



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 15, 2024 – 03:04 PM EDT

PDB ID : 2G39
Title : Crystal structure of coenzyme A transferase from *Pseudomonas aeruginosa*
Authors : Chang, C.; Evdokimova, E.; Kudritska, M.; Savchenko, A.; Edwards, A.;
Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2006-02-17
Resolution : 2.10 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 1.20.1
EDS : 2.37.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.37.1

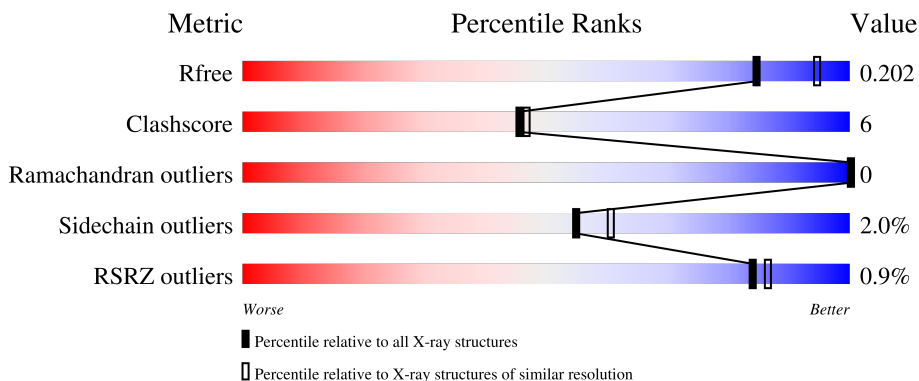
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.10 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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	497	 86% 12% •
1	B	497	 % 88% 10% ••

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
3	ACY	B	902	-	-	X	-
3	ACY	B	905	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8404 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 Acetyl-CoA hydrolase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	495	3792	2376	681	718	3	14	0	4	0
1	B	494	3780	2367	679	717	3	14	0	3	0

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
A	16	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
A	29	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
A	33	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
A	51	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
A	63	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
A	86	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
A	104	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
A	253	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
A	272	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
A	285	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
A	376	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
A	377	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
A	412	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
A	494	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
B	1	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
B	16	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
B	29	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
B	33	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
B	51	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
B	63	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
B	86	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
B	104	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
B	253	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
B	272	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	285	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
B	376	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
B	377	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
B	412	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2
B	494	MSE	MET	MODIFIED RESIDUE	UNP Q9HTC2

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	4	2	2	0	0

- Molecule 3 is ACETIC ACID (three-letter code: ACY) (formula: C₂H₄O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0

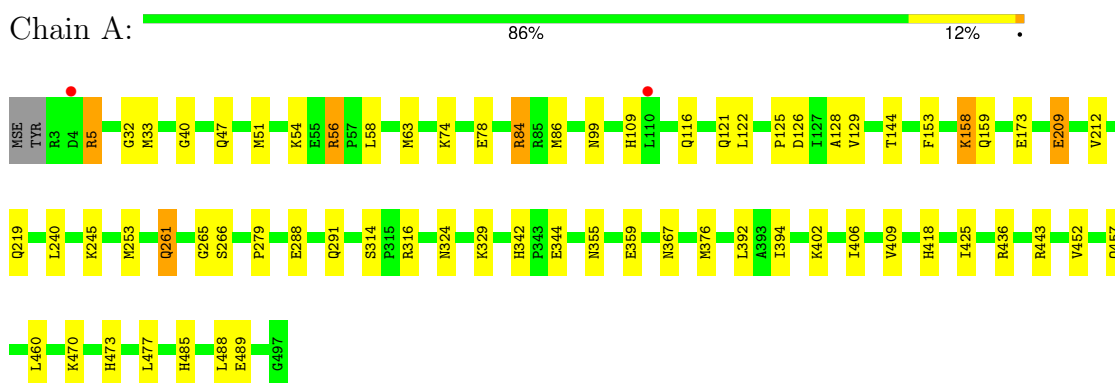
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	411	Total O 411 411	0	0
4	B	385	Total O 385 385	0	0

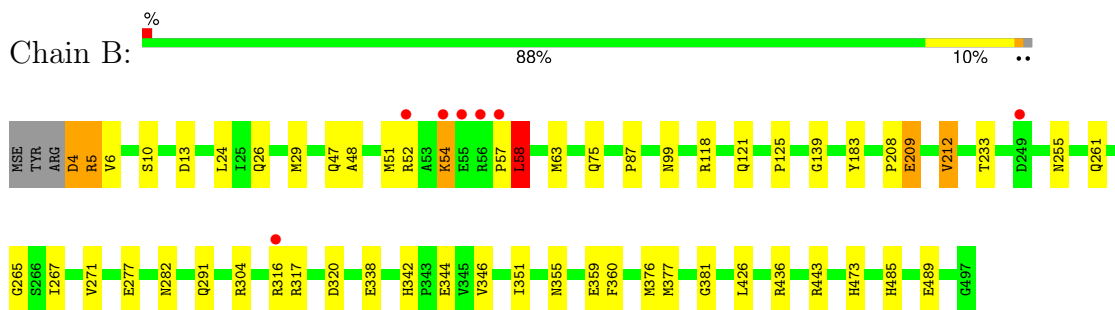
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: Acetyl-CoA hydrolase



- Molecule 1: Acetyl-CoA hydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	171.36Å 69.81Å 113.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.10 44.21 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.0 (50.00-2.10) 99.0 (44.21-2.10)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.19 (at 2.10Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.162 , 0.204 0.163 , 0.202	Depositor DCC
R_{free} test set	3940 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	24.3	Xtrriage
Anisotropy	0.395	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 42.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8404	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: EDO, ACY

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	0.74	0/3857	0.76	7/5208 (0.1%)
1	B	0.73	0/3842	0.78	5/5189 (0.1%)
All	All	0.73	0/7699	0.77	12/10397 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	84	ARG	NE-CZ-NH2	-7.49	116.56	120.30
1	A	443[A]	ARG	NE-CZ-NH2	-7.38	116.61	120.30
1	A	443[B]	ARG	NE-CZ-NH2	-7.38	116.61	120.30
1	A	5	ARG	NE-CZ-NH2	7.21	123.90	120.30
1	B	58	LEU	N-CA-C	6.81	129.38	111.00
1	A	84	ARG	NE-CZ-NH1	6.06	123.33	120.30
1	A	56	ARG	NE-CZ-NH1	5.73	123.16	120.30
1	B	443[A]	ARG	NE-CZ-NH2	-5.69	117.46	120.30
1	B	443[B]	ARG	NE-CZ-NH2	-5.69	117.46	120.30
1	A	5	ARG	NE-CZ-NH1	-5.64	117.48	120.30
1	B	4	ASP	C-N-CA	5.42	135.26	121.70
1	B	5	ARG	N-CA-C	5.37	125.50	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	57	PRO	Peptide

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	3792	0	3831	53	0
1	B	3780	0	3814	48	1
2	A	4	0	6	0	0
3	A	20	0	15	2	0
3	B	12	0	9	8	0
4	A	411	0	0	13	0
4	B	385	0	0	13	0
All	All	8404	0	7675	98	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (98) 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:51:MSE:HG3	4:A:1149:HOH:O	1.40	1.20
1:B:118:ARG:HE	3:B:902:ACY:H1	0.92	1.04
1:B:317:ARG:HD2	4:B:1208:HOH:O	1.61	1.01
1:B:118:ARG:NE	3:B:902:ACY:H1	1.78	0.99
1:A:86:MSE:HE3	4:A:1282:HOH:O	1.72	0.88
1:B:48:ALA:O	1:B:52:ARG:HG2	1.74	0.88
1:A:261:GLN:HE22	1:A:355:ASN:HD21	1.29	0.80
1:A:342:HIS:HD2	1:A:344:GLU:H	1.28	0.80
1:A:99:ASN:HD21	1:A:376:MSE:H	1.31	0.78
1:B:118:ARG:HE	3:B:902:ACY:CH3	1.87	0.78
1:A:265:GLY:H	1:A:291:GLN:HE22	1.32	0.77
1:B:342:HIS:HD2	1:B:344:GLU:H	1.32	0.75
1:B:99:ASN:HD21	1:B:376:MSE:H	1.34	0.74
1:A:86:MSE:CE	4:A:1282:HOH:O	2.29	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:233:THR:HG22	1:B:267:ILE:HD13	1.72	0.72
1:B:265:GLY:H	1:B:291:GLN:HE22	1.37	0.71
1:A:324:ASN:HB3	4:A:1304:HOH:O	1.92	0.70
1:B:261:GLN:HE22	1:B:355:ASN:HD21	1.41	0.69
1:B:26:GLN:HE21	1:B:29:MSE:HE3	1.57	0.68
3:A:909:ACY:H3	4:A:1178:HOH:O	1.97	0.64
1:A:84:ARG:HD3	4:B:1025:HOH:O	1.98	0.63
1:B:47:GLN:HG3	1:B:75:GLN:OE1	1.99	0.63
1:A:47:GLN:HE22	1:A:219:GLN:HE22	1.47	0.63
1:B:24:LEU:HD13	1:B:212:VAL:HG13	1.82	0.62
1:A:54:LYS:HD2	4:A:1308:HOH:O	1.99	0.62
1:B:183:TYR:OH	3:B:902:ACY:H2	2.01	0.61
1:A:488:LEU:HD23	4:A:1306:HOH:O	2.02	0.60
1:B:4:ASP:HB3	1:B:5:ARG:HB2	1.83	0.59
1:A:436:ARG:HH11	1:B:121:GLN:HE22	1.51	0.59
4:A:974:HOH:O	1:B:473:HIS:HE1	1.87	0.58
1:A:359:GLU:OE1	1:A:485:HIS:HE1	1.87	0.57
1:B:58:LEU:H	1:B:58:LEU:HD12	1.70	0.57
1:B:282:ASN:ND2	1:B:304[B]:ARG:HE	2.03	0.56
1:A:436:ARG:HH11	1:B:121:GLN:NE2	2.03	0.56
1:B:316:ARG:HH11	1:B:316:ARG:HG2	1.70	0.56
1:A:261:GLN:HE22	1:A:355:ASN:ND2	2.01	0.56
1:A:473:HIS:H	1:B:99:ASN:HD22	1.53	0.56
1:B:381:GLY:HA3	3:B:905:ACY:H3	1.86	0.56
1:A:63:MSE:HE3	1:A:153:PHE:CE2	2.41	0.56
1:B:52:ARG:HG3	4:B:1271:HOH:O	2.05	0.55
3:B:905:ACY:H1	4:B:1022:HOH:O	2.08	0.54
1:A:342:HIS:HE1	4:A:922:HOH:O	1.91	0.53
3:B:905:ACY:CH3	4:B:1022:HOH:O	2.56	0.53
1:A:63:MSE:HE1	1:A:125:PRO:HD3	1.91	0.52
1:A:109:HIS:HD2	4:A:911:HOH:O	1.92	0.52
1:B:183:TYR:OH	3:B:902:ACY:CH3	2.58	0.52
1:A:342:HIS:CD2	1:A:344:GLU:HB3	2.45	0.51
1:A:470:LYS:HD2	1:A:477:LEU:HD23	1.92	0.51
4:A:933:HOH:O	1:B:473:HIS:HD2	1.94	0.50
1:A:158[A]:LYS:HG3	4:A:1139:HOH:O	2.11	0.50
1:A:314:SER:HB3	1:A:316[B]:ARG:HH22	1.76	0.50
1:A:99:ASN:HD21	1:A:376:MSE:N	2.07	0.49
1:A:240:LEU:HD11	1:A:394:ILE:HD13	1.94	0.49
1:A:121:GLN:NE2	1:B:436:ARG:HH11	2.09	0.49
1:A:418:HIS:HD2	4:B:929:HOH:O	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:54:LYS:HD2	4:B:1249:HOH:O	2.14	0.48
1:A:126:ASP:O	1:A:158[B]:LYS:HG3	2.13	0.48
1:B:342:HIS:HE1	4:B:916:HOH:O	1.96	0.47
1:A:99:ASN:HD22	1:B:473:HIS:H	1.61	0.47
1:B:316:ARG:HG2	1:B:316:ARG:NH1	2.30	0.47
1:A:209:GLU:H	1:A:209:GLU:CD	2.17	0.47
1:A:452:VAL:HG21	1:A:460:LEU:HD23	1.96	0.47
1:B:255:ASN:ND2	4:B:1210:HOH:O	2.48	0.46
1:B:377:MSE:CE	4:B:1253:HOH:O	2.64	0.46
1:A:159:GLN:NE2	1:A:212:VAL:HG11	2.30	0.46
1:A:359:GLU:HG2	1:A:367:ASN:HB3	1.98	0.46
1:A:253:MSE:HE1	1:A:392:LEU:CD1	2.46	0.45
1:A:394:ILE:HG12	1:A:425:ILE:HB	1.98	0.45
1:A:402:LYS:HD3	1:A:406:ILE:HD12	1.98	0.45
1:B:346:VAL:HA	1:B:351:ILE:HD12	1.98	0.45
1:A:121:GLN:HE22	1:B:436:ARG:HH11	1.65	0.45
1:B:359:GLU:OE1	1:B:485:HIS:HE1	1.99	0.45
1:B:63:MSE:HE1	1:B:125:PRO:HD3	1.98	0.44
1:A:265:GLY:H	1:A:291:GLN:NE2	2.08	0.44
1:B:209:GLU:H	1:B:209:GLU:HG3	1.50	0.44
1:B:24:LEU:HD13	1:B:212:VAL:CG1	2.46	0.44
1:A:33:MSE:HG3	1:A:129:VAL:HB	2.00	0.43
1:B:267:ILE:O	1:B:271:VAL:HG23	2.18	0.43
1:B:342:HIS:CD2	1:B:344:GLU:HB3	2.53	0.43
1:B:139:GLY:HA3	1:B:208:PRO:HG3	2.00	0.42
1:B:360:PHE:CD1	1:B:426:LEU:HD13	2.54	0.42
1:A:245:LYS:HE2	1:A:279:PRO:HD3	2.00	0.42
1:A:288:GLU:OE1	3:A:901:ACY:OXT	2.37	0.42
1:B:4:ASP:HB2	4:B:1286:HOH:O	2.18	0.42
1:A:173:GLU:O	1:A:329:LYS:NZ	2.51	0.42
1:B:4:ASP:N	1:B:6:VAL:H	2.18	0.42
1:B:338:GLU:O	1:B:342:HIS:HB2	2.19	0.42
1:A:116:GLN:HE21	1:A:122:LEU:HD22	1.84	0.42
1:A:32:GLY:O	1:A:128:ALA:HA	2.19	0.41
1:A:473:HIS:HE1	4:B:1000:HOH:O	2.01	0.41
1:A:74:LYS:O	1:A:78:GLU:HG3	2.20	0.41
1:A:5:ARG:HD3	4:A:1209:HOH:O	2.21	0.41
1:A:409:VAL:HG11	1:A:452:VAL:HG22	2.03	0.41
1:A:158[B]:LYS:HG3	1:A:158[B]:LYS:H	1.61	0.41
1:A:40:GLY:HA2	1:A:144:THR:OG1	2.21	0.41
1:A:316[B]:ARG:HE	1:A:316[B]:ARG:HB3	1.58	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:489:GLU:HG3	4:B:1289:HOH:O	2.20	0.41
1:B:10:SER:O	1:B:13:ASP:HB2	2.21	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:13:ASP:OD1	1:B:316:ARG:NH2[3_555]	2.18	0.02

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	497/497 (100%)	486 (98%)	11 (2%)	0	100	100
1	B	495/497 (100%)	478 (97%)	17 (3%)	0	100	100
All	All	992/994 (100%)	964 (97%)	28 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	411/395 (104%)	402 (98%)	9 (2%)	52	57

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	410/395 (104%)	402 (98%)	8 (2%)	55	60
All	All	821/790 (104%)	804 (98%)	17 (2%)	55	59

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	ARG
1	A	58	LEU
1	A	158[A]	LYS
1	A	158[B]	LYS
1	A	209	GLU
1	A	261	GLN
1	A	266	SER
1	A	457	GLN
1	A	489	GLU
1	B	51	MSE
1	B	54	LYS
1	B	58	LEU
1	B	87	PRO
1	B	209	GLU
1	B	212	VAL
1	B	277	GLU
1	B	320	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	47	GLN
1	A	99	ASN
1	A	109	HIS
1	A	116	GLN
1	A	121	GLN
1	A	234	GLN
1	A	255	ASN
1	A	261	GLN
1	A	291	GLN
1	A	342	HIS
1	A	415	HIS
1	A	418	HIS
1	A	473	HIS

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Mol	Chain	Res	Type
1	A	485	HIS
1	B	26	GLN
1	B	99	ASN
1	B	109	HIS
1	B	116	GLN
1	B	121	GLN
1	B	138	GLN
1	B	255	ASN
1	B	261	GLN
1	B	282	ASN
1	B	291	GLN
1	B	342	HIS
1	B	418	HIS
1	B	473	HIS
1	B	485	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

9 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ACY	A	907	-	3,3,3	0.92	0	3,3,3	0.80	0
3	ACY	B	905	-	3,3,3	0.69	0	3,3,3	1.06	0
2	EDO	A	908	-	3,3,3	0.88	0	2,2,2	0.58	0
3	ACY	B	904	-	3,3,3	1.04	0	3,3,3	0.53	0
3	ACY	A	901	-	3,3,3	0.98	0	3,3,3	0.43	0
3	ACY	B	902	-	3,3,3	0.91	0	3,3,3	1.25	0
3	ACY	A	909	-	3,3,3	0.84	0	3,3,3	0.74	0
3	ACY	A	906	-	3,3,3	1.14	0	3,3,3	0.63	0
3	ACY	A	903	-	3,3,3	1.10	0	3,3,3	0.70	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	A	908	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	908	EDO	O1-C1-C2-O2

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	905	ACY	3	0
3	A	901	ACY	1	0
3	B	902	ACY	5	0
3	A	909	ACY	1	0

5.7 Other polymers [i](#)

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	481/497 (96%)	-0.51	2 (0%) 92 93	9, 16, 26, 41	0
1	B	480/497 (96%)	-0.55	7 (1%) 73 77	8, 17, 28, 47	0
All	All	961/994 (96%)	-0.53	9 (0%) 84 86	8, 17, 27, 47	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	57	PRO	3.7
1	B	55	GLU	2.8
1	B	54	LYS	2.7
1	B	316	ARG	2.6
1	B	249	ASP	2.5
1	A	4	ASP	2.4
1	B	56	ARG	2.1
1	B	52	ARG	2.1
1	A	110	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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	B-factors(Å ²)	Q<0.9
2	EDO	A	908	4/4	0.89	0.12	25,25,26,27	0
3	ACY	A	909	4/4	0.92	0.14	35,35,35,35	0
3	ACY	B	905	4/4	0.92	0.13	13,16,17,20	0
3	ACY	B	904	4/4	0.94	0.13	17,19,19,19	0
3	ACY	B	902	4/4	0.94	0.10	17,19,20,22	0
3	ACY	A	907	4/4	0.97	0.13	24,24,25,25	0
3	ACY	A	906	4/4	0.98	0.11	15,15,15,16	0
3	ACY	A	901	4/4	0.98	0.09	13,14,15,15	0
3	ACY	A	903	4/4	0.98	0.06	14,15,15,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.