409:HOH:O	2.08	0.53
1:A:255:HIS:CD2	1:A:257:ASN:H	2.25	0.53
1:B:548:HIS:HE1	1:B:572:ASP:OD1	1.91	0.52
1:A:349:PRO:HA	1:A:352:VAL:HG22	1.91	0.52
1:B:410:MET:HE3	1:B:411:MET:O	2.09	0.52
1:A:269:ASP:OD2	1:A:271:VAL:HB	2.10	0.52
1:A:486:THR:HG21	1:B:196:SER:HB3	1.91	0.52
1:A:568:PRO:HD2	1:A:592:TYR:CZ	2.45	0.52
1:A:193:PHE:CE2	1:A:281:LYS:HG2	2.45	0.52
1:A:132:LEU:HD22	1:A:132:LEU:N	2.25	0.52
1:A:115:VAL:HG12	1:A:116:ILE:HG13	1.90	0.52
1:A:291:GLN:HB2	1:A:333:GLU:HG3	1.92	0.52
1:B:308:MET:CE	1:B:330:ILE:HB	2.37	0.51
1:B:35:LEU:HD11	1:B:169:CYS:HA	1.93	0.51
1:B:132:LEU:HD22	1:B:132:LEU:N	2.11	0.51
1:B:90:ARG:HE	1:B:94:MET:HE1	1.74	0.51
1:A:152:MET:HA	1:A:157:THR:OG1	2.11	0.51
1:A:455:ALA:HB2	1:A:468:ILE:HG13	1.93	0.51
1:A:44:GLN:HA	1:A:74:GLN:HE22	1.76	0.51
1:B:539:MET:HE3	3:B:6211:HOH:O	2.10	0.51
1:A:225:THR:HG23	1:A:228:GLN:HE21	1.75	0.51
1:B:104:LYS:HD2	1:B:109:GLN:OE1	2.12	0.50
2:B:6001:152:O3	2:B:6001:152:H5B3	2.10	0.50
1:B:115:VAL:HG12	1:B:116:ILE:HG13	1.94	0.50
1:A:368:VAL:HG23	1:A:368:VAL:O	2.12	0.50
1:A:177:LEU:HD23	1:A:327:LEU:HD12	1.93	0.50
1:A:457:LEU:HD21	1:A:466:ASP:HB2	1.94	0.49
1:A:132:LEU:CD2	1:A:132:LEU:H	2.25	0.49
1:B:202:HIS:HD2	3:B:6406:HOH:O	1.95	0.49
1:B:204:THR:HG21	1:B:279:ILE:HA	1.93	0.49
1:B:421:PHE:CZ	1:B:615:ARG:HG3	2.46	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:326:THR:HG23	1:A:345:ALA:HB3	1.93	0.49
1:B:82:ARG:HB3	3:B:6590:HOH:O	2.11	0.49
1:A:234:GLU:HB3	3:A:5184:HOH:O	2.13	0.49
1:A:59:SER:OG	1:A:62:GLU:HG3	2.12	0.49
1:A:123:ILE:HG13	1:A:565:PHE:HE2	1.78	0.49
1:A:81:GLU:O	1:A:85:LYS:HG2	2.11	0.49
1:B:339:MET:CE	1:B:355:VAL:HG22	2.43	0.49
1:A:163:LEU:HB2	1:A:168:LEU:HD21	1.94	0.49
1:A:322:TRP:H	1:A:328:GLN:HE22	1.59	0.49
1:A:366:GLU:HG2	1:A:369:ARG:NH2	2.29	0.48
1:B:493:VAL:HG12	1:B:497:ARG:NH1	2.28	0.48
1:B:142:LEU:C	1:B:142:LEU:HD23	2.34	0.48
1:A:366:GLU:HA	1:A:369:ARG:HH21	1.79	0.48
1:B:122:VAL:HG22	1:B:564:MET:HG2	1.94	0.48
1:A:204:THR:HG21	1:A:279:ILE:HA	1.95	0.48
1:A:366:GLU:CB	1:A:369:ARG:HH21	2.26	0.48
1:A:95:GLU:OE1	1:A:461:HIS:HE1	1.96	0.48
1:A:35:LEU:HD11	1:A:169:CYS:HA	1.96	0.48
1:A:493:VAL:HG22	1:A:617:LEU:HD11	1.96	0.48
1:A:578:TYR:HB3	1:A:587:PHE:CD1	2.48	0.48
1:B:192:ASN:C	1:B:192:ASN:HD22	2.17	0.48
1:B:84:GLN:O	1:B:88:GLU:HG3	2.14	0.48
1:A:217:TYR:CE1	1:A:223:PRO:HB3	2.49	0.47
1:A:32:GLN:NE2	1:A:170:MET:H	2.12	0.47
1:B:382:ARG:CZ	1:B:382:ARG:HB2	2.44	0.47
1:A:496:LEU:C	1:A:496:LEU:HD23	2.34	0.47
1:B:455:ALA:HB2	1:B:468:ILE:HG13	1.96	0.47
1:B:308:MET:CE	1:B:339:MET:N	2.77	0.47
1:A:548:HIS:HE1	1:A:572:ASP:OD1	1.97	0.47
1:A:176:ILE:HB	1:A:327:LEU:HD11	1.97	0.46
1:B:496:LEU:C	1:B:496:LEU:HD23	2.35	0.46
1:B:416:HIS:CG	1:B:612:LEU:HD21	2.50	0.46
1:B:166:GLN:HA	1:B:167:PRO:HD3	1.84	0.46
1:A:225:THR:HG23	1:A:228:GLN:NE2	2.31	0.46
1:B:568:PRO:HB2	1:B:592:TYR:CE2	2.51	0.46
1:A:380:LYS:HD2	3:A:5425:HOH:O	2.14	0.46
1:A:143:ILE:O	1:A:147:LEU:HG	2.16	0.46
1:B:578:TYR:HB3	1:B:587:PHE:CD1	2.51	0.46
1:A:207:HIS:NE2	1:A:208:ASN:ND2	2.64	0.46
1:B:349:PRO:HB2	1:B:350:PRO:HD3	1.99	0.45
1:B:43:LEU:O	1:B:47:LEU:HG	2.16	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:367:LEU:O	1:A:369:ARG:HG2	2.16	0.45
1:B:130:VAL:HA	3:B:6223:HOH:O	2.16	0.45
1:B:73:PHE:C	1:B:75:THR:H	2.18	0.45
1:A:326:THR:HG23	1:A:345:ALA:CB	2.47	0.45
1:A:263:TYR:O	1:A:267:ILE:HG12	2.17	0.45
1:A:140:ALA:HB2	1:A:233:LEU:HD12	1.99	0.45
1:A:322:TRP:H	1:A:328:GLN:NE2	2.15	0.45
1:A:401:SER:O	1:A:405:GLN:HG3	2.17	0.45
1:B:410:MET:HG2	1:B:601:ALA:HA	1.97	0.45
1:A:166:GLN:N	1:A:166:GLN:NE2	2.65	0.45
1:B:255:HIS:CD2	1:B:257:ASN:H	2.33	0.45
1:B:354:LEU:HD23	1:B:354:LEU:C	2.38	0.44
1:A:421:PHE:CZ	1:A:615:ARG:HG3	2.52	0.44
1:A:349:PRO:O	1:A:352:VAL:HG22	2.16	0.44
1:A:225:THR:N	1:A:228:GLN:HE21	2.14	0.44
1:A:172:GLN:OE1	1:A:350:PRO:HG2	2.18	0.44
1:A:65:HIS:CE1	1:A:69:LEU:HD11	2.53	0.44
1:A:150:LYS:HE3	1:A:283:ILE:HD13	2.00	0.44
1:B:124:LEU:HD11	1:B:339:MET:CE	2.47	0.44
1:A:166:GLN:HG3	1:A:459:MET:HE1	2.00	0.44
1:A:150:LYS:HG2	1:A:154:ASP:OD2	2.18	0.44
1:A:294:ARG:HH11	1:A:294:ARG:HG3	1.83	0.43
1:B:295:VAL:HG11	1:B:303:HIS:CD2	2.54	0.43
1:A:534:MET:HE1	1:A:538:PHE:CD1	2.54	0.43
1:B:512:ARG:NH1	1:B:512:ARG:HG3	2.31	0.43
1:A:176:ILE:HA	1:A:326:THR:HG21	2.00	0.43
1:B:320:ASN:C	1:B:321:ARG:HG2	2.38	0.43
1:B:539:MET:HE1	3:B:6523:HOH:O	2.18	0.43
1:B:439:LEU:HA	1:B:479:VAL:HG12	2.01	0.43
1:B:459:MET:HE3	1:B:511:ILE:HG22	2.00	0.43
1:B:242:GLN:HB3	1:B:244:ASN:OD1	2.19	0.43
1:B:204:THR:HG23	1:B:282:SER:HB3	2.00	0.43
1:A:489:GLU:HB3	1:A:621:HIS:NE2	2.33	0.42
1:A:474:ASP:OD1	1:A:498:LYS:HE2	2.19	0.42
1:B:50:TYR:CZ	1:B:54:LEU:HD11	2.55	0.42
1:B:61:GLU:H	1:B:61:GLU:CD	2.22	0.42
1:B:123:ILE:HG13	1:B:565:PHE:CE2	2.55	0.42
1:B:308:MET:CE	1:B:338:GLY:C	2.88	0.42
1:B:322:TRP:H	1:B:328:GLN:HE22	1.66	0.42
1:A:565:PHE:HB3	1:A:588:SER:HB2	2.01	0.42
1:B:82:ARG:HD2	3:B:6590:HOH:O	2.18	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:147:LEU:HD11	1:B:216:VAL:HB	2.02	0.42
1:B:163:LEU:HB2	1:B:168:LEU:HD11	2.02	0.42
1:A:304:VAL:O	1:A:308:MET:HG2	2.20	0.42
1:B:310:HIS:HE1	3:B:6396:HOH:O	2.03	0.42
1:A:358:VAL:O	1:A:362:THR:HG23	2.20	0.42
1:B:208:ASN:O	1:B:209:TYR:HB2	2.19	0.42
1:A:534:MET:HE2	1:A:535:PRO:CD	2.47	0.42
1:B:32:GLN:NE2	1:B:170:MET:H	2.11	0.42
1:A:235:LYS:NZ	1:A:375:LEU:O	2.53	0.41
1:B:225:THR:H	1:B:228:GLN:NE2	2.17	0.41
1:A:212:PHE:CZ	1:A:240:SER:HB3	2.54	0.41
1:A:192:ASN:C	1:A:192:ASN:HD22	2.22	0.41
1:A:482:MET:HE1	1:A:492:LYS:HB3	1.98	0.41
2:A:5001:152:H5A2	2:A:5001:152:H3	1.82	0.41
1:A:354:LEU:C	1:A:354:LEU:HD23	2.41	0.41
1:A:482:MET:HB2	1:A:482:MET:HE3	1.93	0.41
1:A:166:GLN:HA	1:A:167:PRO:HD3	1.89	0.41
1:A:128:ASP:O	1:A:129:PHE:C	2.59	0.41
1:A:209:TYR:CD2	1:A:209:TYR:N	2.89	0.41
1:A:420:ASP:HB3	1:A:582:GLU:OE2	2.21	0.41
1:A:310:HIS:HE1	3:A:5315:HOH:O	2.02	0.41
1:B:420:ASP:OD1	1:B:421:PHE:N	2.54	0.41
1:B:82:ARG:HH11	1:B:82:ARG:HG3	1.86	0.41
1:B:67:LYS:HE2	1:B:71:ASP:OD2	2.20	0.41
1:B:621:HIS:N	1:B:622:PRO:HD3	2.35	0.41
1:B:308:MET:HE1	1:B:338:GLY:C	2.41	0.40
1:A:192:ASN:ND2	1:A:194:LEU:H	2.18	0.40
1:A:372:MET:HB2	1:A:373:VAL:H	1.64	0.40
1:A:255:HIS:CD2	1:A:257:ASN:HB2	2.56	0.40
1:B:294:ARG:HG3	1:B:294:ARG:NH1	2.35	0.40
1:A:389:ILE:O	1:A:393:ILE:HG13	2.22	0.40
1:A:373:VAL:O	1:A:375:LEU:N	2.44	0.40
1:A:421:PHE:CE1	1:A:615:ARG:HG3	2.57	0.40
1:B:199:PRO:HA	1:B:200:PRO:HD3	1.93	0.40
1:A:149:PHE:HE1	1:A:153:ILE:HD11	1.87	0.40
1:A:379:LYS:HG3	3:A:5451:HOH:O	2.22	0.40
1:B:476:LEU:HA	1:B:479:VAL:HG22	2.03	0.40
1:B:237:TRP:CE2	1:B:241:LEU:HD21	2.57	0.40
1:B:359:MET:O	1:B:363:LYS:HG2	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	594/596 (100%)	569 (96%)	21 (4%)	4 (1%)	26	14
1	B	594/596 (100%)	575 (97%)	18 (3%)	1 (0%)	52	42
All	All	1188/1192 (100%)	1144 (96%)	39 (3%)	5 (0%)	39	27

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	372	MET
1	B	624	ALA
1	A	164	GLY
1	A	375	LEU
1	A	368	VAL

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	524/524 (100%)	518 (99%)	6 (1%)	80	79
1	B	524/524 (100%)	517 (99%)	7 (1%)	76	73
All	All	1048/1048 (100%)	1035 (99%)	13 (1%)	78	76

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	137	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	166	GLN
1	A	192	ASN
1	A	343	HIS
1	A	372	MET
1	A	506	TYR
1	B	132	LEU
1	B	141	LYS
1	B	192	ASN
1	B	198	ARG
1	B	237	TRP
1	B	343	HIS
1	B	506	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (51) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	HIS
1	A	32	GLN
1	A	74	GLN
1	A	84	GLN
1	A	112	GLN
1	A	135	GLN
1	A	187	GLN
1	A	192	ASN
1	A	202	HIS
1	A	208	ASN
1	A	210	GLN
1	A	228	GLN
1	A	255	HIS
1	A	277	ASN
1	A	307	GLN
1	A	310	HIS
1	A	328	GLN
1	A	461	HIS
1	A	491	GLN
1	A	503	HIS
1	A	548	HIS
1	A	550	ASN
1	A	555	GLN
1	A	586	ASN
1	A	619	GLN
1	A	620	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	32	GLN
1	B	44	GLN
1	B	68	GLN
1	B	84	GLN
1	B	109	GLN
1	B	112	GLN
1	B	135	GLN
1	B	166	GLN
1	B	192	ASN
1	B	202	HIS
1	B	207	HIS
1	B	228	GLN
1	B	255	HIS
1	B	307	GLN
1	B	310	HIS
1	B	328	GLN
1	B	357	HIS
1	B	461	HIS
1	B	503	HIS
1	B	548	HIS
1	B	550	ASN
1	B	555	GLN
1	B	586	ASN
1	B	619	GLN
1	B	620	ASN

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

2 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	152	A	5001	-	7,10,10	1.25	1 (14%)	9,14,14	1.12	1 (11%)
2	152	B	6001	-	7,10,10	1.21	1 (14%)	9,14,14	1.01	1 (11%)

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	152	A	5001	-	-	0/7/9/9	0/0/0/0
2	152	B	6001	-	-	0/7/9/9	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	6001	152	C4-N5	-2.27	1.47	1.51
2	A	5001	152	C4-N5	-2.21	1.47	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	5001	152	C3-C4-N5	-3.26	110.75	117.24
2	B	6001	152	C3-C4-N5	-2.89	111.49	117.24

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	A	5001	152	2	0
2	B	6001	152	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	596/596 (100%)	0.49	70 (11%) 6 7	9, 23, 57, 90	0
1	B	596/596 (100%)	-0.05	21 (3%) 48 51	7, 17, 39, 73	0
All	All	1192/1192 (100%)	0.22	91 (7%) 17 18	7, 19, 53, 90	0

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	624	ALA	17.2
1	A	368	VAL	11.5
1	A	625	LYS	10.8
1	A	367	LEU	10.2
1	A	373	VAL	9.8
1	A	370	SER	8.9
1	B	77	GLY	8.7
1	B	76	SER	8.3
1	B	625	LYS	7.5
1	A	624	ALA	7.3
1	A	371	PRO	5.8
1	A	198	ARG	5.7
1	A	237	TRP	5.5
1	A	369	ARG	5.5
1	B	78	GLY	5.3
1	A	163	LEU	5.3
1	A	165	GLY	5.1
1	A	375	LEU	5.0
1	A	372	MET	5.0
1	A	365	PRO	4.9
1	A	366	GLU	4.8
1	B	33	ASP	4.7
1	B	370	SER	4.4
1	A	227	ASP	4.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	294	ARG	4.1
1	A	374	PRO	4.0
1	A	166	GLN	3.8
1	A	130	VAL	3.8
1	B	198	ARG	3.7
1	A	231	VAL	3.7
1	B	31	HIS	3.6
1	B	30	ALA	3.6
1	A	128	ASP	3.6
1	A	220	ASP	3.6
1	B	369	ARG	3.6
1	B	75	THR	3.5
1	A	293	PRO	3.4
1	A	242	GLN	3.4
1	B	623	ARG	3.4
1	A	243	SER	3.3
1	A	316	PHE	3.2
1	B	164	GLY	3.2
1	A	233	LEU	3.1
1	A	197	LYS	3.1
1	A	229	ILE	3.1
1	A	216	VAL	3.1
1	A	214	LEU	3.1
1	B	165	GLY	3.1
1	A	390	LYS	3.0
1	A	30	ALA	3.0
1	A	221	GLY	3.0
1	A	162	PHE	3.0
1	A	160	VAL	3.0
1	A	139	ALA	2.9
1	A	222	THR	2.9
1	A	217	TYR	2.9
1	A	134	GLY	2.8
1	A	623	ARG	2.8
1	A	379	LYS	2.8
1	A	241	LEU	2.8
1	A	236	ILE	2.7
1	B	197	LYS	2.6
1	A	129	PHE	2.6
1	B	163	LEU	2.6
1	A	157	THR	2.6
1	A	143	ILE	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	361	TYR	2.5
1	B	486	THR	2.5
1	B	368	VAL	2.5
1	A	533	SER	2.4
1	A	376	PRO	2.4
1	A	133	GLN	2.4
1	A	131	ASP	2.4
1	A	377	MET	2.4
1	A	405	GLN	2.4
1	A	297	ASP	2.4
1	A	291	GLN	2.3
1	A	164	GLY	2.3
1	A	288	LEU	2.3
1	A	364	LYS	2.2
1	A	315	LYS	2.2
1	A	244	ASN	2.2
1	A	230	PHE	2.2
1	A	207	HIS	2.1
1	A	295	VAL	2.1
1	A	206	VAL	2.1
1	B	294	ARG	2.1
1	A	205	VAL	2.1
1	B	68	GLN	2.1
1	A	378	PRO	2.0
1	A	234	GLU	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 ⓘ

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	152	B	6001	11/11	0.96	0.07	-0.19	12,14,17,18	0
2	152	A	5001	11/11	0.96	0.08	-0.20	9,14,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.