



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:48 AM GMT

PDB ID : 2YEV
Title : Structure of caa3-type cytochrome oxidase
Authors : Lyons, J.A.; Aragao, D.; Soulimane, T.; Caffrey, M.
Deposited on : 2011-03-31
Resolution : 2.36 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

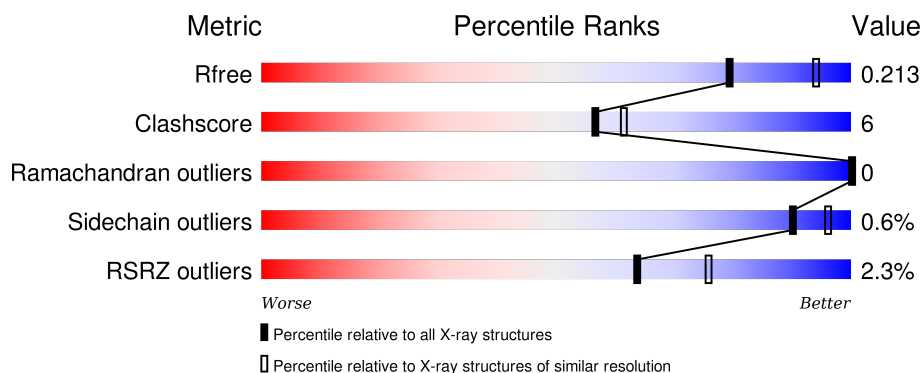
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.36 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	1352 (2.38-2.34)
Clashscore	102246	1456 (2.38-2.34)
Ramachandran outliers	100387	1435 (2.38-2.34)
Sidechain outliers	100360	1436 (2.38-2.34)
RSRZ outliers	91569	1358 (2.38-2.34)

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	791	<div> <div>2%</div> <div>85%</div> <div>13%</div> <div>•</div> </div>
1	D	791	<div> <div>2%</div> <div>85%</div> <div>14%</div> <div>•</div> </div>
2	B	337	<div> <div>•</div> <div>84%</div> <div>11%</div> <div>5%</div> </div>
2	E	337	<div> <div>4%</div> <div>83%</div> <div>11%</div> <div>•</div> <div>5%</div> </div>
3	C	66	<div> <div>2%</div> <div>88%</div> <div>9%</div> <div>•</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	66	

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
12	7E9	B	701	-	-	-	X
7	4AG	A	1200	-	-	-	X
8	7E8	A	1300	X	-	-	X
8	7E8	A	1301	-	-	-	X
9	MG	A	1801	-	-	-	X

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 19587 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 CYTOCHROME C OXIDASE POLYPEPTIDE I+III.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	780	Total	C	N	O	S	0	1	0
			6281	4244	992	1022	23			
1	D	780	Total	C	N	O	S	0	0	0
			6276	4241	992	1020	23			

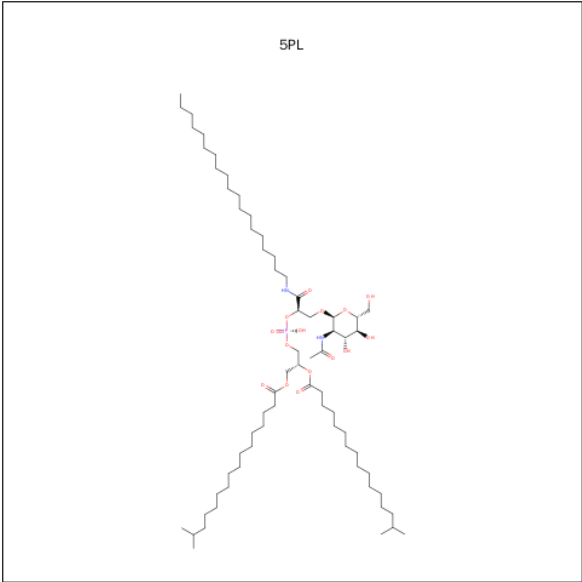
- Molecule 2 is a protein called CYTOCHROME C OXIDASE SUBUNIT 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	319	Total	C	N	O	S	0	0	0
			2507	1637	422	439	9			
2	E	319	Total	C	N	O	S	0	0	0
			2507	1637	422	439	9			

- Molecule 3 is a protein called CAA3-TYPE CYTOCHROME OXIDASE SUBUNIT IV.

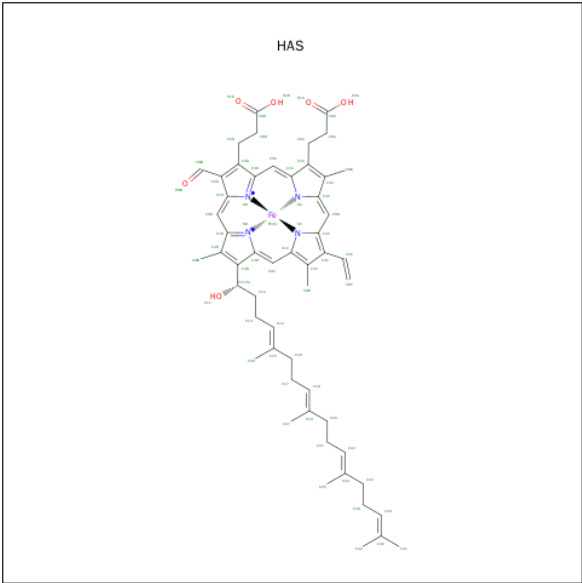
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	64	Total	C	N	O	S	0	0	0
			502	344	76	79	3			
3	F	63	Total	C	N	O	S	0	0	0
			492	338	73	78	3			

- Molecule 4 is (1R,4S,6R)-6-({[2-(ACETYLAMINO)-2-DEOXY-ALPHA-D-GLUCOPYRANOSYL]OXY}METHYL)-4-HYDROXY-1-{{[(15-METHYLHEXADECANOYL)OXY]METHYL}-4-OXIDO-7-OXO-3,5-DIOXA-8-AZA-4-PHOSPHAHEPTACOS-1-YL 15-METHYLHEXADECANOATE (three-letter code: 5PL) (formula: C₆₇H₁₂₉N₂O₁₅P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			85	67	2	15	1		
4	D	1	Total	C	N	O	P	0	0
			85	67	2	15	1		

- Molecule 5 is HEME-AS (three-letter code: HAS) (formula: $C_{54}H_{64}FeN_4O_6$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	Fe	N	O	0	0
			65	54	1	4	6		
5	A	1	Total	C	Fe	N	O	0	0
			65	54	1	4	6		

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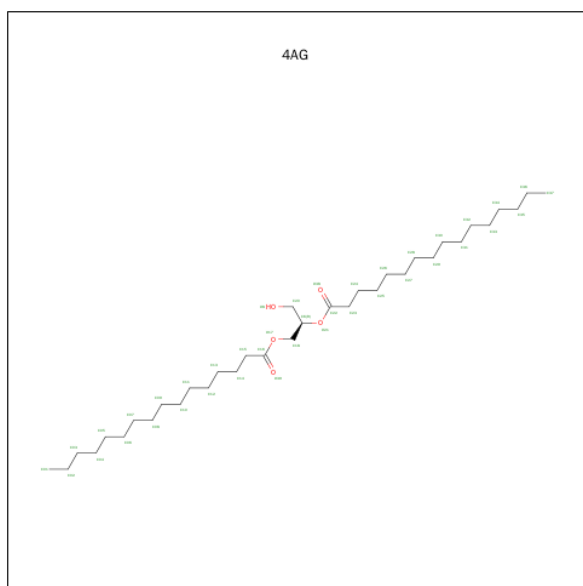
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	D	1	Total	C	Fe	N	O	
			65	54	1	4	6	
5	D	1	Total	C	Fe	N	O	
			65	54	1	4	6	

- Molecule 6 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

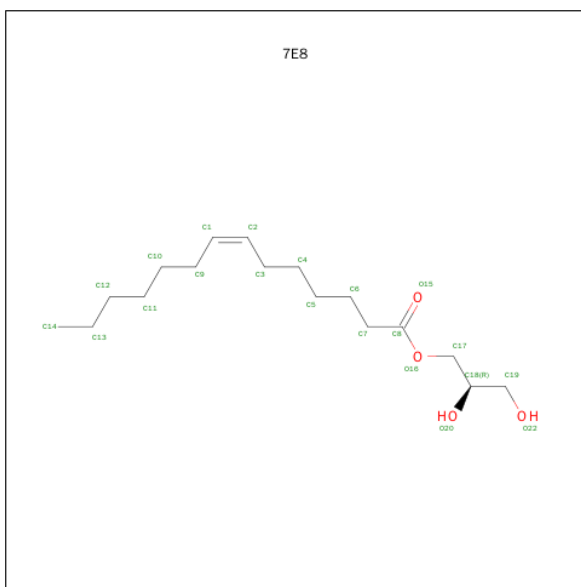
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Cu		
			1	1	0	0
6	D	1	Total	Cu		
			1	1	0	0

- Molecule 7 is (2R)-3-HYDROXYPROPANE-1,2-DIYL DIHEXADECANOATE (three-letter code: 4AG) (formula: C₃₅H₆₈O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O		
			40	35	5	0	0

- Molecule 8 is (2R)-2,3-DIHYDROXYPROPYL (7Z)-TETRADEC-7-ENOATE (three-letter code: 7E8) (formula: C₁₇H₃₂O₄).

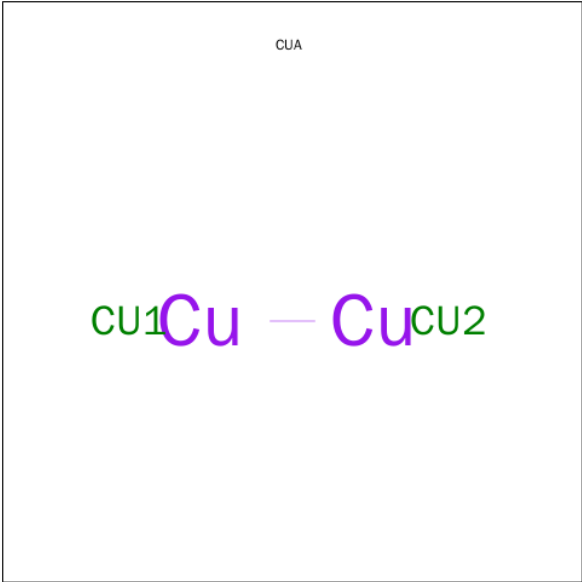


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C O 21 17 4	0	0
8	A	1	Total C O 21 17 4	0	0

- Molecule 9 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

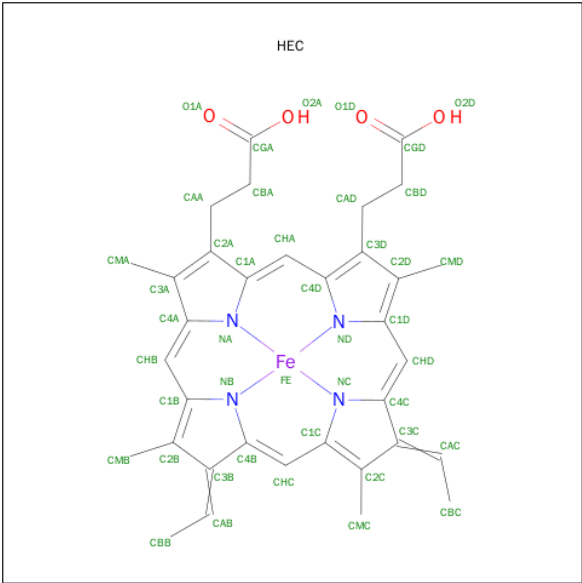
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	1	Total Mg 1 1	0	0
9	D	1	Total Mg 1 1	0	0

- Molecule 10 is DINUCLEAR COPPER ION (three-letter code: CUA) (formula: Cu₂).



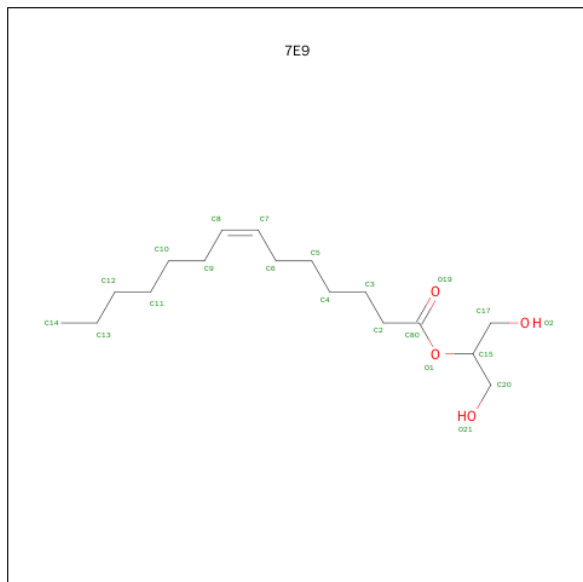
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	B	1	Total 2	Cu 2	0	0
10	E	1	Total 2	Cu 2	0	0

- Molecule 11 is HEME C (three-letter code: HEC) (formula: C₃₄H₃₄FeN₄O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
11	B	1	Total	C	Fe	N	O	0
			43	34	1	4	4	
11	E	1	Total	C	Fe	N	O	0
			43	34	1	4	4	

- Molecule 12 is 1,3-DIHYDROXYPROPAN-2-YL (Z)-TETRADEC-7-ENOATE (three-letter code: 7E9) (formula: C₁₇H₃₂O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	B	1	Total	C	O	0	0
			21	17	4		

- Molecule 13 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	D	1	Total	Cl	0	0
			1	1		
13	C	1	Total	Cl	0	0
			1	1		
13	F	1	Total	Cl	0	0
			1	1		

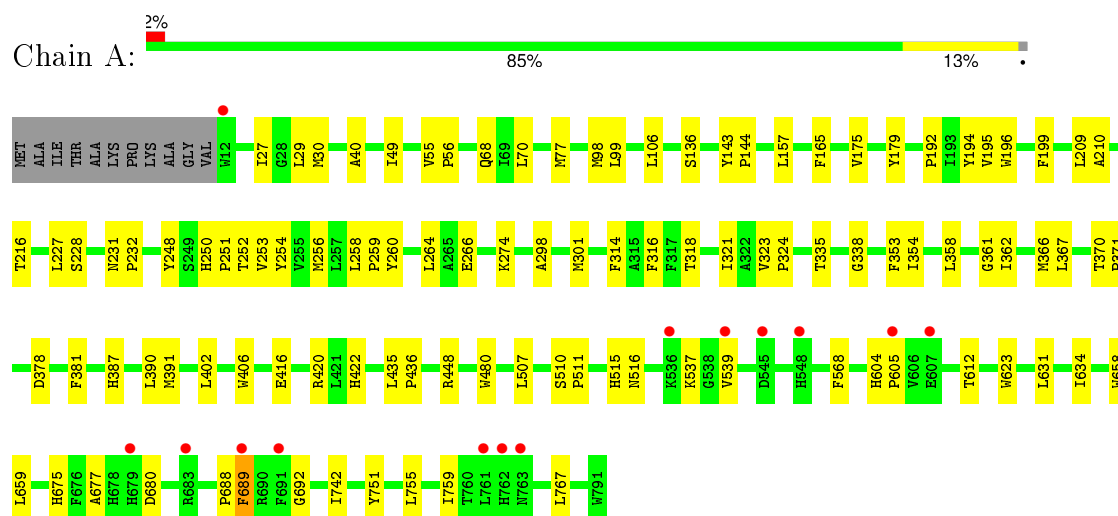
- Molecule 14 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
14	A	157	Total	O	0	0
			157	157		
14	B	90	Total	O	0	0
			90	90		
14	D	111	Total	O	0	0
			111	111		
14	E	34	Total	O	0	0
			34	34		

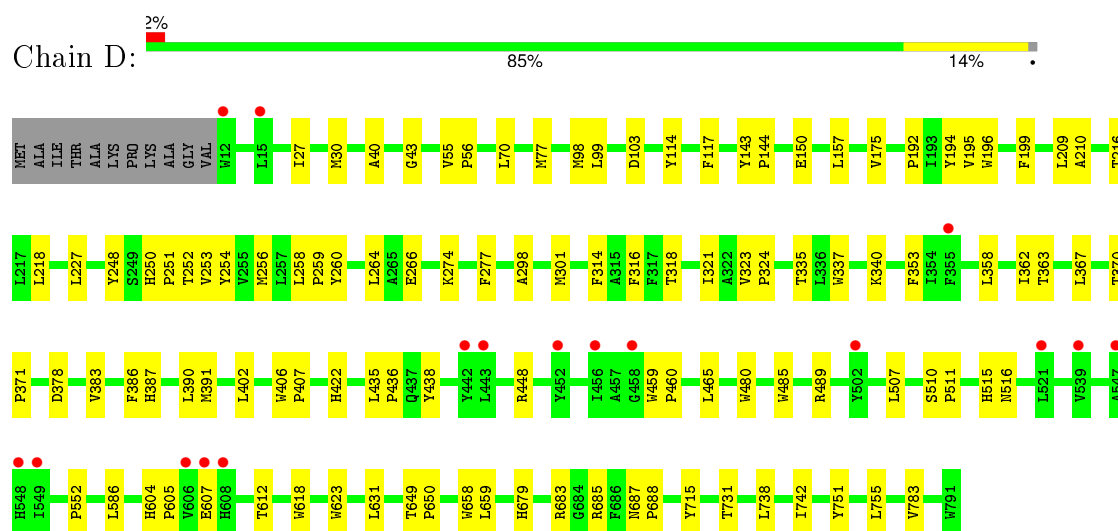
3 Residue-property plots [i](#)

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.

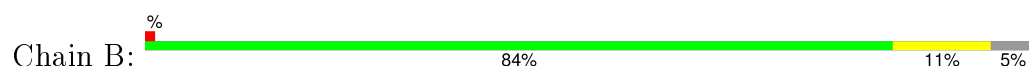
• Molecule 1: CYTOCHROME C OXIDASE POLYPEPTIDE I+III

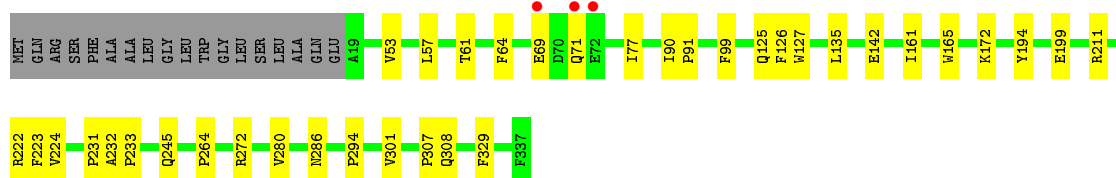


• Molecule 1: CYTOCHROME C OXIDASE POLYPEPTIDE I+III

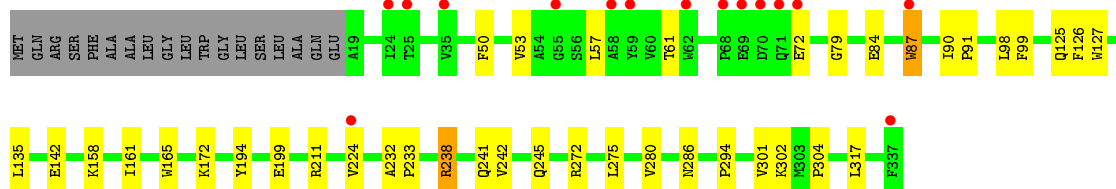
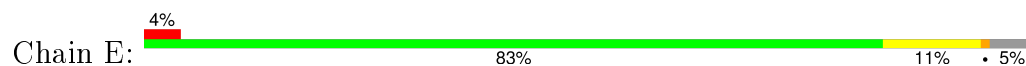


• Molecule 2: CYTOCHROME C OXIDASE SUBUNIT 2

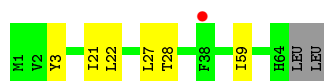
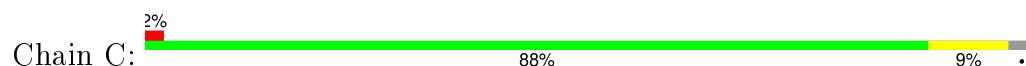




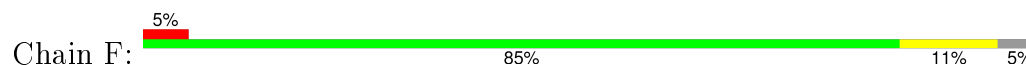
• Molecule 2: CYTOCHROME C OXIDASE SUBUNIT 2



• Molecule 3: CAA3-TYPE CYTOCHROME OXIDASE SUBUNIT IV



• Molecule 3: CAA3-TYPE CYTOCHROME OXIDASE SUBUNIT IV



4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	127.25Å 76.03Å 300.27Å 90.00° 92.21° 90.00°	Depositor
Resolution (Å)	77.18 – 2.36 115.49 – 2.36	Depositor EDS
% Data completeness (in resolution range)	100.0 (77.18-2.36) 100.0 (115.49-2.36)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.37Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.6.4_486)	Depositor
R, R_{free}	0.171 , 0.218 0.168 , 0.213	Depositor DCC
R_{free} test set	5970 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	50.2	Xtriage
Anisotropy	0.263	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 56.5	EDS
Estimated twinning fraction	0.017 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 118259 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	19587	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.82% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 5PL, 7E8, CL, MG, HAS, 4AG, CUA, HEC, 7E9, FME, CU

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	0.32	0/6523	0.44	0/8921
1	D	0.31	0/6515	0.44	0/8910
2	B	0.32	0/2586	0.46	0/3523
2	E	0.31	0/2586	0.46	0/3523
3	C	0.26	0/502	0.38	0/683
3	F	0.25	0/491	0.39	0/668
All	All	0.31	0/19203	0.44	0/26228

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 [i](#)

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	6281	0	6210	82	0
1	D	6276	0	6206	97	0
2	B	2507	0	2473	26	0
2	E	2507	0	2473	27	0
3	C	502	0	545	4	0
3	F	492	0	538	6	0
4	A	85	0	128	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	85	0	128	4	0
5	A	130	0	124	18	0
5	D	130	0	124	23	0
6	A	1	0	0	0	0
6	D	1	0	0	0	0
7	A	40	0	68	1	0
8	A	42	0	64	2	0
9	A	1	0	0	0	0
9	D	1	0	0	0	0
10	B	2	0	0	0	0
10	E	2	0	0	0	0
11	B	43	0	30	3	0
11	E	43	0	30	3	0
12	B	21	0	32	0	0
13	C	1	0	0	0	0
13	D	1	0	0	0	0
13	F	1	0	0	0	0
14	A	157	0	0	4	0
14	B	90	0	0	1	0
14	D	111	0	0	0	0
14	E	34	0	0	0	0
All	All	19587	0	19173	232	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 6.

All (232) 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:77:MET:HB3	5:A:1015:HAS:CAC	1.99	0.93
1:D:77:MET:HB3	5:D:1015:HAS:CAC	2.00	0.91
1:D:367:LEU:CB	5:D:1016:HAS:HMD	2.15	0.77
1:D:70:LEU:HD13	5:D:1015:HAS:HBD1	1.66	0.77
1:A:77:MET:HB3	5:A:1015:HAS:HAC	1.68	0.73
3:C:3:TYR:HB3	3:C:59:ILE:HD11	1.72	0.71
1:D:367:LEU:HB2	5:D:1016:HAS:HMD	1.72	0.70
3:F:3:TYR:HB3	3:F:59:ILE:HD11	1.73	0.69
2:B:245:GLN:NE2	1:D:150:GLU:HG3	2.08	0.69
1:D:358:LEU:O	1:D:362:ILE:HG12	1.94	0.68
1:A:367:LEU:HB3	5:A:1016:HAS:HMD	1.74	0.68
1:A:367:LEU:CB	5:A:1016:HAS:HMD	2.25	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:338:GLY:HA2	2:B:64:PHE:CE1	2.31	0.66
1:D:277:PHE:CE2	2:E:79:GLY:HA2	2.31	0.65
1:A:358:LEU:O	1:A:362:ILE:HG12	1.96	0.64
2:B:307:PRO:HG2	2:E:304:PRO:HG3	1.80	0.63
2:B:245:GLN:HE22	1:D:150:GLU:HG3	1.63	0.63
1:D:70:LEU:HD21	1:D:448:ARG:HG2	1.81	0.63
1:A:40:ALA:HB1	5:A:1015:HAS:H211	1.81	0.63
1:A:70:LEU:HD21	1:A:448:ARG:HG2	1.82	0.62
1:D:298:ALA:HB3	1:D:314:PHE:CG	2.35	0.62
14:A:2082:HOH:O	2:B:77:ILE:HD11	2.01	0.61
5:D:1016:HAS:HMC1	5:D:1016:HAS:HBC1	1.82	0.60
5:D:1016:HAS:H323	2:E:98:LEU:HD11	1.84	0.60
1:A:323:VAL:HB	1:A:324:PRO:HD3	1.84	0.60
2:B:127:TRP:CD1	11:B:587:HEC:HAD1	2.38	0.59
1:D:604:HIS:HB3	1:D:605:PRO:HD2	1.85	0.59
1:A:298:ALA:HB3	1:A:314:PHE:CG	2.38	0.58
2:B:135:LEU:HD13	2:B:224:VAL:HG11	1.84	0.58
2:E:127:TRP:CD1	11:E:587:HEC:HAD1	2.38	0.58
2:E:238:ARG:O	2:E:238:ARG:HD2	2.03	0.58
1:A:568:PHE:CZ	4:A:900:5PL:H202	2.39	0.58
1:D:323:VAL:HB	1:D:324:PRO:HD3	1.86	0.57
1:D:783:VAL:HG23	3:F:43:LEU:HD11	1.87	0.57
1:D:55:VAL:HB	1:D:56:PRO:HD2	1.88	0.56
1:D:77:MET:HB3	5:D:1015:HAS:HAC	1.85	0.55
1:D:367:LEU:HD13	5:D:1016:HAS:HBD2	1.87	0.55
1:D:298:ALA:HB3	1:D:314:PHE:CD2	2.42	0.55
1:A:165:PHE:HE1	4:A:900:5PL:HBK2	1.72	0.55
1:D:143:TYR:CG	1:D:144:PRO:HA	2.42	0.54
1:D:250:HIS:HB3	1:D:251:PRO:HD3	1.89	0.54
1:D:618:TRP:HH2	3:F:27:LEU:HD22	1.73	0.54
1:D:362:ILE:HG23	5:D:1016:HAS:H313	1.90	0.54
1:A:258:LEU:HB2	1:A:259:PRO:HD3	1.89	0.54
1:A:157:LEU:HD22	1:A:216:THR:HG23	1.89	0.54
2:E:135:LEU:HD22	2:E:224:VAL:HG11	1.90	0.53
1:A:537:LYS:HB2	1:A:539:VAL:HG23	1.90	0.53
1:D:258:LEU:HB2	1:D:259:PRO:HD3	1.91	0.52
2:B:161:ILE:HG22	2:B:199:GLU:HG2	1.91	0.52
1:D:783:VAL:HG23	3:F:43:LEU:CD1	2.39	0.52
1:A:49:ILE:CG2	5:A:1015:HAS:HMD	2.40	0.52
1:A:143:TYR:CG	1:A:144:PRO:HA	2.45	0.52
1:D:256:MET:O	1:D:260:TYR:HD2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:157:LEU:HD22	1:D:216:THR:HG23	1.92	0.51
2:E:135:LEU:HD13	2:E:224:VAL:HG11	1.93	0.51
2:E:57:LEU:O	2:E:61:THR:HG23	2.10	0.51
1:A:366:MET:SD	7:A:1200:4AG:H352	2.50	0.51
2:B:57:LEU:O	2:B:61:THR:HG23	2.11	0.50
1:A:250:HIS:HB3	1:A:251:PRO:HD3	1.93	0.50
1:D:679:HIS:CE1	1:D:683:ARG:CZ	2.94	0.50
1:A:298:ALA:HB3	1:A:314:PHE:CD2	2.46	0.50
1:A:55:VAL:HB	1:A:56:PRO:HD2	1.93	0.50
1:A:209:LEU:HA	1:A:248:TYR:CD1	2.46	0.50
1:D:227:LEU:C	1:D:227:LEU:HD23	2.32	0.50
1:D:40:ALA:HB1	5:D:1015:HAS:H253	1.93	0.50
1:D:435:LEU:HB2	1:D:436:PRO:HD3	1.93	0.50
11:B:587:HEC:HBB3	11:B:587:HEC:HMB1	1.94	0.50
1:A:106:LEU:HD13	4:A:900:5PL:HBV1	1.93	0.50
1:A:227:LEU:C	1:A:227:LEU:HD23	2.33	0.49
2:E:142:GLU:HG2	2:E:211:ARG:HB2	1.94	0.49
1:A:256:MET:O	1:A:260:TYR:HD2	1.96	0.49
1:D:387:HIS:O	1:D:391:MET:HB3	2.12	0.49
1:A:612:THR:HA	3:C:28:THR:HG22	1.95	0.49
1:A:98:MET:HB3	1:A:192:PRO:HG2	1.95	0.49
1:D:277:PHE:CD2	2:E:79:GLY:HA2	2.48	0.49
1:A:658:TRP:CH2	1:A:659:LEU:HD13	2.46	0.49
1:D:209:LEU:HA	1:D:248:TYR:CD1	2.48	0.49
2:E:161:ILE:HG22	2:E:199:GLU:HG2	1.95	0.49
1:A:298:ALA:HA	1:A:301:MET:HE2	1.95	0.48
1:A:677:ALA:HB2	1:A:692:GLY:HA3	1.95	0.48
2:E:317:LEU:HD21	11:E:587:HEC:HMB1	1.94	0.48
1:D:658:TRP:CH2	1:D:659:LEU:HD13	2.48	0.48
1:A:70:LEU:HD13	5:A:1015:HAS:HBD1	1.95	0.48
1:D:607:GLU:O	1:D:607:GLU:HG3	2.13	0.48
1:A:318:THR:HG22	5:A:1016:HAS:HMB2	1.96	0.48
2:B:222:ARG:NH1	14:B:2044:HOH:O	2.46	0.48
1:D:367:LEU:C	5:D:1016:HAS:HMD	2.34	0.48
1:A:435:LEU:HB2	1:A:436:PRO:HD3	1.95	0.48
1:A:49:ILE:HG21	5:A:1015:HAS:HMD	1.95	0.48
1:A:675:HIS:HD2	14:A:2157:HOH:O	1.97	0.48
1:A:274:LYS:HE2	1:A:335:THR:O	2.14	0.48
1:D:252:THR:O	1:D:256:MET:HG3	2.13	0.47
1:A:316:PHE:HB2	2:B:99:PHE:CE1	2.49	0.47
2:B:231:PRO:HG3	1:D:715:TYR:OH	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:367:LEU:HB3	5:D:1016:HAS:HMD	1.95	0.47
2:B:90:ILE:HB	2:B:91:PRO:HD3	1.97	0.47
1:D:386:PHE:HB2	5:D:1016:HAS:CMA	2.45	0.47
2:B:142:GLU:HG2	2:B:211:ARG:HB2	1.97	0.47
5:A:1016:HAS:HMC1	5:A:1016:HAS:HBC1	1.97	0.47
3:C:21:ILE:HG23	3:C:22:LEU:HG	1.97	0.47
1:A:370:THR:HB	1:A:371:PRO:HD3	1.97	0.47
1:A:318:THR:CG2	5:A:1016:HAS:HMB2	2.44	0.47
1:D:340:LYS:HE2	2:E:72:GLU:O	2.14	0.47
1:D:316:PHE:HB2	2:E:99:PHE:CE1	2.49	0.47
1:A:175:VAL:HG11	1:A:623:TRP:CE2	2.51	0.46
1:D:103:ASP:OD2	1:D:552:PRO:HG2	2.15	0.46
1:D:390:LEU:HG	5:D:1016:HAS:HAC	1.97	0.46
1:D:256:MET:O	1:D:260:TYR:CD2	2.68	0.46
1:D:264:LEU:HD11	1:D:353:PHE:CG	2.51	0.46
1:A:264:LEU:HD11	1:A:353:PHE:CG	2.51	0.46
1:A:353:PHE:CD1	1:A:353:PHE:C	2.89	0.46
1:D:370:THR:HB	1:D:371:PRO:HD3	1.97	0.46
2:E:280:VAL:HG13	2:E:286:ASN:ND2	2.30	0.46
1:A:252:THR:O	1:A:256:MET:HG3	2.16	0.46
1:A:228:SER:HA	14:A:2066:HOH:O	2.16	0.46
2:E:90:ILE:HB	2:E:91:PRO:HD3	1.98	0.46
1:D:321:ILE:O	1:D:324:PRO:HD2	2.16	0.45
1:D:679:HIS:CE1	1:D:683:ARG:NH1	2.84	0.45
1:D:683:ARG:HG2	1:D:685:ARG:HH21	1.81	0.45
1:A:387:HIS:O	1:A:391:MET:HB3	2.16	0.45
1:A:358:LEU:HD21	2:B:53:VAL:HG11	1.97	0.45
1:D:274:LYS:HE2	1:D:335:THR:O	2.16	0.45
1:D:318:THR:CG2	5:D:1016:HAS:HMB2	2.47	0.45
2:B:264:PRO:HD2	11:B:587:HEC:C3D	2.46	0.45
1:A:143:TYR:CE1	1:A:144:PRO:HB3	2.52	0.45
1:A:179:TYR:CE2	3:C:27:LEU:HD12	2.51	0.45
1:A:256:MET:O	1:A:260:TYR:CD2	2.70	0.45
2:B:135:LEU:HD22	2:B:224:VAL:HG11	1.98	0.45
1:D:353:PHE:C	1:D:353:PHE:CD1	2.90	0.45
1:A:250:HIS:CE1	1:A:254:TYR:OH	2.68	0.45
1:D:515:HIS:O	1:D:516:ASN:HB2	2.17	0.45
2:E:241:GLN:O	2:E:245:GLN:HG3	2.17	0.45
2:B:125:GLN:HA	2:B:126:PHE:HA	1.76	0.45
5:A:1015:HAS:H281	5:A:1015:HAS:H251	1.65	0.45
1:A:250:HIS:O	1:A:253:VAL:HG22	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:402:LEU:O	1:A:406:TRP:HB2	2.17	0.44
4:D:900:5PL:OCO	4:D:900:5PL:HAY1	2.16	0.44
2:E:294:PRO:HB2	2:E:301:VAL:HG11	1.98	0.44
1:A:416:GLU:O	1:A:420:ARG:HG3	2.18	0.44
2:E:238:ARG:O	2:E:242:VAL:HG23	2.18	0.44
1:D:659:LEU:HA	1:D:659:LEU:HD12	1.85	0.44
1:A:422:HIS:CE1	1:A:480:TRP:HB2	2.53	0.44
1:A:210:ALA:HB1	1:A:631:LEU:HA	2.00	0.44
1:A:689:PHE:CE1	1:A:767:LEU:HD22	2.53	0.44
3:F:21:ILE:HG23	3:F:22:LEU:HG	2.00	0.44
1:D:383:VAL:HA	1:D:386:PHE:CE2	2.53	0.44
1:D:298:ALA:HA	1:D:301:MET:HE2	2.00	0.44
1:A:507:LEU:HG	1:A:507:LEU:O	2.17	0.44
1:D:210:ALA:HB1	1:D:631:LEU:HA	2.00	0.44
1:A:29:LEU:HD22	8:A:1300:7E8:H42C	2.00	0.43
1:D:43:GLY:HA3	5:D:1015:HAS:H161	2.00	0.43
1:D:367:LEU:HB2	5:D:1016:HAS:CMD	2.45	0.43
1:D:738:LEU:HD11	4:D:900:5PL:H202	2.00	0.43
1:D:194:TYR:HB2	1:D:266:GLU:HG3	1.99	0.43
1:D:422:HIS:CE1	1:D:480:TRP:HB2	2.53	0.43
1:A:604:HIS:HB3	1:A:605:PRO:HD2	1.99	0.43
1:D:250:HIS:CE1	1:D:254:TYR:OH	2.71	0.43
2:E:84:GLU:HA	2:E:87:TRP:NE1	2.33	0.43
2:B:294:PRO:HB2	2:B:301:VAL:HG11	2.00	0.43
1:D:250:HIS:CD2	1:D:250:HIS:C	2.91	0.43
1:D:250:HIS:O	1:D:253:VAL:HG22	2.18	0.43
1:A:510:SER:HA	1:A:511:PRO:HA	1.88	0.43
4:A:900:5PL:OCO	4:A:900:5PL:HAY1	2.19	0.43
2:E:165:TRP:CG	2:E:172:LYS:HE3	2.53	0.43
1:D:386:PHE:CG	5:D:1016:HAS:HMA2	2.54	0.43
1:D:143:TYR:CE1	1:D:144:PRO:HB3	2.53	0.43
1:D:27:ILE:HD13	1:D:30:MET:HE3	2.01	0.43
8:A:1301:7E8:H191	14:A:2010:HOH:O	2.18	0.43
1:A:515:HIS:O	1:A:516:ASN:HB2	2.19	0.43
1:D:485:TRP:O	1:D:489:ARG:HG2	2.18	0.43
5:A:1015:HAS:HHA	5:A:1015:HAS:HAD2	1.88	0.42
1:A:40:ALA:HB1	5:A:1015:HAS:C21	2.47	0.42
1:A:742:ILE:HG23	4:A:900:5PL:CBR	2.49	0.42
1:A:367:LEU:HD13	5:A:1016:HAS:CBD	2.50	0.42
1:A:298:ALA:CB	1:A:301:MET:HE2	2.49	0.42
1:D:742:ILE:HG23	4:D:900:5PL:CBR	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:98:MET:HB3	1:D:192:PRO:HG2	2.00	0.42
1:D:99:LEU:HD21	1:D:195:VAL:HG11	2.01	0.42
1:D:27:ILE:HD13	1:D:30:MET:CE	2.49	0.42
1:D:175:VAL:HG11	1:D:623:TRP:CE2	2.54	0.42
1:D:586:LEU:HA	1:D:586:LEU:HD12	1.93	0.42
1:A:634:ILE:HD12	1:A:634:ILE:HA	1.94	0.42
1:D:402:LEU:O	1:D:406:TRP:HB2	2.19	0.42
2:B:165:TRP:CG	2:B:172:LYS:HE3	2.55	0.42
1:D:264:LEU:HD11	1:D:353:PHE:CD1	2.55	0.42
1:D:612:THR:HA	3:F:28:THR:HG22	2.01	0.42
2:E:125:GLN:HA	2:E:126:PHE:HA	1.75	0.42
1:A:390:LEU:HD23	5:A:1016:HAS:HBC2	2.02	0.42
1:D:358:LEU:HD21	2:E:53:VAL:HG11	2.02	0.42
1:A:321:ILE:O	1:A:324:PRO:HD2	2.19	0.41
1:A:196:TRP:O	1:A:199:PHE:HB3	2.19	0.41
2:B:308:GLN:HG3	2:E:302:LYS:HA	2.01	0.41
1:D:751:TYR:CE2	1:D:755:LEU:HD11	2.55	0.41
1:D:260:TYR:O	1:D:264:LEU:HG	2.21	0.41
1:A:680:ASP:OD2	1:A:688:PRO:HB2	2.19	0.41
1:D:218:LEU:HD21	1:D:731:THR:HG21	2.03	0.41
1:A:361:GLY:HA3	5:A:1016:HAS:C15	2.51	0.41
1:A:354:ILE:CG2	2:B:57:LEU:HD22	2.51	0.41
4:D:900:5PL:OCO	4:D:900:5PL:HBH	2.19	0.41
1:D:687:ASN:HB2	1:D:688:PRO:HD3	2.03	0.41
1:D:43:GLY:HA3	5:D:1015:HAS:C16	2.51	0.41
1:D:386:PHE:CD2	5:D:1016:HAS:HAA1	2.56	0.41
1:A:27:ILE:HD13	1:A:30:MET:CE	2.50	0.41
2:E:275:LEU:HD12	11:E:587:HEC:HAA1	2.02	0.41
1:A:250:HIS:C	1:A:250:HIS:CD2	2.94	0.41
1:A:264:LEU:HD11	1:A:353:PHE:CD1	2.55	0.41
1:D:406:TRP:HB3	1:D:407:PRO:HD3	2.01	0.41
2:B:223:PHE:HA	2:B:329:PHE:CZ	2.55	0.41
1:D:274:LYS:NZ	1:D:337:TRP:O	2.51	0.41
1:A:751:TYR:CE2	1:A:755:LEU:HD11	2.55	0.41
2:E:232:ALA:HA	2:E:233:PRO:HD3	1.95	0.41
1:D:507:LEU:O	1:D:507:LEU:HG	2.21	0.41
1:A:99:LEU:HD21	1:A:195:VAL:HG11	2.03	0.41
1:D:390:LEU:HD11	5:D:1016:HAS:HMA1	2.02	0.41
1:D:318:THR:HG22	5:D:1016:HAS:HMB2	2.02	0.41
1:A:367:LEU:HD21	1:A:381:PHE:HD1	1.86	0.41
2:B:232:ALA:HA	2:B:233:PRO:HD3	1.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:GLN:HG3	1:A:136:SER:HB3	2.03	0.41
1:A:367:LEU:CB	5:A:1016:HAS:CMD	2.98	0.41
1:D:196:TRP:O	1:D:199:PHE:HB3	2.21	0.41
1:D:510:SER:HA	1:D:511:PRO:HA	1.88	0.40
2:E:158:LYS:HA	2:E:158:LYS:HD3	1.92	0.40
1:D:114:TYR:O	1:D:117:PHE:HB3	2.20	0.40
1:A:194:TYR:HB2	1:A:266:GLU:HG3	2.03	0.40
2:B:69:GLU:O	2:B:71:GLN:HG2	2.21	0.40
5:D:1015:HAS:H251	5:D:1015:HAS:H281	1.35	0.40
1:D:459:TRP:HB2	1:D:460:PRO:HD3	2.02	0.40
1:D:363:THR:HG21	1:D:438:TYR:CE1	2.56	0.40
1:A:231:ASN:HA	1:A:232:PRO:HD3	1.92	0.40
2:B:280:VAL:HG13	2:B:286:ASN:ND2	2.36	0.40
1:D:465:LEU:HA	1:D:465:LEU:HD23	1.95	0.40
1:D:649:THR:HA	1:D:650:PRO:HD3	1.96	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	779/791 (98%)	762 (98%)	17 (2%)	0	100	100
1	D	778/791 (98%)	759 (98%)	19 (2%)	0	100	100
2	B	317/337 (94%)	308 (97%)	9 (3%)	0	100	100
2	E	317/337 (94%)	307 (97%)	10 (3%)	0	100	100
3	C	62/66 (94%)	62 (100%)	0	0	100	100
3	F	61/66 (92%)	61 (100%)	0	0	100	100
All	All	2314/2388 (97%)	2259 (98%)	55 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	640/646 (99%)	637 (100%)	3 (0%)	92	97
1	D	639/646 (99%)	638 (100%)	1 (0%)	95	99
2	B	261/274 (95%)	259 (99%)	2 (1%)	86	94
2	E	261/274 (95%)	256 (98%)	5 (2%)	65	79
3	C	51/53 (96%)	51 (100%)	0	100	100
3	F	50/53 (94%)	50 (100%)	0	100	100
All	All	1902/1946 (98%)	1891 (99%)	11 (1%)	90	96

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	378	ASP
1	A	689	PHE
1	A	759	ILE
2	B	194	TYR
2	B	272	ARG
1	D	378	ASP
2	E	50	PHE
2	E	87	TRP
2	E	194	TYR
2	E	238	ARG
2	E	272	ARG

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	171	ASN
1	A	205	ASN
2	B	245	GLN
3	C	63	HIS
1	D	171	ASN
1	D	205	ASN

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Mol	Chain	Res	Type
1	D	764	HIS
3	F	63	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	FME	C	1	3	8,9,10	0.32	0	6,9,11	1.50	1 (16%)
3	FME	F	1	3	8,9,10	0.34	0	6,9,11	0.88	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
3	FME	C	1	3	-	0/6/9/11	0/0/0/0
3	FME	F	1	3	-	0/6/9/11	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	C	1	FME	CA-N-CN	-2.92	118.33	122.82

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 21 ligands modelled in this entry, 7 are monoatomic - leaving 14 for Mogul analysis.

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	$\# Z > 2$
5	HAS	A	1015	1	45,72,72	1.11	3 (6%)	47,109,109	2.74	19 (40%)
5	HAS	A	1016	1,14	45,72,72	1.17	3 (6%)	47,109,109	2.91	19 (40%)
7	4AG	A	1200	-	39,39,39	1.02	2 (5%)	41,41,41	1.22	3 (7%)
8	7E8	A	1300	-	20,20,20	1.02	1 (5%)	21,21,21	0.99	1 (4%)
8	7E8	A	1301	-	20,20,20	0.98	1 (5%)	21,21,21	1.05	2 (9%)
4	5PL	A	900	-	83,85,85	0.78	2 (2%)	94,101,101	1.07	5 (5%)
10	CUA	B	585	2	0,1,1	0.00	-	0,0,0	0.00	-
11	HEC	B	587	2	24,50,50	2.46	3 (12%)	19,82,82	3.19	6 (31%)
12	7E9	B	701	-	20,20,20	1.02	1 (5%)	20,21,21	1.15	1 (5%)
5	HAS	D	1015	1	45,72,72	1.16	3 (6%)	47,109,109	2.83	23 (48%)
5	HAS	D	1016	1	45,72,72	1.07	3 (6%)	47,109,109	3.03	21 (44%)
4	5PL	D	900	-	83,85,85	0.78	2 (2%)	94,101,101	1.08	6 (6%)
10	CUA	E	585	2	0,1,1	0.00	-	0,0,0	0.00	-
11	HEC	E	587	2	24,50,50	2.50	3 (12%)	19,82,82	3.41	8 (42%)

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
5	HAS	A	1015	1	-	0/30/82/82	0/0/8/8
5	HAS	A	1016	1,14	-	0/30/82/82	0/0/8/8
7	4AG	A	1200	-	-	0/41/41/41	0/0/0/0
8	7E8	A	1300	-	1/1/2/4	0/20/20/20	0/0/0/0
8	7E8	A	1301	-	-	0/20/20/20	0/0/0/0
4	5PL	A	900	-	-	0/87/107/107	0/1/1/1
10	CUA	B	585	2	-	0/0/0/0	0/0/0/0
11	HEC	B	587	2	-	0/6/54/54	0/0/8/8
12	7E9	B	701	-	-	0/21/21/21	0/0/0/0
5	HAS	D	1015	1	-	0/30/82/82	0/0/8/8
5	HAS	D	1016	1	-	0/30/82/82	0/0/8/8
4	5PL	D	900	-	-	0/87/107/107	0/1/1/1
10	CUA	E	585	2	-	0/0/0/0	0/0/0/0
11	HEC	E	587	2	-	0/6/54/54	0/0/8/8

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	E	587	HEC	C3B-C2B	-7.22	1.33	1.40
11	B	587	HEC	C3B-C2B	-6.93	1.33	1.40
11	B	587	HEC	C3C-C2C	-6.67	1.33	1.40
11	E	587	HEC	C3C-C2C	-6.52	1.33	1.40
5	D	1015	HAS	C2A-C3A	2.81	1.46	1.37
5	A	1016	HAS	C2A-C3A	3.06	1.46	1.37
5	D	1016	HAS	C2A-C3A	3.07	1.46	1.37
5	A	1015	HAS	C2A-C3A	3.08	1.46	1.37
4	D	900	5PL	OCL-CBL	3.82	1.44	1.33
5	D	1016	HAS	C3C-C2C	3.92	1.45	1.40
5	A	1015	HAS	C3C-C2C	3.99	1.45	1.40
7	A	1200	4AG	O21-C22	4.00	1.46	1.34
12	B	701	7E9	O1-C80	4.01	1.46	1.34
5	A	1016	HAS	C3C-C2C	4.07	1.45	1.40
8	A	1301	7E8	O16-C8	4.14	1.45	1.33
4	A	900	5PL	OCK-CBU	4.15	1.46	1.34
5	D	1015	HAS	C3C-C2C	4.19	1.45	1.40
5	A	1015	HAS	C2D-C3D	4.21	1.45	1.40
5	D	1016	HAS	C2D-C3D	4.25	1.45	1.40
4	A	900	5PL	OCL-CBL	4.29	1.46	1.33
8	A	1300	7E8	O16-C8	4.32	1.46	1.33
7	A	1200	4AG	O17-C16	4.39	1.46	1.33
4	D	900	5PL	OCK-CBU	4.39	1.47	1.34
11	B	587	HEC	C3D-C2D	4.92	1.52	1.37
5	A	1016	HAS	C2D-C3D	5.02	1.46	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	1015	HAS	C2D-C3D	5.10	1.47	1.40
11	E	587	HEC	C3D-C2D	5.21	1.53	1.37

All (114) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	B	587	HEC	CBB-CAB-C3B	-8.89	107.59	127.35
11	E	587	HEC	CBB-CAB-C3B	-8.79	107.81	127.35
5	A	1016	HAS	CAD-CBD-CGD	-8.55	97.07	112.75
11	B	587	HEC	CBD-CAD-C3D	-7.86	98.45	112.53
11	E	587	HEC	CBD-CAD-C3D	-7.72	98.69	112.53
5	D	1016	HAS	CAD-CBD-CGD	-7.26	99.43	112.75
5	D	1016	HAS	CMA-C3A-C4A	-6.48	117.65	128.36
5	A	1016	HAS	C4D-C3D-C2D	-6.27	100.79	107.07
5	D	1016	HAS	C4D-C3D-C2D	-6.25	100.82	107.07
5	A	1015	HAS	C4D-C3D-C2D	-6.22	100.84	107.07
5	D	1015	HAS	CAD-CBD-CGD	-6.12	101.53	112.75
11	E	587	HEC	CBC-CAC-C3C	-5.94	114.16	127.35
5	A	1016	HAS	CAA-CBA-CGA	-5.84	102.04	112.75
5	D	1016	HAS	C13-C12-C11	-5.81	106.78	114.51
5	D	1016	HAS	CMC-C2C-C1C	-5.77	118.82	128.36
5	D	1015	HAS	C4D-C3D-C2D	-5.65	101.41	107.07
5	A	1016	HAS	CMC-C2C-C1C	-5.56	119.16	128.36
5	A	1015	HAS	CAD-CBD-CGD	-5.54	102.60	112.75
5	A	1016	HAS	CMA-C3A-C4A	-5.49	119.28	128.36
5	A	1015	HAS	CMA-C3A-C4A	-5.42	119.39	128.36
5	D	1015	HAS	CMC-C2C-C1C	-5.27	119.64	128.36
5	D	1015	HAS	CBA-CAA-C2A	-5.25	103.11	112.53
5	A	1016	HAS	C13-C12-C11	-4.73	108.23	114.51
5	D	1015	HAS	C13-C12-C11	-4.65	108.33	114.51
5	A	1015	HAS	CMC-C2C-C1C	-4.50	120.92	128.36
5	D	1015	HAS	CMA-C3A-C4A	-4.48	120.95	128.36
5	A	1015	HAS	OMD-CMD-C2D	-4.41	116.20	125.11
5	D	1016	HAS	CAA-CBA-CGA	-4.33	104.80	112.75
5	D	1016	HAS	C13-C14-C15	-4.29	118.43	127.76
5	A	1015	HAS	CBA-CAA-C2A	-4.13	105.12	112.53
11	E	587	HEC	CBA-CAA-C2A	-4.09	105.20	112.53
5	A	1015	HAS	C4B-C3B-C11	-4.05	122.61	127.01
5	D	1015	HAS	C4B-C3B-C11	-4.00	122.66	127.01
5	A	1015	HAS	C3C-CAC-CBC	-3.84	118.46	126.32
11	B	587	HEC	CMC-C2C-C1C	-3.79	122.09	128.36
5	D	1015	HAS	C21-C22-C23	-3.76	119.59	127.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	1015	HAS	C13-C14-C15	-3.73	119.65	127.76
5	A	1016	HAS	C13-C14-C15	-3.70	119.72	127.76
5	A	1015	HAS	C13-C12-C11	-3.55	109.80	114.51
5	A	1016	HAS	OMD-CMD-C2D	-3.46	118.12	125.11
11	B	587	HEC	CBC-CAC-C3C	-3.46	119.67	127.35
4	D	900	5PL	CBH-OCK-CBU	-3.45	109.61	117.89
4	A	900	5PL	CBH-OCK-CBU	-3.23	110.15	117.89
5	A	1016	HAS	C21-C22-C23	-3.13	120.95	127.76
5	D	1016	HAS	C3C-CAC-CBC	-3.12	119.94	126.32
5	D	1015	HAS	C3C-CAC-CBC	-3.11	119.95	126.32
5	D	1015	HAS	C17-C18-C19	-3.11	121.00	127.76
5	D	1016	HAS	C21-C22-C23	-2.97	121.31	127.76
5	D	1016	HAS	C4B-C3B-C11	-2.89	123.87	127.01
5	A	1015	HAS	C17-C18-C19	-2.86	121.53	127.76
5	D	1016	HAS	OMD-CMD-C2D	-2.83	119.40	125.11
4	A	900	5PL	CAY-CAX-NCB	-2.81	103.97	112.19
5	D	1015	HAS	OMD-CMD-C2D	-2.80	119.46	125.11
5	A	1015	HAS	C21-C22-C23	-2.69	121.91	127.76
5	A	1016	HAS	C4B-C3B-C11	-2.69	124.09	127.01
5	D	1016	HAS	C17-C18-C19	-2.67	121.95	127.76
5	A	1016	HAS	C17-C18-C19	-2.67	121.96	127.76
4	D	900	5PL	C3-C2-N2	-2.63	105.21	110.66
5	A	1015	HAS	C13-C14-C15	-2.59	122.14	127.76
11	E	587	HEC	CMC-C2C-C1C	-2.57	124.11	128.36
7	A	1200	4AG	C6-O21-C22	-2.43	112.06	117.89
5	D	1016	HAS	C26-C15-C14	-2.37	118.85	123.50
5	A	1015	HAS	C26-C15-C14	-2.30	118.98	123.50
8	A	1301	7E8	O16-C8-O15	-2.29	117.58	123.49
5	A	1016	HAS	C3C-CAC-CBC	-2.18	121.86	126.32
11	E	587	HEC	CAA-C2A-C1A	-2.16	124.67	127.01
11	B	587	HEC	CBA-CAA-C2A	-2.06	108.83	112.53
5	D	1015	HAS	C20-C21-C22	-2.06	106.30	111.69
5	D	1015	HAS	C28-C29-C30	-2.00	120.02	127.73
5	D	1015	HAS	CAA-C2A-C1A	2.01	129.19	127.01
11	E	587	HEC	C4B-C3B-C2B	2.03	108.55	106.35
11	B	587	HEC	C4B-C3B-C2B	2.07	108.59	106.35
5	D	1016	HAS	C32-C30-C31	2.12	119.86	114.64
5	A	1016	HAS	C32-C30-C31	2.13	119.88	114.64
5	A	1015	HAS	CBD-CAD-C3D	2.14	116.36	112.53
4	D	900	5PL	O5-C5-C4	2.18	113.78	109.68
11	E	587	HEC	CAD-CBD-CGD	2.22	116.81	112.75
5	A	1016	HAS	C25-C23-C24	2.25	118.85	115.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	1015	HAS	CBD-CAD-C3D	2.26	116.57	112.53
5	D	1015	HAS	C32-C30-C31	2.26	120.19	114.64
5	A	1016	HAS	C26-C15-C16	2.31	118.94	115.41
5	D	1016	HAS	C25-C23-C24	2.34	118.99	115.41
4	D	900	5PL	O1-C1-C2	2.52	110.92	107.57
4	A	900	5PL	OCL-CBL-CBM	2.63	119.91	111.90
5	D	1015	HAS	C26-C15-C16	2.65	119.45	115.41
8	A	1300	7E8	O16-C8-C7	2.72	120.18	111.90
7	A	1200	4AG	O17-C16-C15	2.80	120.42	111.90
5	A	1016	HAS	C27-C19-C20	2.83	119.73	115.41
4	D	900	5PL	OCL-CBL-CBM	2.85	120.58	111.90
5	D	1015	HAS	C25-C23-C24	2.89	119.82	115.41
5	D	1015	HAS	C27-C19-C20	2.91	119.86	115.41
5	D	1016	HAS	C26-C15-C16	3.04	120.05	115.41
5	D	1016	HAS	C27-C19-C20	3.07	120.09	115.41
5	A	1015	HAS	C25-C23-C24	3.09	120.13	115.41
5	A	1016	HAS	CMB-C2B-C3B	3.29	131.87	125.14
8	A	1301	7E8	O16-C8-C7	3.32	122.02	111.90
4	A	900	5PL	O1-C1-C2	3.37	112.06	107.57
5	A	1015	HAS	C26-C15-C16	3.44	120.66	115.41
5	D	1015	HAS	CMA-C3A-C2A	3.54	132.63	125.24
5	A	1015	HAS	CMB-C2B-C3B	3.86	133.03	125.14
12	B	701	7E9	O1-C80-C2	3.99	120.20	111.53
7	A	1200	4AG	O21-C22-C23	4.07	120.37	111.53
5	D	1016	HAS	CAD-C3D-C4D	4.15	131.51	127.01
5	A	1016	HAS	CMA-C3A-C2A	4.16	133.93	125.24
4	A	900	5PL	OCK-CBU-CBV	4.17	120.58	111.53
4	D	900	5PL	OCK-CBU-CBV	4.26	120.78	111.53
5	A	1015	HAS	CMA-C3A-C2A	4.41	134.46	125.24
5	D	1015	HAS	CMB-C2B-C3B	4.66	134.67	125.14
5	D	1016	HAS	CMA-C3A-C2A	4.93	135.54	125.24
5	D	1016	HAS	CMB-C2B-C3B	4.98	135.31	125.14
5	A	1015	HAS	CMC-C2C-C3C	5.06	134.98	125.09
5	D	1015	HAS	CMC-C2C-C3C	5.48	135.80	125.09
5	A	1016	HAS	CMC-C2C-C3C	5.69	136.22	125.09
5	D	1016	HAS	CMC-C2C-C3C	5.75	136.34	125.09

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
8	A	1300	7E8	C18

There are no torsion outliers.

There are no ring outliers.

11 monomers are involved in 59 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1015	HAS	9	0
5	A	1016	HAS	9	0
7	A	1200	4AG	1	0
8	A	1300	7E8	1	0
8	A	1301	7E8	1	0
4	A	900	5PL	5	0
11	B	587	HEC	3	0
5	D	1015	HAS	7	0
5	D	1016	HAS	16	0
4	D	900	5PL	4	0
11	E	587	HEC	3	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	780/791 (98%)	0.21	14 (1%) 71 81	32, 48, 91, 174	0
1	D	780/791 (98%)	0.22	17 (2%) 65 77	33, 53, 89, 153	0
2	B	319/337 (94%)	0.26	3 (0%) 85 92	33, 50, 85, 163	0
2	E	319/337 (94%)	0.37	15 (4%) 35 50	38, 67, 105, 189	0
3	C	63/66 (95%)	0.16	1 (1%) 74 84	44, 62, 112, 152	0
3	F	62/66 (93%)	0.17	3 (4%) 34 49	45, 65, 104, 153	0
All	All	2323/2388 (97%)	0.24	53 (2%) 64 76	32, 53, 94, 189	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	539	VAL	5.6
2	E	69	GLU	5.6
2	E	68	PRO	5.3
2	E	72	GLU	4.5
1	D	548	HIS	4.3
1	D	547	ALA	4.0
1	A	548	HIS	4.0
2	B	69	GLU	3.8
1	D	452	TYR	3.4
1	A	607	GLU	3.1
1	A	12	TRP	3.1
2	E	35	VAL	3.0
1	A	762	HIS	3.0
2	E	25	THR	3.0
1	D	355	PHE	3.0
2	E	58	ALA	3.0
1	D	607	GLU	3.0
2	E	224	VAL	2.9
1	A	683	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	608	HIS	2.9
2	E	71	GLN	2.8
1	A	761	LEU	2.7
1	D	15	LEU	2.7
1	D	456	ILE	2.7
1	A	691	PHE	2.6
2	E	337	PHE	2.6
1	D	539	VAL	2.6
2	B	72	GLU	2.6
2	E	70	ASP	2.6
2	E	55	GLY	2.5
2	E	87	TRP	2.5
2	E	24	ILE	2.5
1	A	536	LYS	2.5
3	C	38	PHE	2.5
1	D	606	VAL	2.5
1	D	12	TRP	2.4
1	A	545	ASP	2.4
1	D	443	LEU	2.4
2	E	59	TYR	2.4
1	A	605	PRO	2.3
2	B	71	GLN	2.3
1	D	521	LEU	2.3
1	A	679	HIS	2.3
1	D	458	GLY	2.3
1	D	502	TYR	2.3
1	A	689	PHE	2.3
3	F	27	LEU	2.2
1	A	763	ASN	2.2
3	F	24	PRO	2.2
3	F	25	ARG	2.1
1	D	442	TYR	2.1
1	D	549	ILE	2.1
2	E	62	TRP	2.0

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

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
3	FME	F	1	10/11	0.92	0.14	-	72,91,112,160	0
3	FME	C	1	10/11	0.93	0.13	-	69,92,107,171	0

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
8	7E8	A	1300	21/21	0.76	0.31	8.44	48,79,102,115	0
7	4AG	A	1200	40/40	0.89	0.27	7.40	33,68,115,123	0
8	7E8	A	1301	21/21	0.86	0.25	6.36	44,80,96,118	0
9	MG	A	1801	1/1	0.98	0.19	4.96	39,39,39,39	0
12	7E9	B	701	21/21	0.84	0.26	4.67	58,90,108,109	0
10	CUA	B	585	2/2	1.00	0.17	1.54	35,35,35,36	0
5	HAS	A	1015	65/65	0.96	0.18	1.45	24,35,102,151	0
4	5PL	D	900	85/85	0.90	0.19	1.38	37,62,107,128	0
4	5PL	A	900	85/85	0.92	0.18	1.19	35,65,105,134	0
5	HAS	D	1015	65/65	0.95	0.17	1.08	30,55,128,138	0
13	CL	D	1701	1/1	0.96	0.17	1.05	86,86,86,86	0
11	HEC	E	587	43/43	0.97	0.17	1.05	21,54,80,91	0
10	CUA	E	585	2/2	0.99	0.16	0.90	43,43,43,44	0
5	HAS	D	1016	65/65	0.97	0.16	0.56	24,47,86,103	0
11	HEC	B	587	43/43	0.98	0.15	-0.06	22,34,48,52	0
5	HAS	A	1016	65/65	0.97	0.15	-0.42	22,38,59,62	0
9	MG	D	1801	1/1	0.98	0.12	-1.56	50,50,50,50	0
13	CL	F	101	1/1	0.94	0.27	-	70,70,70,70	0
6	CU	A	1017	1/1	1.00	0.20	-	37,37,37,37	0
6	CU	D	1017	1/1	1.00	0.15	-	42,42,42,42	0
13	CL	C	101	1/1	0.98	0.28	-	60,60,60,60	0

6.5 Other polymers

There are no such residues in this entry.