



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 06:30 pm BST

PDB ID : 4I26
Title : 2.20 Angstroms X-ray crystal structure of 2-aminomuconate 6-semialdehyde dehydrogenase from *Pseudomonas fluorescens*
Authors : Davis, I.; Huo, L.; Chen, L.; Liu, A.
Deposited on : 2012-11-21
Resolution : 2.20 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

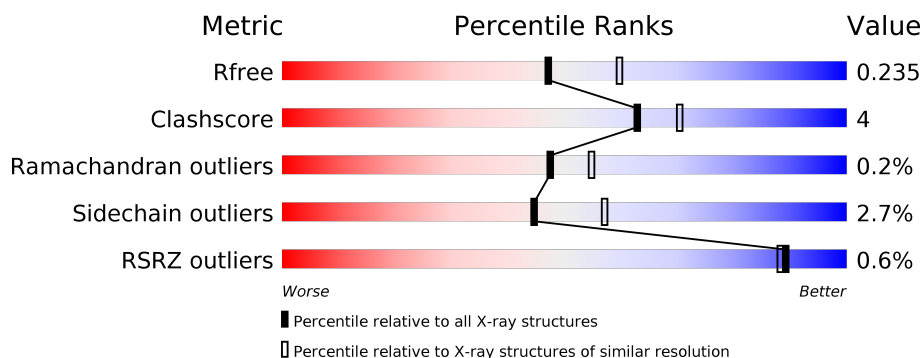
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 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	500	<div> <div>86%</div> <div>9%</div> <div>• •</div> </div>
1	B	500	<div> <div>%</div> <div>85%</div> <div>11%</div> <div>•</div> </div>
1	C	500	<div> <div>%</div> <div>86%</div> <div>10%</div> <div>•</div> </div>
1	D	500	<div> <div>%</div> <div>87%</div> <div>8%</div> <div>• •</div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	EDO	A	601	-	-	X	-
2	EDO	B	602	-	-	X	-

2 Entry composition [i](#)

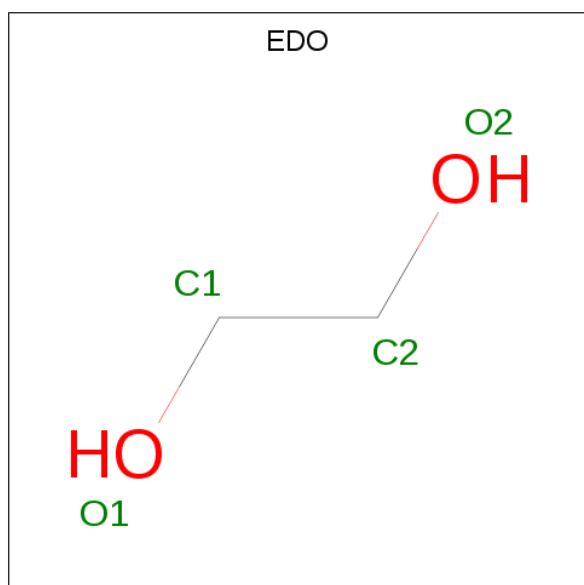
There are 4 unique types of molecules in this entry. The entry contains 15376 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 2-aminomuconate 6-semialdehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	482	Total	C	N	O	S	0	0	0
			3665	2318	636	699	12			
1	B	482	Total	C	N	O	S	0	0	0
			3665	2318	636	699	12			
1	C	481	Total	C	N	O	S	0	0	0
			3656	2313	634	697	12			
1	D	482	Total	C	N	O	S	0	0	0
			3665	2318	636	699	12			

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



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

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

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Na	0	0
			1	1		
3	D	1	Total	Na	0	0
			1	1		
3	C	1	Total	Na	0	0
			1	1		

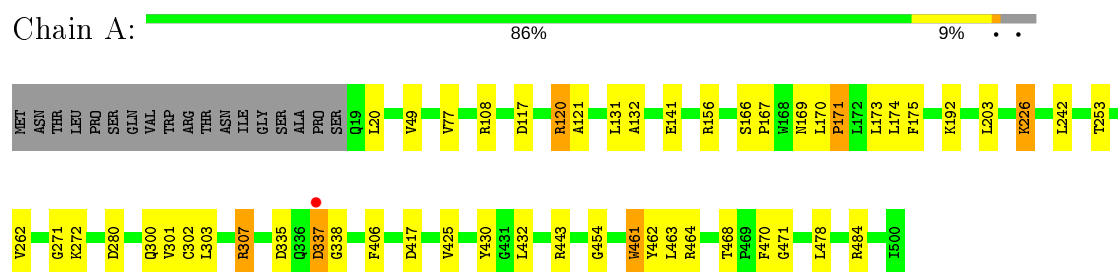
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	175	Total	O	0	0
			175	175		
4	B	168	Total	O	0	0
			168	168		
4	C	149	Total	O	0	0
			149	149		
4	D	210	Total	O	0	0
			210	210		

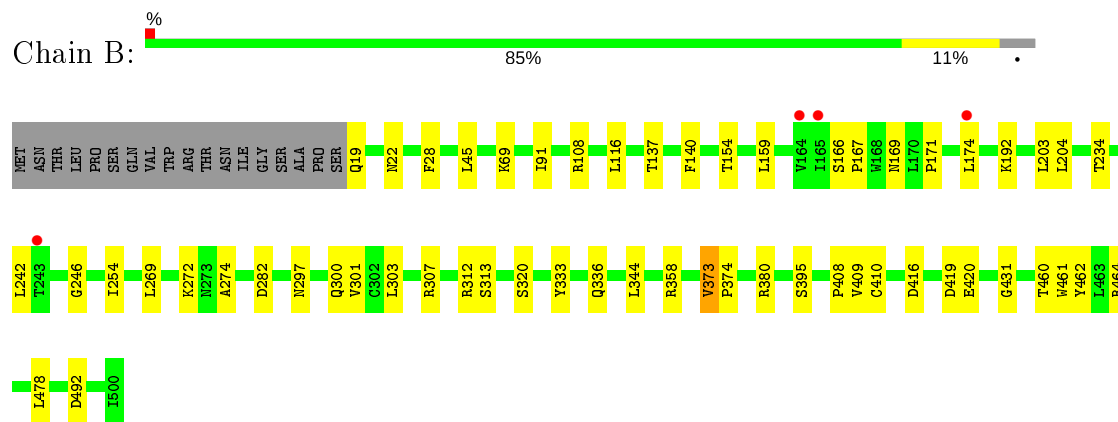
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

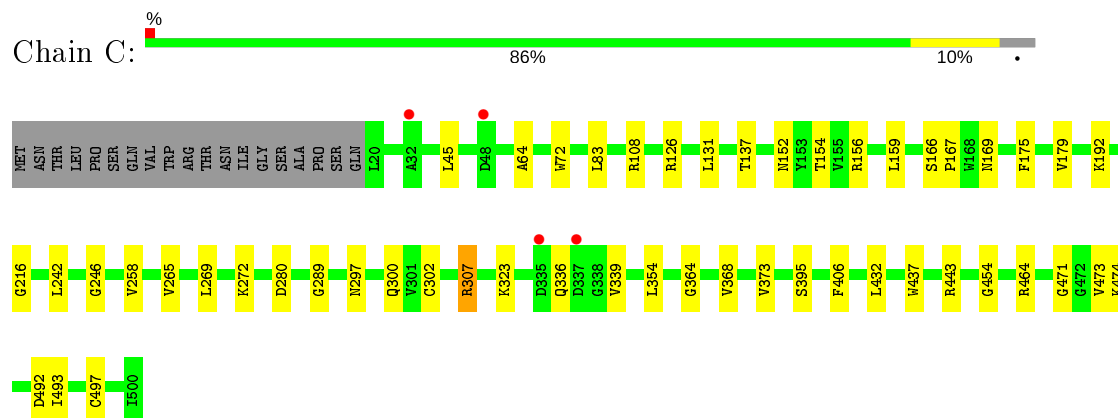
- Molecule 1: 2-aminomuconate 6-semialdehyde dehydrogenase



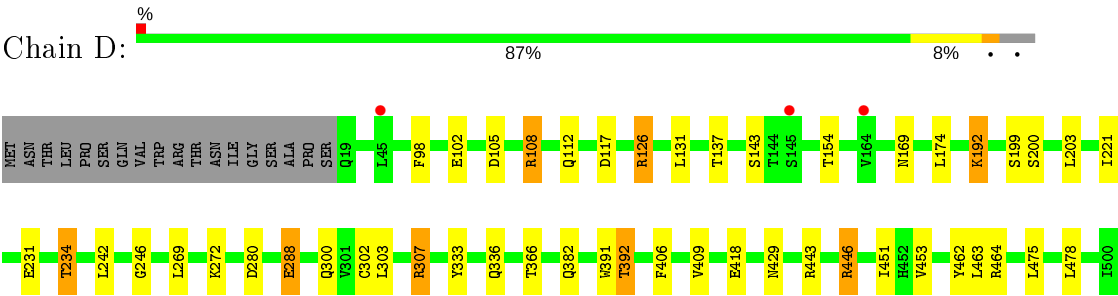
- Molecule 1: 2-aminomuconate 6-semialdehyde dehydrogenase



- Molecule 1: 2-aminomuconate 6-semialdehyde dehydrogenase



● Molecule 1: 2-aminomuconate 6-semialdehyde dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.27Å 141.89Å 172.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.72 – 2.20 31.72 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.6 (31.72-2.20) 96.7 (31.72-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.74 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.7.3 _928	Depositor
R, R_{free}	0.186 , 0.239 0.184 , 0.235	Depositor DCC
R_{free} test set	5472 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	36.9	Xtriage
Anisotropy	0.176	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 39.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15376	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.43% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA, EDO

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	0.65	1/3744 (0.0%)	0.71	3/5090 (0.1%)
1	B	0.64	0/3744	0.74	0/5090
1	C	0.63	0/3735	0.69	1/5078 (0.0%)
1	D	0.67	0/3744	0.75	5/5090 (0.1%)
All	All	0.65	1/14967 (0.0%)	0.72	9/20348 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	121	ALA	CA-CB	5.33	1.63	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	446	ARG	NE-CZ-NH2	-10.89	114.86	120.30
1	D	446	ARG	NE-CZ-NH1	9.21	124.90	120.30
1	D	126	ARG	NE-CZ-NH2	-7.49	116.56	120.30
1	A	120	ARG	NE-CZ-NH1	-6.87	116.87	120.30
1	D	126	ARG	NE-CZ-NH1	6.37	123.48	120.30
1	C	307	ARG	NE-CZ-NH2	-5.55	117.52	120.30
1	A	307	ARG	NE-CZ-NH2	-5.53	117.53	120.30
1	D	307	ARG	NE-CZ-NH2	-5.13	117.73	120.30
1	A	156	ARG	NE-CZ-NH1	-5.08	117.76	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3665	0	3600	43	0
1	B	3665	0	3600	37	0
1	C	3656	0	3592	23	0
1	D	3665	0	3600	24	0
2	A	4	0	6	10	0
2	B	8	0	12	5	0
2	C	4	0	6	1	0
2	D	4	0	6	2	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	175	0	0	3	0
4	B	168	0	0	4	0
4	C	149	0	0	1	0
4	D	210	0	0	3	0
All	All	15376	0	14422	126	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 4.

All (126) 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:C:108:ARG:NH2	1:C:169:ASN:O	2.05	0.89
1:B:19:GLN:OE1	1:B:19:GLN:N	2.10	0.85
1:A:302:CYS:SG	4:A:741:HOH:O	2.35	0.84
1:A:280:ASP:OD2	1:A:443:ARG:NH1	2.13	0.82
1:D:302:CYS:SG	4:D:856:HOH:O	2.40	0.79
1:D:280:ASP:OD2	1:D:443:ARG:NH1	2.20	0.74
1:D:272:LYS:HG3	1:D:307:ARG:HD2	1.68	0.74
1:B:242:LEU:HD23	1:B:254:ILE:CG2	2.17	0.74
1:D:366:THR:HB	1:D:392:THR:HG22	1.70	0.74
1:B:242:LEU:CD2	1:B:254:ILE:HG23	2.18	0.73
1:A:464:ARG:NH1	2:A:601:EDO:C1	2.52	0.72
1:A:464:ARG:NH1	2:A:601:EDO:H12	2.05	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:303:LEU:O	1:A:461:TRP:HZ3	1.73	0.70
1:A:272:LYS:HG3	1:A:307:ARG:HD2	1.72	0.70
1:C:272:LYS:HG3	1:C:307:ARG:HD2	1.76	0.68
1:D:108:ARG:NH1	1:D:112:GLN:OE1	2.27	0.67
1:A:120:ARG:HH12	2:A:601:EDO:H11	1.60	0.67
1:A:337:ASP:CG	1:A:338:GLY:H	1.98	0.66
1:B:312:ARG:HD2	1:B:416:ASP:OD1	1.96	0.65
1:A:478:LEU:HD21	2:B:602:EDO:C2	2.27	0.65
1:A:141:GLU:OE1	4:A:721:HOH:O	2.14	0.64
1:D:418:GLU:OE2	1:D:446:ARG:HD2	1.99	0.63
1:B:358:ARG:NH2	4:B:793:HOH:O	2.32	0.62
1:B:242:LEU:HD23	1:B:254:ILE:HG23	1.81	0.61
1:B:242:LEU:CD2	1:B:254:ILE:CG2	2.78	0.61
1:A:478:LEU:HD21	2:B:602:EDO:H12	1.84	0.60
1:A:272:LYS:HG3	1:A:307:ARG:CD	2.32	0.60
1:A:120:ARG:HH12	2:A:601:EDO:C1	2.14	0.60
1:A:462:TYR:CE1	2:A:601:EDO:H21	2.38	0.58
1:B:272:LYS:HG3	1:B:307:ARG:HD2	1.85	0.58
1:A:272:LYS:CG	1:A:307:ARG:HD2	2.34	0.57
1:A:478:LEU:HD21	2:B:602:EDO:C1	2.36	0.56
1:D:108:ARG:NH1	1:D:117:ASP:OD1	2.39	0.55
1:A:173:LEU:C	1:A:173:LEU:HD23	2.25	0.55
1:D:464:ARG:HD3	2:D:601:EDO:H12	1.89	0.55
1:C:169:ASN:HB2	1:C:300:GLN:O	2.06	0.55
1:A:462:TYR:OH	2:A:601:EDO:H21	2.07	0.54
1:A:464:ARG:NH1	1:A:470:PHE:CD2	2.76	0.53
1:D:174:LEU:HD21	2:D:601:EDO:H22	1.91	0.53
1:D:382:GLN:NE2	4:D:776:HOH:O	2.35	0.52
1:C:323:LYS:HE2	1:C:368:VAL:O	2.10	0.52
1:A:464:ARG:NH2	4:A:713:HOH:O	2.43	0.52
1:D:231:GLU:O	1:D:234:THR:HG22	2.09	0.52
1:B:464:ARG:HE	2:B:601:EDO:H12	1.75	0.51
1:C:137:THR:HB	1:C:154:THR:OG1	2.11	0.51
1:B:344:LEU:HD21	1:B:408:PRO:HD3	1.93	0.50
1:D:137:THR:HB	1:D:154:THR:OG1	2.12	0.49
1:D:429:ASN:HB3	1:D:475:LEU:HD12	1.95	0.49
1:C:364:GLY:O	4:C:844:HOH:O	2.20	0.48
1:B:108:ARG:HH21	1:B:116:LEU:HD23	1.79	0.48
1:B:242:LEU:HD23	1:B:254:ILE:HG21	1.93	0.48
1:C:454:GLY:HA3	1:C:471:GLY:O	2.14	0.48
1:C:108:ARG:NE	1:C:297:ASN:OD1	2.39	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:280:ASP:OD2	1:C:443:ARG:NH1	2.48	0.47
1:C:152:ASN:OD1	1:C:497:CYS:HB2	2.15	0.47
1:A:335:ASP:N	1:A:335:ASP:OD1	2.48	0.46
1:D:169:ASN:HB2	1:D:300:GLN:O	2.15	0.46
1:A:337:ASP:CG	1:A:338:GLY:N	2.67	0.46
1:B:169:ASN:HB2	1:B:300:GLN:O	2.14	0.46
1:A:166:SER:HB2	1:A:167:PRO:HD2	1.97	0.46
1:A:468:THR:HG22	1:B:140:PHE:CD1	2.51	0.46
1:A:108:ARG:NH1	1:A:117:ASP:OD2	2.49	0.46
1:A:226:LYS:HE3	1:A:253:THR:HG21	1.97	0.46
1:D:451:ILE:HG22	1:D:453:VAL:HG13	1.98	0.46
1:A:169:ASN:HB2	1:A:300:GLN:O	2.15	0.46
1:C:64:ALA:HA	1:C:216:GLY:O	2.16	0.46
1:B:159:LEU:HD12	1:B:492:ASP:OD1	2.16	0.45
1:C:464:ARG:HD3	2:C:601:EDO:O2	2.15	0.45
1:D:333:TYR:HB2	1:D:336:GLN:HB2	1.98	0.45
1:B:108:ARG:NH2	1:B:116:LEU:HD23	2.32	0.45
1:B:19:GLN:CB	4:B:867:HOH:O	2.63	0.45
1:D:105:ASP:OD1	1:D:199:SER:OG	2.33	0.45
1:D:246:GLY:O	1:D:269:LEU:HA	2.17	0.45
1:A:77:VAL:HG11	1:A:132:ALA:O	2.17	0.45
1:A:337:ASP:OD1	1:A:338:GLY:N	2.50	0.44
1:B:312:ARG:CG	1:B:313:SER:N	2.80	0.44
1:B:108:ARG:HD2	1:B:297:ASN:OD1	2.18	0.44
1:A:303:LEU:C	1:A:461:TRP:HZ3	2.21	0.44
1:C:159:LEU:HD12	1:C:492:ASP:OD1	2.18	0.44
1:A:464:ARG:HH11	2:A:601:EDO:H22	1.83	0.44
1:A:120:ARG:NH1	2:A:601:EDO:H11	2.30	0.43
1:C:72:TRP:HZ3	1:C:83:LEU:HD23	1.83	0.43
1:D:288:GLU:HG2	4:D:822:HOH:O	2.18	0.43
1:B:301:VAL:HG12	1:B:303:LEU:H	1.81	0.43
1:B:274:ALA:HA	1:B:307:ARG:O	2.18	0.43
1:A:462:TYR:CZ	2:A:601:EDO:H21	2.53	0.43
1:C:156:ARG:HG2	1:C:493:ILE:CD1	2.49	0.43
1:A:462:TYR:HE1	2:A:601:EDO:H21	1.82	0.43
1:A:20:LEU:HD12	1:A:49:VAL:HG22	2.01	0.43
1:B:137:THR:HB	1:B:154:THR:OG1	2.18	0.43
1:B:174:LEU:HA	1:B:174:LEU:HD23	1.86	0.43
1:C:166:SER:HB2	1:C:167:PRO:HD2	2.01	0.43
1:B:419:ASP:OD1	1:B:420:GLU:N	2.53	0.42
1:C:289:GLY:HA3	1:C:437:TRP:CZ2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:91:ILE:HD11	1:B:204:LEU:HD13	2.00	0.42
1:C:175:PHE:O	1:C:179:VAL:HG23	2.19	0.42
1:D:192:LYS:HA	1:D:221:ILE:O	2.19	0.42
1:B:478:LEU:HD21	2:B:602:EDO:H11	2.00	0.42
1:A:271:GLY:HA2	1:A:430:TYR:HB3	2.02	0.42
1:B:166:SER:HB2	1:B:167:PRO:HD2	2.02	0.42
1:B:19:GLN:HG3	4:B:867:HOH:O	2.20	0.42
1:B:373:VAL:HG23	1:B:374:PRO:O	2.20	0.41
1:B:460:THR:OG1	1:B:461:TRP:N	2.54	0.41
1:C:473:VAL:HG23	1:C:474:LYS:HG3	2.02	0.41
1:A:262:VAL:O	1:A:262:VAL:HG12	2.20	0.41
1:A:170:LEU:HD11	1:A:301:VAL:HG13	2.03	0.41
1:A:454:GLY:HA3	1:A:471:GLY:O	2.20	0.41
1:B:282:ASP:OD1	1:B:282:ASP:C	2.59	0.41
1:B:19:GLN:HB3	4:B:867:HOH:O	2.21	0.41
1:C:336:GLN:HB3	1:C:339:VAL:HB	2.02	0.41
1:D:391:TRP:CH2	1:D:409:VAL:HG21	2.56	0.41
1:A:174:LEU:HD23	1:A:174:LEU:HA	1.94	0.41
1:A:307:ARG:NH2	1:A:425:VAL:HA	2.36	0.41
1:B:22:ASN:O	1:B:28:PHE:HA	2.21	0.41
1:B:246:GLY:O	1:B:269:LEU:HA	2.20	0.41
1:C:246:GLY:O	1:C:269:LEU:HA	2.21	0.41
1:C:354:LEU:HA	1:C:354:LEU:HD23	1.91	0.41
1:D:98:PHE:CZ	1:D:203:LEU:HB3	2.56	0.41
1:B:333:TYR:HB2	1:B:336:GLN:HB2	2.03	0.40
1:B:409:VAL:HG22	1:B:410:CYS:N	2.36	0.40
1:D:102:GLU:OE2	1:D:200:SER:OG	2.28	0.40
1:B:333:TYR:CE2	1:B:380:ARG:NE	2.89	0.40
1:C:258:VAL:HG21	1:C:265:VAL:CG1	2.51	0.40
1:D:303:LEU:HD12	1:D:462:TYR:CD1	2.56	0.40
1:A:166:SER:HB2	1:A:167:PRO:CD	2.51	0.40
1:A:171:PRO:O	1:A:175:PHE:HB2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

of similar resolution.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	480/500 (96%)	458 (95%)	20 (4%)	2 (0%)	34	37
1	B	480/500 (96%)	460 (96%)	18 (4%)	2 (0%)	34	37
1	C	479/500 (96%)	459 (96%)	20 (4%)	0	100	100
1	D	480/500 (96%)	463 (96%)	17 (4%)	0	100	100
All	All	1919/2000 (96%)	1840 (96%)	75 (4%)	4 (0%)	47	55

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	337	ASP
1	B	431	GLY
1	B	171	PRO
1	A	171	PRO

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	391/407 (96%)	380 (97%)	11 (3%)	43	56
1	B	391/407 (96%)	382 (98%)	9 (2%)	50	63
1	C	390/407 (96%)	380 (97%)	10 (3%)	46	58
1	D	391/407 (96%)	379 (97%)	12 (3%)	40	51
All	All	1563/1628 (96%)	1521 (97%)	42 (3%)	44	57

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	131	LEU
1	A	192	LYS

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Mol	Chain	Res	Type
1	A	203	LEU
1	A	226	LYS
1	A	242	LEU
1	A	406	PHE
1	A	417	ASP
1	A	432	LEU
1	A	461	TRP
1	A	463	LEU
1	A	484	ARG
1	B	45	LEU
1	B	69	LYS
1	B	192	LYS
1	B	203	LEU
1	B	234	THR
1	B	320	SER
1	B	373	VAL
1	B	395	SER
1	B	462	TYR
1	C	45	LEU
1	C	126	ARG
1	C	131	LEU
1	C	192	LYS
1	C	242	LEU
1	C	302	CYS
1	C	373	VAL
1	C	395	SER
1	C	406	PHE
1	C	432	LEU
1	D	108	ARG
1	D	126	ARG
1	D	131	LEU
1	D	143	SER
1	D	192	LYS
1	D	234	THR
1	D	242	LEU
1	D	288	GLU
1	D	392	THR
1	D	406	PHE
1	D	463	LEU
1	D	478	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	347	HIS
1	B	19	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 3 are monoatomic - leaving 5 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	$\# Z > 2$
2	EDO	C	601	-	3,3,3	0.47	0	2,2,2	1.03	0
2	EDO	B	602	-	3,3,3	0.90	0	2,2,2	0.22	0
2	EDO	B	601	-	3,3,3	0.85	0	2,2,2	0.76	0
2	EDO	D	601	-	3,3,3	0.45	0	2,2,2	0.60	0
2	EDO	A	601	-	3,3,3	0.36	0	2,2,2	0.80	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	C	601	-	-	1/1/1/1	-
2	EDO	B	602	-	-	1/1/1/1	-
2	EDO	B	601	-	-	1/1/1/1	-
2	EDO	D	601	-	-	0/1/1/1	-
2	EDO	A	601	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	601	EDO	O1-C1-C2-O2
2	B	602	EDO	O1-C1-C2-O2
2	C	601	EDO	O1-C1-C2-O2

There are no ring outliers.

5 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	601	EDO	1	0
2	B	602	EDO	4	0
2	B	601	EDO	1	0
2	D	601	EDO	2	0
2	A	601	EDO	10	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	482/500 (96%)	-0.41	1 (0%) 95 94	23, 42, 54, 68	0
1	B	482/500 (96%)	-0.35	4 (0%) 86 85	24, 41, 53, 67	0
1	C	481/500 (96%)	-0.31	4 (0%) 86 85	25, 43, 59, 81	0
1	D	482/500 (96%)	-0.43	3 (0%) 89 88	23, 36, 49, 63	0
All	All	1927/2000 (96%)	-0.37	12 (0%) 89 88	23, 40, 55, 81	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	337	ASP	3.9
1	C	337	ASP	3.7
1	C	32	ALA	3.3
1	B	164	VAL	2.8
1	D	164	VAL	2.5
1	C	335	ASP	2.3
1	B	243	THR	2.2
1	B	165	ILE	2.2
1	D	45	LEU	2.1
1	D	145	SER	2.1
1	B	174	LEU	2.0
1	C	48	ASP	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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	B	602	4/4	0.85	0.20	37,39,42,51	0
2	EDO	A	601	4/4	0.87	0.16	32,37,37,44	0
2	EDO	B	601	4/4	0.90	0.19	32,35,38,45	0
2	EDO	C	601	4/4	0.91	0.18	33,37,37,40	0
2	EDO	D	601	4/4	0.93	0.21	29,34,35,44	0
3	NA	C	602	1/1	0.96	0.06	33,33,33,33	0
3	NA	D	602	1/1	0.96	0.08	34,34,34,34	0
3	NA	B	603	1/1	0.98	0.04	37,37,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.