



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 04:18 PM GMT

PDB ID : 2H3U
Title : Crystal structure of murine carnitine acetyltransferase in complex with carnitine and CoA
Authors : Hsiao, Y.S.; Jogl, G.; Tong, L.
Deposited on : 2006-05-23
Resolution : 1.90 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

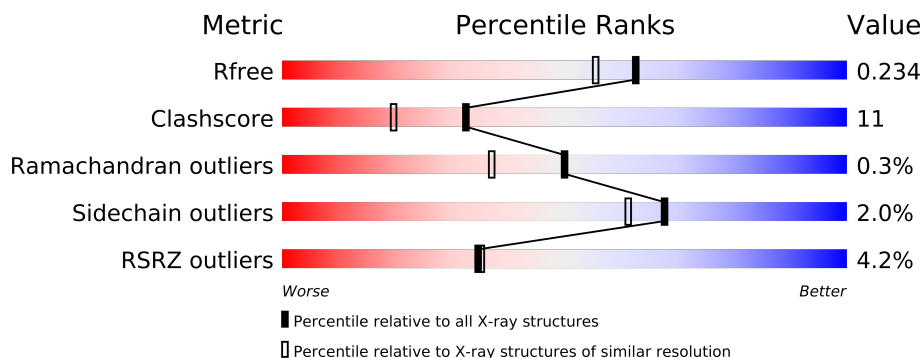
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	3684 (1.90-1.90)
Clashscore	79885	4465 (1.90-1.90)
Ramachandran outliers	78287	4413 (1.90-1.90)
Sidechain outliers	78261	4414 (1.90-1.90)
RSRZ outliers	66119	3686 (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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	599	
1	B	599	

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 11366 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

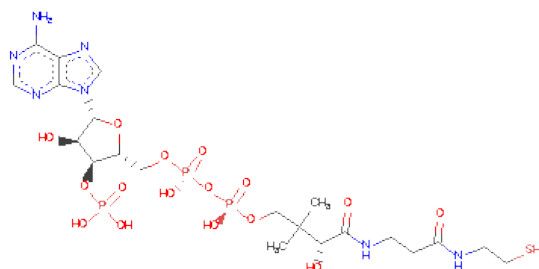
- Molecule 1 is a protein called carnitine acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	599	Total	C	N	O	S	0	0	0
			4781	3048	825	879	29			
1	B	599	Total	C	N	O	S	0	0	0
			4781	3048	825	879	29			

There are 6 discrepancies between the modelled and reference sequences:

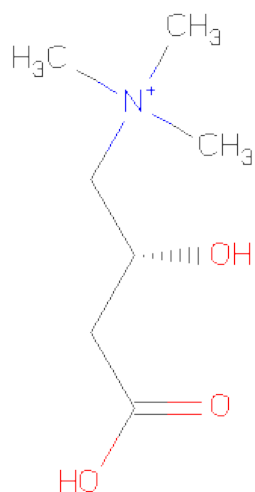
Chain	Residue	Modelled	Actual	Comment	Reference
A	27	SER	-	CLONING ARTIFACT	UNP Q3V1Y3
A	28	HIS	-	CLONING ARTIFACT	UNP Q3V1Y3
A	29	MET	-	CLONING ARTIFACT	UNP Q3V1Y3
B	27	SER	-	CLONING ARTIFACT	UNP Q3V1Y3
B	28	HIS	-	CLONING ARTIFACT	UNP Q3V1Y3
B	29	MET	-	CLONING ARTIFACT	UNP Q3V1Y3

- Molecule 2 is COENZYME A (three-letter code: COA) (formula: $C_{21}H_{36}N_7O_{16}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		
2	B	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

- Molecule 3 is CARNITINE (three-letter code: 152) (formula: $C_7H_{16}NO_3$).



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

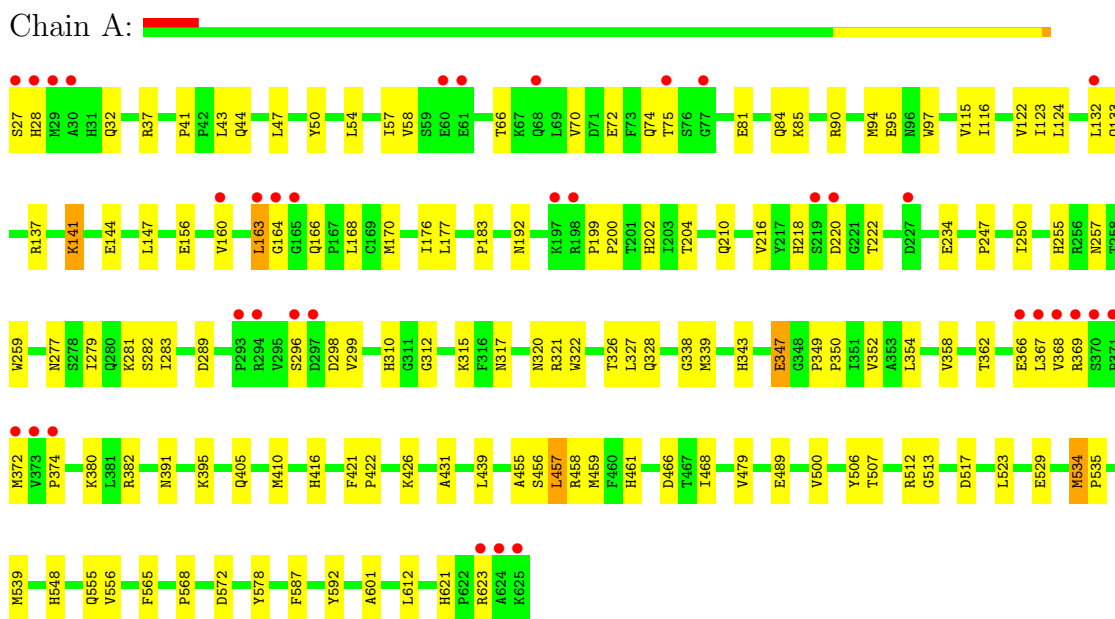
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	737	Total	O	0	0
			737	737		
4	B	949	Total	O	0	0
			949	949		

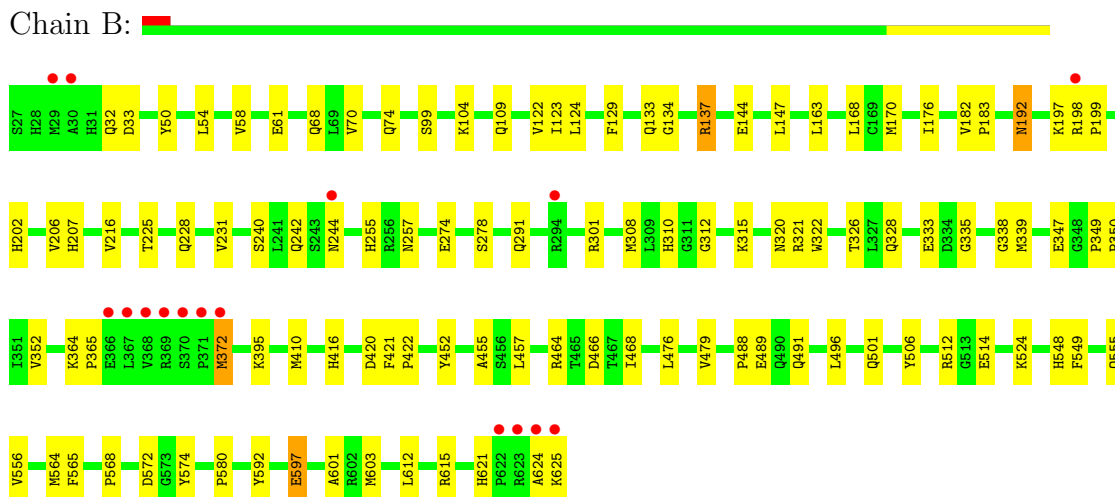
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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	164.51Å 89.74Å 122.95Å 90.00° 128.96° 90.00°	Depositor
Resolution (Å)	29.21 – 1.90 29.21 – 1.90	Depositor EDS
% Data completeness (in resolution range)	92.5 (29.21-1.90) 92.7 (29.21-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.96 (at 1.91Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.193 , 0.235 0.193 , 0.234	Depositor DCC
R_{free} test set	7625 reflections (7.53%)	DCC
Wilson B-factor (Å ²)	16.1	Xtriage
Anisotropy	0.533	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 44.5	EDS
Estimated twinning fraction	0.012 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	4 of 106811 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11366	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 3.99% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: COA, 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	0/4897	0.53	0/6633
1	B	0.35	0/4897	0.58	0/6633
All	All	0.33	0/9794	0.55	0/13266

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4781	0	4754	128	0
1	B	4781	0	4754	86	0
2	A	48	0	32	4	0
2	B	48	0	32	2	0
3	A	11	0	15	1	0
3	B	11	0	15	2	0
4	A	737	0	0	46	0
4	B	949	0	0	24	0
All	All	11366	0	9602	217	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 11.

All (217) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:141:LYS:HD2	1:A:362:THR:HA	1.46	0.96
3:B:1901:152:H5A1	4:B:2511:HOH:O	1.70	0.92
1:A:457:LEU:HD13	4:A:2014:HOH:O	1.68	0.92
1:A:534:MET:HE2	1:A:535:PRO:HD2	1.48	0.91
1:A:90:ARG:HG2	4:A:2038:HOH:O	1.76	0.84
1:B:310:HIS:HD2	1:B:312:GLY:H	1.26	0.84
1:A:289:ASP:HA	4:A:2047:HOH:O	1.77	0.83
1:A:310:HIS:HD2	1:A:312:GLY:H	1.28	0.80
1:A:457:LEU:HD11	1:A:466:ASP:HB2	1.64	0.80
1:B:315:LYS:HG2	4:B:2683:HOH:O	1.81	0.80
1:B:291:GLN:HA	4:B:2747:HOH:O	1.83	0.79
1:A:320:ASN:HB3	4:A:2217:HOH:O	1.82	0.78
1:B:32:GLN:HE22	1:B:170:MET:H	1.33	0.76
1:A:457:LEU:CD1	1:A:466:ASP:HB2	2.15	0.76
1:A:349:PRO:HB2	1:A:350:PRO:HD3	1.68	0.76
1:A:461:HIS:HB2	4:A:2045:HOH:O	1.86	0.75
1:B:310:HIS:CD2	1:B:312:GLY:H	2.05	0.74
1:A:200:PRO:HB2	4:A:2157:HOH:O	1.87	0.74
1:B:501:GLN:HG2	4:B:2397:HOH:O	1.90	0.71
1:B:349:PRO:HD3	2:B:1902:COA:H31	1.70	0.71
1:B:333:GLU:HB3	4:B:2747:HOH:O	1.90	0.71
1:A:32:GLN:HE22	1:A:170:MET:H	1.36	0.70
1:B:349:PRO:HB2	1:B:350:PRO:HD3	1.73	0.69
1:A:426:LYS:HG2	1:A:623:ARG:HH22	1.57	0.69
1:A:349:PRO:HD3	2:A:1802:COA:H31	1.75	0.68
1:A:283:ILE:HG12	4:A:2157:HOH:O	1.92	0.68
1:B:464:ARG:HD3	4:B:2553:HOH:O	1.93	0.68
1:B:207:HIS:HD2	1:B:240:SER:OG	1.78	0.67
1:A:459:MET:HG2	4:A:2176:HOH:O	1.95	0.67
1:A:426:LYS:HB2	1:A:623:ARG:HH12	1.59	0.66
1:A:310:HIS:CD2	1:A:312:GLY:H	2.12	0.66
1:A:317:ASN:HB3	4:A:2217:HOH:O	1.96	0.66
1:A:457:LEU:HD23	4:A:2176:HOH:O	1.95	0.65
1:B:255:HIS:HD2	1:B:257:ASN:H	1.44	0.65
1:A:97:TRP:HA	4:A:2038:HOH:O	1.95	0.65
1:A:391:ASN:O	1:A:395:LYS:HG2	1.97	0.65
1:B:322:TRP:H	1:B:328:GLN:NE2	1.96	0.64
1:A:43:LEU:HD13	4:A:2015:HOH:O	1.96	0.64
1:A:534:MET:CE	1:A:535:PRO:HD2	2.27	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:321:ARG:HH11	1:A:321:ARG:HG2	1.63	0.63
1:A:534:MET:HA	1:A:534:MET:HE3	1.80	0.63
1:B:491:GLN:HG3	4:B:2552:HOH:O	1.98	0.62
1:B:556:VAL:HG22	2:B:1902:COA:H32	1.82	0.62
1:A:50:TYR:CZ	1:A:54:LEU:HD11	2.35	0.62
1:A:320:ASN:O	1:A:321:ARG:HG2	2.00	0.62
1:A:421:PHE:HB3	1:A:422:PRO:HD3	1.80	0.61
1:A:321:ARG:HB3	1:A:328:GLN:HE22	1.67	0.60
1:A:72:GLU:O	1:A:75:THR:HG22	2.01	0.60
1:B:421:PHE:HB3	1:B:422:PRO:HD3	1.83	0.60
1:A:322:TRP:H	1:A:328:GLN:NE2	2.00	0.59
1:A:366:GLU:HA	1:A:369:ARG:HH21	1.68	0.59
1:B:68:GLN:HG2	4:B:2344:HOH:O	2.01	0.58
1:A:132:LEU:HD22	1:A:132:LEU:H	1.67	0.58
1:B:147:LEU:HD11	1:B:216:VAL:HB	1.85	0.58
1:A:43:LEU:N	4:A:2039:HOH:O	2.35	0.57
1:B:372:MET:HA	4:B:2351:HOH:O	2.04	0.57
1:B:163:LEU:HB3	1:B:168:LEU:HD21	1.86	0.57
1:B:122:VAL:HB	1:B:339:MET:HG2	1.87	0.57
1:B:123:ILE:HG13	1:B:565:PHE:CE2	2.40	0.57
1:A:568:PRO:HD2	1:A:592:TYR:CZ	2.40	0.57
1:B:133:GLN:HE21	1:B:137:ARG:HH11	1.51	0.56
1:A:133:GLN:HG2	4:A:2000:HOH:O	2.06	0.56
1:A:512:ARG:HH11	1:A:512:ARG:HG3	1.70	0.56
1:A:366:GLU:HB3	1:A:369:ARG:HE	1.70	0.55
1:B:395:LYS:HD3	4:B:2826:HOH:O	2.05	0.55
1:A:177:LEU:HD23	1:A:327:LEU:HD12	1.87	0.55
1:B:555:GLN:NE2	1:B:580:PRO:HG2	2.22	0.54
1:A:489:GLU:HB3	1:A:621:HIS:HE2	1.71	0.54
1:A:322:TRP:H	1:A:328:GLN:HE22	1.55	0.54
1:A:455:ALA:HB2	1:A:468:ILE:HG13	1.89	0.54
1:A:210:GLN:NE2	1:A:380:LYS:HE3	2.23	0.54
1:A:84:GLN:CD	4:A:2039:HOH:O	2.46	0.53
1:A:85:LYS:HD3	4:A:1985:HOH:O	2.08	0.53
1:A:539:MET:HE1	4:A:2538:HOH:O	2.07	0.53
1:A:44:GLN:HB2	4:A:2200:HOH:O	2.08	0.53
1:B:457:LEU:HD21	1:B:466:ASP:HB2	1.91	0.52
1:A:296:SER:HB2	1:A:298:ASP:OD1	2.10	0.52
1:B:568:PRO:HD2	1:B:592:TYR:CZ	2.45	0.52
1:B:322:TRP:H	1:B:328:GLN:HE22	1.56	0.52
1:A:176:ILE:HA	1:A:326:THR:HG21	1.91	0.52
1:A:255:HIS:CD2	1:A:257:ASN:H	2.27	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:28:HIS:HA	1:A:156:GLU:OE1	2.10	0.52
1:A:320:ASN:ND2	4:A:2047:HOH:O	2.43	0.51
1:B:225:THR:H	1:B:228:GLN:NE2	2.08	0.51
1:B:244:ASN:HB2	4:B:2330:HOH:O	2.08	0.51
1:A:529:GLU:HG2	4:A:1962:HOH:O	2.09	0.51
1:A:84:GLN:NE2	4:A:2015:HOH:O	2.44	0.51
1:A:84:GLN:NE2	4:A:2039:HOH:O	2.43	0.51
1:A:58:VAL:HG21	4:A:2017:HOH:O	2.10	0.51
1:A:457:LEU:HD12	1:A:466:ASP:HB2	1.92	0.51
1:A:32:GLN:NE2	1:A:170:MET:H	2.08	0.51
1:A:548:HIS:HE1	1:A:572:ASP:OD1	1.94	0.50
1:B:301:ARG:HD3	4:B:2612:HOH:O	2.11	0.50
1:B:176:ILE:HA	1:B:326:THR:HG21	1.93	0.50
3:A:1801:152:O3	3:A:1801:152:H5B3	2.12	0.49
1:A:368:VAL:HG22	4:A:2020:HOH:O	2.12	0.49
1:A:328:GLN:O	1:A:339:MET:HB2	2.13	0.49
1:B:255:HIS:CD2	1:B:257:ASN:H	2.27	0.49
1:A:218:HIS:HA	4:A:2396:HOH:O	2.13	0.49
1:A:277:ASN:OD1	1:A:281:LYS:HE2	2.12	0.49
1:A:41:PRO:HG3	1:A:513:GLY:O	2.11	0.49
1:B:32:GLN:NE2	1:B:170:MET:H	2.08	0.49
3:B:1901:152:O3	3:B:1901:152:H5B3	2.12	0.49
1:B:242:GLN:HB2	4:B:2583:HOH:O	2.13	0.49
1:B:310:HIS:HE1	4:B:2252:HOH:O	1.95	0.48
1:A:456:SER:HB2	2:A:1802:COA:H71	1.94	0.48
1:B:349:PRO:HA	1:B:352:VAL:HG22	1.95	0.48
1:B:122:VAL:HG22	1:B:564:MET:HG2	1.94	0.48
1:B:123:ILE:HG13	1:B:565:PHE:HE2	1.78	0.48
1:B:99:SER:HB3	4:B:2538:HOH:O	2.12	0.48
1:A:367:LEU:HD12	4:A:2196:HOH:O	2.14	0.48
1:A:296:SER:OG	1:A:299:VAL:HG22	2.14	0.48
1:B:70:VAL:O	1:B:74:GLN:HG2	2.14	0.48
1:B:104:LYS:HG2	1:B:109:GLN:CD	2.34	0.48
1:A:163:LEU:HB2	1:A:168:LEU:HD21	1.95	0.48
1:A:352:VAL:HG11	1:A:556:VAL:HG12	1.96	0.48
1:A:405:GLN:HG3	4:A:2388:HOH:O	2.13	0.48
1:A:122:VAL:HB	1:A:339:MET:HG2	1.96	0.47
1:A:426:LYS:CB	1:A:623:ARG:HH12	2.23	0.47
1:B:496:LEU:HD23	1:B:496:LEU:C	2.34	0.47
1:B:452:TYR:HB2	1:B:549:PHE:CD1	2.49	0.47
1:A:416:HIS:CG	1:A:612:LEU:HD21	2.50	0.47
1:B:198:ARG:HG3	1:B:198:ARG:O	2.14	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:321:ARG:NH1	4:A:2217:HOH:O	2.48	0.47
1:B:333:GLU:HB2	4:B:2338:HOH:O	2.13	0.47
1:A:166:GLN:HB2	4:A:2249:HOH:O	2.15	0.47
1:A:132:LEU:HD22	1:A:132:LEU:N	2.28	0.47
1:A:426:LYS:CG	1:A:623:ARG:HH12	2.27	0.46
1:A:183:PRO:HG3	1:A:259:TRP:NE1	2.30	0.46
1:B:129:PHE:CZ	1:B:335:GLY:HA2	2.50	0.46
1:B:274:GLU:HG3	4:B:2378:HOH:O	2.14	0.46
1:A:507:THR:HG23	2:A:1802:COA:H143	1.97	0.46
1:A:439:LEU:HA	1:A:479:VAL:HG12	1.98	0.46
1:B:410:MET:HG2	1:B:601:ALA:HA	1.99	0.46
1:A:489:GLU:HB3	1:A:621:HIS:NE2	2.31	0.45
1:B:202:HIS:HD2	4:B:1914:HOH:O	1.99	0.45
1:A:255:HIS:HE1	4:A:1986:HOH:O	1.98	0.45
1:B:197:LYS:O	1:B:199:PRO:HD3	2.17	0.45
1:B:512:ARG:HG3	1:B:512:ARG:HH11	1.81	0.45
1:A:124:LEU:HG	1:A:338:GLY:HA2	1.98	0.45
1:B:476:LEU:HA	1:B:479:VAL:HG22	1.98	0.45
1:A:458:ARG:HD3	4:A:1935:HOH:O	2.17	0.45
1:A:204:THR:HG23	1:A:282:SER:HB3	1.97	0.45
1:B:124:LEU:HG	1:B:338:GLY:HA2	1.98	0.45
1:A:366:GLU:CA	1:A:369:ARG:HH21	2.29	0.45
1:B:548:HIS:HD2	4:B:2258:HOH:O	2.00	0.45
1:A:354:LEU:O	1:A:358:VAL:HG23	2.17	0.45
1:B:207:HIS:CD2	1:B:240:SER:OG	2.64	0.44
1:B:420:ASP:OD1	1:B:421:PHE:N	2.50	0.44
1:A:144:GLU:HG3	4:A:1953:HOH:O	2.17	0.44
1:B:50:TYR:CZ	1:B:54:LEU:HD11	2.53	0.44
1:A:160:VAL:HG13	4:A:2018:HOH:O	2.18	0.44
1:B:182:VAL:HA	1:B:183:PRO:HD3	1.88	0.44
1:A:347:GLU:HB3	1:A:349:PRO:HD2	2.00	0.44
1:A:94:MET:HE2	4:A:2144:HOH:O	2.17	0.44
1:B:548:HIS:HE1	1:B:572:ASP:OD1	2.01	0.44
1:B:514:GLU:HG2	4:B:2375:HOH:O	2.17	0.44
1:A:512:ARG:NH1	1:A:512:ARG:HG3	2.31	0.44
1:B:320:ASN:C	1:B:321:ARG:HG2	2.39	0.44
1:A:27:SER:OG	1:A:32:GLN:NE2	2.51	0.43
1:A:81:GLU:HG3	4:A:2026:HOH:O	2.18	0.43
1:B:574:TYR:CE1	1:B:603:MET:HE3	2.52	0.43
1:A:354:LEU:C	1:A:354:LEU:HD23	2.38	0.43
1:B:54:LEU:O	1:B:58:VAL:HG22	2.18	0.43
1:A:37:ARG:HG3	1:A:95:GLU:O	2.17	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:310:HIS:HD2	1:B:312:GLY:N	2.06	0.43
1:A:47:LEU:HD13	1:A:70:VAL:HG13	2.00	0.43
1:A:255:HIS:HD2	1:A:257:ASN:H	1.66	0.43
1:B:488:PRO:HD2	1:B:491:GLN:CD	2.38	0.43
1:A:204:THR:HG21	1:A:279:ILE:HA	2.01	0.43
1:A:115:VAL:HG12	1:A:116:ILE:HG13	2.00	0.43
1:A:578:TYR:HB3	1:A:587:PHE:CD1	2.54	0.43
1:A:57:ILE:HG13	1:A:58:VAL:HG13	1.99	0.43
1:B:129:PHE:HB2	1:B:134:GLY:HA3	2.00	0.43
1:A:123:ILE:HG13	1:A:565:PHE:CE2	2.54	0.43
1:A:461:HIS:N	4:A:2045:HOH:O	2.49	0.43
1:A:147:LEU:HD11	1:A:216:VAL:HB	2.01	0.43
1:A:199:PRO:HA	1:A:200:PRO:HD3	1.93	0.42
1:B:197:LYS:C	1:B:199:PRO:HD3	2.40	0.42
1:A:47:LEU:HD12	1:A:74:GLN:NE2	2.34	0.42
1:A:431:ALA:HB3	1:A:500:VAL:HG13	2.00	0.42
1:A:220:ASP:OD1	1:A:222:THR:HG23	2.19	0.42
1:A:523:LEU:HD21	4:A:2015:HOH:O	2.18	0.42
1:B:421:PHE:CZ	1:B:615:ARG:HG3	2.55	0.42
1:B:597:GLU:HG2	4:B:2522:HOH:O	2.18	0.42
1:A:132:LEU:CD2	1:A:132:LEU:H	2.30	0.42
1:B:489:GLU:HB3	1:B:621:HIS:NE2	2.35	0.42
1:A:202:HIS:HD2	4:A:2069:HOH:O	2.02	0.42
1:B:364:LYS:HA	1:B:365:PRO:HD3	1.94	0.42
1:B:556:VAL:HG21	1:B:564:MET:HE2	2.02	0.42
1:A:321:ARG:NH1	1:A:321:ARG:HG2	2.32	0.41
1:A:426:LYS:HB2	1:A:623:ARG:NH1	2.32	0.41
1:A:382:ARG:HG3	4:A:2031:HOH:O	2.19	0.41
1:B:621:HIS:HB3	4:B:2282:HOH:O	2.19	0.41
1:A:234:GLU:HB3	4:A:2140:HOH:O	2.19	0.41
1:B:416:HIS:CG	1:B:612:LEU:HD21	2.56	0.41
1:B:455:ALA:HB2	1:B:468:ILE:HG13	2.02	0.41
1:B:192:ASN:C	1:B:192:ASN:HD22	2.22	0.41
1:B:144:GLU:HG3	4:B:2087:HOH:O	2.20	0.41
1:A:556:VAL:HG22	2:A:1802:COA:H32	2.02	0.41
1:A:85:LYS:HE3	4:A:2026:HOH:O	2.20	0.41
1:A:315:LYS:HE3	4:A:2022:HOH:O	2.20	0.41
1:B:524:LYS:HD3	4:B:2322:HOH:O	2.21	0.41
1:A:457:LEU:HA	4:A:2176:HOH:O	2.21	0.41
1:B:163:LEU:CB	1:B:168:LEU:HD21	2.51	0.41
1:A:548:HIS:HD2	4:A:2181:HOH:O	2.03	0.41
1:B:202:HIS:HE1	1:B:278:SER:O	2.04	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:66:THR:O	1:A:70:VAL:HG23	2.21	0.41
1:A:410:MET:HG2	1:A:601:ALA:HA	2.02	0.41
1:A:374:PRO:HA	4:A:2004:HOH:O	2.21	0.41
1:B:163:LEU:HB2	1:B:168:LEU:HD11	2.02	0.40
1:B:512:ARG:NH1	1:B:512:ARG:HG3	2.37	0.40
1:B:231:VAL:HG21	1:B:372:MET:SD	2.61	0.40
1:B:489:GLU:HB3	1:B:621:HIS:HE2	1.86	0.40
1:A:555:GLN:NE2	4:A:1885:HOH:O	2.53	0.40
1:A:247:PRO:O	1:A:250:ILE:HG22	2.22	0.40
1:B:225:THR:H	1:B:228:GLN:HE21	1.70	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	597/599 (100%)	578 (97%)	18 (3%)	1 (0%)	56	44
1	B	597/599 (100%)	580 (97%)	15 (2%)	2 (0%)	50	37
All	All	1194/1198 (100%)	1158 (97%)	33 (3%)	3 (0%)	50	37

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	164	GLY
1	B	372	MET
1	B	624	ALA

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	527/527 (100%)	516 (98%)	11 (2%)	66	59
1	B	527/527 (100%)	517 (98%)	10 (2%)	69	63
All	All	1054/1054 (100%)	1033 (98%)	21 (2%)	68	61

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	137	ARG
1	A	141	LYS
1	A	163	LEU
1	A	192	ASN
1	A	343	HIS
1	A	347	GLU
1	A	372	MET
1	A	457	LEU
1	A	506	TYR
1	A	517	ASP
1	A	534	MET
1	B	33	ASP
1	B	61	GLU
1	B	137	ARG
1	B	192	ASN
1	B	206	VAL
1	B	308	MET
1	B	347	GLU
1	B	506	TYR
1	B	597	GLU
1	B	625	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (48) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	74	GLN
1	A	84	GLN
1	A	109	GLN
1	A	135	GLN
1	A	155	ASN
1	A	192	ASN
1	A	202	HIS
1	A	207	HIS

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Mol	Chain	Res	Type
1	A	210	GLN
1	A	228	GLN
1	A	255	HIS
1	A	307	GLN
1	A	310	HIS
1	A	328	GLN
1	A	447	GLN
1	A	461	HIS
1	A	503	HIS
1	A	548	HIS
1	A	586	ASN
1	A	619	GLN
1	A	620	ASN
1	B	32	GLN
1	B	65	HIS
1	B	74	GLN
1	B	84	GLN
1	B	133	GLN
1	B	135	GLN
1	B	155	ASN
1	B	187	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	405	GLN
1	B	461	HIS
1	B	490	GLN
1	B	501	GLN
1	B	503	HIS
1	B	548	HIS
1	B	586	ASN
1	B	619	GLN
1	B	620	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

4 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$
3	152	A	1801	-	10,10,10	1.14	1 (10%)	14,14,14	0.86	1 (7%)
2	COA	A	1802	-	50,50,50	1.16	3 (6%)	75,75,75	1.61	7 (9%)
3	152	B	1901	-	10,10,10	1.05	0	14,14,14	1.06	1 (7%)
2	COA	B	1902	-	50,50,50	1.10	3 (6%)	75,75,75	1.65	7 (9%)

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
3	152	A	1801	-	-	0/9/9/9	0/0/0/0
2	COA	A	1802	-	-	0/48/64/64	0/1/3/3
3	152	B	1901	-	-	0/9/9/9	0/0/0/0
2	COA	B	1902	-	-	0/48/64/64	0/1/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1802	COA	O4B-C1B	4.07	1.47	1.41
2	B	1902	COA	O4B-C1B	3.61	1.46	1.41
2	A	1802	COA	P3B-O7A	3.00	1.61	1.51
2	A	1802	COA	C4A-N9A	-2.98	1.33	1.37
2	B	1902	COA	C4A-N9A	-2.84	1.33	1.37
2	B	1902	COA	P3B-O7A	2.83	1.60	1.51
3	A	1801	152	C4-N5	-2.19	1.47	1.51

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1902	COA	N3A-C2A-N1A	-9.89	120.44	128.71
2	A	1802	COA	N3A-C2A-N1A	-9.68	120.61	128.71
2	B	1902	COA	O4B-C1B-N9A	4.52	112.64	108.44
2	B	1902	COA	N3A-C4A-N9A	4.07	132.78	125.43
2	A	1802	COA	N3A-C4A-N9A	4.06	132.77	125.43
2	A	1802	COA	O4B-C1B-N9A	3.85	112.02	108.44
2	B	1902	COA	P2A-O3A-P1A	-3.08	122.64	131.68
3	B	1901	152	C3-C4-N5	-3.05	111.01	117.21
2	A	1802	COA	P2A-O3A-P1A	-3.01	122.84	131.68
2	A	1802	COA	O9A-P3B-O8A	2.90	118.91	107.61
2	B	1902	COA	O9A-P3B-O8A	2.87	118.78	107.61
2	B	1902	COA	C5A-C4A-N3A	-2.69	119.84	125.70
2	A	1802	COA	C5A-C4A-N3A	-2.69	119.85	125.70
3	A	1801	152	C3-C4-N5	-2.61	111.91	117.21
2	A	1802	COA	C2A-N3A-C4A	2.37	120.75	114.01
2	B	1902	COA	C2A-N3A-C4A	2.36	120.72	114.01

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	599/599 (100%)	0.25	35 (5%) 22 22	11, 22, 47, 75	0
1	B	599/599 (100%)	-0.10	16 (2%) 52 53	7, 15, 36, 75	0
All	All	1198/1198 (100%)	0.08	51 (4%) 35 34	7, 19, 43, 75	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	368	VAL	12.2
1	B	367	LEU	10.3
1	A	368	VAL	8.9
1	A	367	LEU	8.4
1	B	371	PRO	7.9
1	B	369	ARG	7.2
1	B	370	SER	7.1
1	A	625	LYS	7.1
1	A	624	ALA	7.0
1	B	624	ALA	6.3
1	A	370	SER	6.2
1	A	371	PRO	5.9
1	B	625	LYS	5.2
1	A	374	PRO	5.1
1	A	369	ARG	4.7
1	B	372	MET	4.4
1	A	163	LEU	4.3
1	A	30	ALA	3.9
1	B	244	ASN	3.8
1	A	373	VAL	3.8
1	A	623	ARG	3.8
1	B	294	ARG	3.8
1	B	366	GLU	3.8
1	B	623	ARG	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	77	GLY	3.7
1	A	132	LEU	3.3
1	A	366	GLU	3.3
1	A	198	ARG	3.2
1	B	29	MET	3.1
1	A	29	MET	3.1
1	A	164	GLY	3.0
1	A	27	SER	2.9
1	A	372	MET	2.8
1	A	197	LYS	2.8
1	A	160	VAL	2.7
1	A	61	GLU	2.6
1	B	622	PRO	2.6
1	A	28	HIS	2.6
1	A	165	GLY	2.6
1	B	198	ARG	2.6
1	A	68	GLN	2.5
1	A	227	ASP	2.5
1	A	220	ASP	2.5
1	A	294	ARG	2.4
1	A	296	SER	2.3
1	A	297	ASP	2.2
1	A	293	PRO	2.2
1	A	219	SER	2.2
1	A	75	THR	2.1
1	B	30	ALA	2.1
1	A	60	GLU	2.1

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	152	A	1801	11/11	0.12	1.55	13,15,17,18	0
2	COA	B	1902	48/48	0.11	0.20	14,24,31,32	0
2	COA	A	1802	48/48	0.10	-0.21	19,24,33,35	0
3	152	B	1901	11/11	0.09	-0.46	8,9,11,13	0

6.5 Other polymers ⓘ

There are no such residues in this entry.