



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:30 PM UTC

PDB ID : 4NYY / pdb_00004nyy
Title : Structure of *Vibrio cholerae* chitin de-N-acetylase in complex with acetate ion (ACT) in P 2 21 21
Authors : Albesa-Jove, D.; Andres, E.; Biarnes, X.; Planas, A.; Guerin, M.E.
Deposited on : 2013-12-11
Resolution : 2.65 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

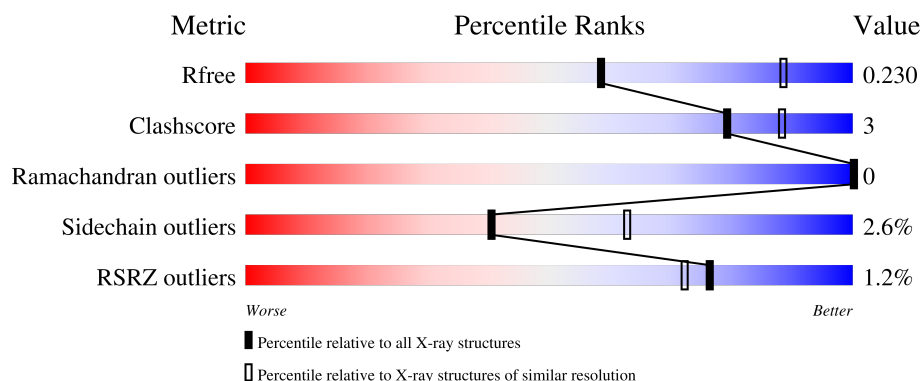
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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}	180053	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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	419	0% 88% 9% .
1	B	419	0% 90% 7% .
1	C	419	0% 88% 8% .
1	D	419	2% 86% 10% .

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 13312 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 Deacetylase DA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	405	Total	C	N	O	S	0	7	0
			3224	2043	540	631	10			
1	B	405	Total	C	N	O	S	0	2	2
			3171	2016	532	613	10			
1	C	404	Total	C	N	O	S	0	5	1
			3178	2018	534	616	10			
1	D	405	Total	C	N	O	S	0	1	0
			3165	2008	533	615	9			

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	26	MET	-	expression tag	UNP M7M1G8
A	432	GLU	-	expression tag	UNP M7M1G8
A	433	LEU	-	expression tag	UNP M7M1G8
A	434	ARG	-	expression tag	UNP M7M1G8
A	435	ARG	-	expression tag	UNP M7M1G8
A	436	GLN	-	expression tag	UNP M7M1G8
A	437	TRP	-	expression tag	UNP M7M1G8
A	438	SER	-	expression tag	UNP M7M1G8
A	439	HIS	-	expression tag	UNP M7M1G8
A	440	PRO	-	expression tag	UNP M7M1G8
A	441	GLN	-	expression tag	UNP M7M1G8
A	442	PHE	-	expression tag	UNP M7M1G8
A	443	GLU	-	expression tag	UNP M7M1G8
A	444	LYS	-	expression tag	UNP M7M1G8
B	26	MET	-	expression tag	UNP M7M1G8
B	432	GLU	-	expression tag	UNP M7M1G8
B	433	LEU	-	expression tag	UNP M7M1G8
B	434	ARG	-	expression tag	UNP M7M1G8
B	435	ARG	-	expression tag	UNP M7M1G8
B	436	GLN	-	expression tag	UNP M7M1G8
B	437	TRP	-	expression tag	UNP M7M1G8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	438	SER	-	expression tag	UNP M7M1G8
B	439	HIS	-	expression tag	UNP M7M1G8
B	440	PRO	-	expression tag	UNP M7M1G8
B	441	GLN	-	expression tag	UNP M7M1G8
B	442	PHE	-	expression tag	UNP M7M1G8
B	443	GLU	-	expression tag	UNP M7M1G8
B	444	LYS	-	expression tag	UNP M7M1G8
C	26	MET	-	expression tag	UNP M7M1G8
C	432	GLU	-	expression tag	UNP M7M1G8
C	433	LEU	-	expression tag	UNP M7M1G8
C	434	ARG	-	expression tag	UNP M7M1G8
C	435	ARG	-	expression tag	UNP M7M1G8
C	436	GLN	-	expression tag	UNP M7M1G8
C	437	TRP	-	expression tag	UNP M7M1G8
C	438	SER	-	expression tag	UNP M7M1G8
C	439	HIS	-	expression tag	UNP M7M1G8
C	440	PRO	-	expression tag	UNP M7M1G8
C	441	GLN	-	expression tag	UNP M7M1G8
C	442	PHE	-	expression tag	UNP M7M1G8
C	443	GLU	-	expression tag	UNP M7M1G8
C	444	LYS	-	expression tag	UNP M7M1G8
D	26	MET	-	expression tag	UNP M7M1G8
D	432	GLU	-	expression tag	UNP M7M1G8
D	433	LEU	-	expression tag	UNP M7M1G8
D	434	ARG	-	expression tag	UNP M7M1G8
D	435	ARG	-	expression tag	UNP M7M1G8
D	436	GLN	-	expression tag	UNP M7M1G8
D	437	TRP	-	expression tag	UNP M7M1G8
D	438	SER	-	expression tag	UNP M7M1G8
D	439	HIS	-	expression tag	UNP M7M1G8
D	440	PRO	-	expression tag	UNP M7M1G8
D	441	GLN	-	expression tag	UNP M7M1G8
D	442	PHE	-	expression tag	UNP M7M1G8
D	443	GLU	-	expression tag	UNP M7M1G8
D	444	LYS	-	expression tag	UNP M7M1G8

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0

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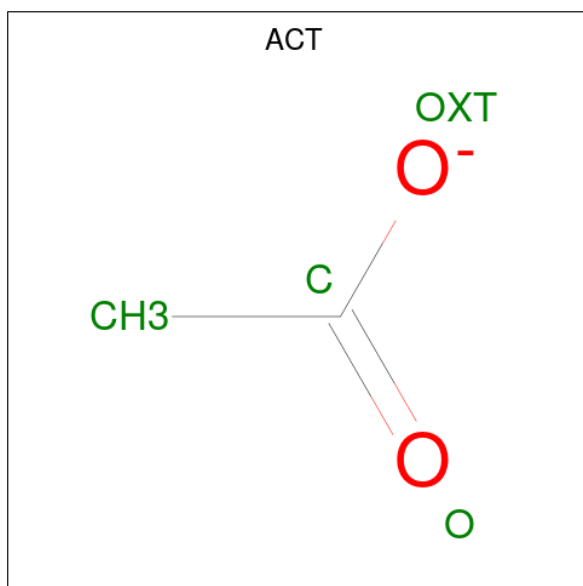
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

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

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



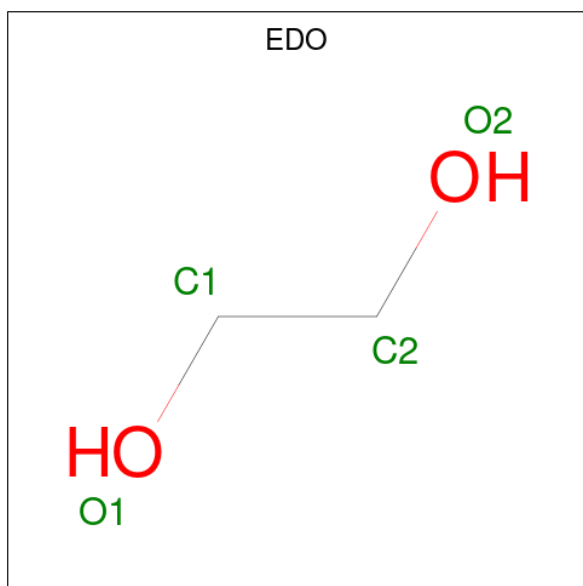
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		

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

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	172	Total	O	0	0
			172	172		
6	B	152	Total	O	0	0
			152	152		
6	C	99	Total	O	0	0
			99	99		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	107	Total 107	O 107	0	0

- Molecule 1: Deacetylase DA1





4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.17Å 165.01Å 203.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.15 – 2.65 49.15 – 2.65	Depositor EDS
% Data completeness (in resolution range)	98.7 (49.15-2.65) 98.7 (49.15-2.65)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.32 (at 2.65Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.202 , 0.227 0.208 , 0.230	Depositor DCC
R_{free} test set	3148 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	43.4	Xtriage
Anisotropy	0.577	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 29.8	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	0.94	EDS
Total number of atoms	13312	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.94% 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: EDO, ACT, NA, ZN

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	0.28	0/3323	0.64	0/4547
1	B	0.29	0/3267	0.67	0/4472
1	C	0.28	0/3278	0.66	2/4484 (0.0%)
1	D	0.28	0/3259	0.65	1/4461 (0.0%)
All	All	0.28	0/13127	0.65	3/17964 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	140	GLU	N-CA-C	-5.22	105.27	111.69
1	C	150	LEU	CA-C-N	5.07	124.73	119.56
1	C	150	LEU	C-N-CA	5.07	124.73	119.56

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	3224	0	3026	18	0
1	B	3171	0	2987	14	0
1	C	3178	0	3000	15	0
1	D	3165	0	2975	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	4	0	3	0	0
4	B	4	0	3	0	0
4	C	4	0	3	0	0
4	D	4	0	3	0	0
5	A	8	0	12	0	0
5	B	8	0	12	0	0
5	C	4	0	6	0	0
6	A	172	0	0	1	0
6	B	152	0	0	0	0
6	C	99	0	0	0	0
6	D	107	0	0	0	0
All	All	13312	0	12030	68	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 3.

All (68) 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:D:331:THR:OG1	1:D:333:ASP:OD1	2.08	0.70
1:A:59:LYS:NZ	6:A:771:HOH:O	2.26	0.64
1:B:240:ILE:HB	1:B:241:PRO:HD2	1.86	0.56
1:A:265[B]:SER:N	1:A:266:PRO:CD	2.70	0.54
1:D:240:ILE:HB	1:D:241:PRO:CD	2.38	0.53
1:A:265[A]:SER:N	1:A:266:PRO:CD	2.72	0.52
1:B:312:LEU:N	1:B:313:PRO:CD	2.75	0.50
1:A:97:HIS:O	1:A:98:SER:OG	2.30	0.50
1:D:285:LEU:HD11	1:D:350:TYR:HB2	1.93	0.50
1:C:331:THR:OG1	1:C:333:ASP:OD1	2.28	0.49
1:D:45:ALA:HB2	1:D:300:GLU:HB2	1.95	0.49
1:A:179:PHE:C	1:A:179:PHE:CD1	2.91	0.49
1:C:97:HIS:O	1:C:98:SER:OG	2.30	0.48
1:C:39:ASP:O	1:C:40:ASP:HB2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:THR:OG1	1:A:333:ASP:OD1	2.25	0.47
1:C:312:LEU:N	1:C:313:PRO:CD	2.78	0.47
1:A:144:ILE:HD11	1:B:140:GLU:HG2	1.97	0.47
1:C:307:GLY:O	1:C:311:ASN:HB2	2.15	0.47
1:C:179:PHE:CD1	1:C:179:PHE:C	2.92	0.46
1:D:265:SER:N	1:D:266:PRO:CD	2.78	0.46
1:B:179:PHE:C	1:B:179:PHE:CD1	2.91	0.46
1:D:39:ASP:O	1:D:40:ASP:HB2	2.15	0.46
1:D:312:LEU:N	1:D:313:PRO:CD	2.79	0.46
1:A:312:LEU:N	1:A:313:PRO:CD	2.78	0.46
1:A:339:TRP:CZ2	1:A:366:ALA:HB1	2.51	0.46
1:B:39:ASP:O	1:B:40:ASP:HB2	2.15	0.46
1:C:251:VAL:N	1:C:252:PRO:HD2	2.31	0.45
1:A:62:PHE:HB3	1:A:64:PHE:CE1	2.52	0.45
1:D:307:GLY:O	1:D:311:ASN:HB2	2.16	0.45
1:D:165:ALA:O	1:D:226:THR:HA	2.18	0.44
1:D:179:PHE:C	1:D:179:PHE:CD1	2.92	0.44
1:A:402:TYR:CE2	1:A:403:LYS:HG3	2.53	0.44
1:C:174:ARG:HD3	1:C:180:GLN:HB3	1.99	0.44
1:A:45:ALA:HB2	1:A:300:GLU:HB2	1.99	0.44
1:A:39:ASP:O	1:A:40:ASP:HB2	2.16	0.44
1:D:97:HIS:O	1:D:98:SER:OG	2.29	0.44
1:B:98:SER:HA	1:B:139:PHE:HA	2.00	0.44
1:D:191:LEU:CD1	1:D:191:LEU:N	2.81	0.44
1:C:62:PHE:HB3	1:C:64:PHE:CE1	2.52	0.44
1:D:251:VAL:N	1:D:252:PRO:HD2	2.33	0.43
1:D:421:PRO:HA	1:D:428:PHE:CD2	2.53	0.43
1:B:171:ASN:OD1	1:B:228:GLY:HA3	2.19	0.43
1:B:307:GLY:O	1:B:311:ASN:HB2	2.18	0.43
1:D:242:MET:HE3	1:D:305:GLY:HA2	1.99	0.43
1:C:29:THR:HA	1:C:30:PRO:HD3	1.82	0.43
1:A:307:GLY:O	1:A:311:ASN:HB2	2.19	0.43
1:A:193:PRO:HA	1:A:198:TYR:CG	2.54	0.43
1:B:331:THR:OG1	1:B:333:ASP:OD1	2.19	0.43
1:D:234:ALA:HB1	1:D:235:PRO:HD2	2.00	0.42
1:A:192:LYS:HD3	1:A:194:TRP:CZ2	2.53	0.42
1:D:190:ASN:HB3	1:D:191:LEU:HD13	2.00	0.42
1:A:251:VAL:N	1:A:252:PRO:HD2	2.35	0.42
1:B:97:HIS:O	1:B:98:SER:OG	2.31	0.42
1:A:144:ILE:CG1	1:B:140:GLU:HG2	2.50	0.42
1:C:265:SER:N	1:C:266:PRO:CD	2.83	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:31:LYS:HE2	1:D:327:TYR:OH	2.21	0.41
1:B:96:ASN:HB2	1:B:161:GLY:O	2.21	0.41
1:B:251:VAL:N	1:B:252:PRO:HD2	2.35	0.41
1:B:265:SER:N	1:B:266:PRO:CD	2.83	0.41
1:D:233:TRP:HB2	1:D:253:PHE:CZ	2.56	0.41
1:D:45:ALA:HB2	1:D:300:GLU:CB	2.51	0.41
1:D:229:TRP:CD2	1:D:293:LEU:HD22	2.56	0.41
1:D:242:MET:CE	1:D:305:GLY:HA2	2.49	0.41
1:C:312:LEU:HB2	1:C:313:PRO:HD3	2.03	0.41
1:D:97:HIS:O	1:D:98:SER:CB	2.67	0.41
1:C:234:ALA:HB1	1:C:235:PRO:HD2	2.02	0.41
1:C:126:ILE:HG22	1:C:127:HIS:CD2	2.57	0.40
1:C:265:SER:HB2	1:C:266:PRO:HD3	2.03	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	410/419 (98%)	394 (96%)	16 (4%)	0	100	100
1	B	405/419 (97%)	392 (97%)	13 (3%)	0	100	100
1	C	407/419 (97%)	396 (97%)	11 (3%)	0	100	100
1	D	404/419 (96%)	389 (96%)	15 (4%)	0	100	100
All	All	1626/1676 (97%)	1571 (97%)	55 (3%)	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	348/358 (97%)	337 (97%)	11 (3%)	34	55
1	B	339/358 (95%)	332 (98%)	7 (2%)	47	69
1	C	341/358 (95%)	328 (96%)	13 (4%)	29	49
1	D	338/358 (94%)	330 (98%)	8 (2%)	43	65
All	All	1366/1432 (95%)	1327 (97%)	39 (3%)	40	58

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	108	GLU
1	A	140[A]	GLU
1	A	140[B]	GLU
1	A	180	GLN
1	A	200	CYS
1	A	272	ILE
1	A	285	LEU
1	A	300	GLU
1	A	340[A]	SER
1	A	340[B]	SER
1	A	425	ASN
1	B	140	GLU
1	B	180	GLN
1	B	200	CYS
1	B	272	ILE
1	B	285	LEU
1	B	300	GLU
1	B	344	THR
1	C	103	ILE
1	C	109[A]	GLU
1	C	109[B]	GLU
1	C	180	GLN
1	C	200	CYS
1	C	272	ILE

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Mol	Chain	Res	Type
1	C	285	LEU
1	C	300	GLU
1	C	310[A]	GLN
1	C	310[B]	GLN
1	C	341	VAL
1	C	389	THR
1	C	425	ASN
1	D	108	GLU
1	D	180	GLN
1	D	191	LEU
1	D	200	CYS
1	D	272	ILE
1	D	300	GLU
1	D	344	THR
1	D	417	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	368	GLN
1	B	354	GLN
1	B	368	GLN
1	C	125	GLN
1	C	354	GLN
1	D	125	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 17 ligands modelled in this entry, 8 are monoatomic - leaving 9 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	EDO	C	504	-	3,3,3	0.43	0	2,2,2	0.32	0
5	EDO	B	504	-	3,3,3	0.45	0	2,2,2	0.33	0
5	EDO	A	504	-	3,3,3	0.43	0	2,2,2	0.35	0
5	EDO	B	505	-	3,3,3	0.44	0	2,2,2	0.35	0
4	ACT	B	503	2	3,3,3	0.78	0	3,3,3	1.42	0
4	ACT	A	503	2	3,3,3	0.79	0	3,3,3	1.39	0
4	ACT	C	503	2	3,3,3	0.82	0	3,3,3	1.36	0
4	ACT	D	503	2	3,3,3	0.79	0	3,3,3	1.36	0
5	EDO	A	505	-	3,3,3	0.42	0	2,2,2	0.40	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
5	EDO	C	504	-	-	1/1/1/1	-
5	EDO	B	504	-	-	0/1/1/1	-
5	EDO	A	504	-	-	0/1/1/1	-
5	EDO	B	505	-	-	1/1/1/1	-
5	EDO	A	505	-	-	1/1/1/1	-

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
5	A	505	EDO	O1-C1-C2-O2
5	B	505	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
5	C	504	EDO	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

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	405/419 (96%)	-0.47	4 (0%) 79 76	1, 14, 41, 76	7 (1%)
1	B	405/419 (96%)	-0.41	3 (0%) 84 82	2, 19, 39, 73	2 (0%)
1	C	404/419 (96%)	-0.25	5 (1%) 76 72	5, 23, 50, 98	5 (1%)
1	D	405/419 (96%)	-0.05	8 (1%) 65 59	4, 25, 62, 92	1 (0%)
All	All	1619/1676 (96%)	-0.29	20 (1%) 76 72	1, 20, 51, 98	15 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	241	PRO	3.5
1	A	433	LEU	3.5
1	C	238	TRP	3.2
1	D	194	TRP	3.1
1	D	240	ILE	2.7
1	C	194	TRP	2.6
1	A	29	THR	2.5
1	C	432	GLU	2.3
1	D	238	TRP	2.3
1	A	238	TRP	2.3
1	D	29	THR	2.2
1	B	190	ASN	2.2
1	D	306	MET	2.2
1	D	246	SER	2.2
1	C	29	THR	2.1
1	A	200	CYS	2.1
1	B	310	GLN	2.1
1	C	239	GLY	2.1
1	B	29	THR	2.1
1	D	243	PRO	2.1

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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
4	ACT	B	503	4/4	0.81	0.16	32,32,32,32	0
5	EDO	B	504	4/4	0.81	0.17	44,44,44,44	0
3	NA	D	502	1/1	0.84	0.22	26,26,26,26	0
5	EDO	B	505	4/4	0.84	0.17	42,42,42,42	0
5	EDO	A	505	4/4	0.86	0.13	31,31,31,31	0
2	ZN	D	501	1/1	0.87	0.11	88,88,88,88	0
4	ACT	D	503	4/4	0.88	0.12	43,43,43,43	0
2	ZN	C	501	1/1	0.88	0.11	80,80,80,80	0
3	NA	B	502	1/1	0.88	0.13	24,24,24,24	0
4	ACT	C	503	4/4	0.88	0.14	49,49,49,49	0
5	EDO	C	504	4/4	0.88	0.17	47,47,47,47	0
4	ACT	A	503	4/4	0.89	0.12	25,25,25,25	0
5	EDO	A	504	4/4	0.91	0.10	24,24,24,24	0
2	ZN	B	501	1/1	0.92	0.09	77,77,77,77	0
3	NA	C	502	1/1	0.94	0.04	17,17,17,17	0
2	ZN	A	501	1/1	0.97	0.05	68,68,68,68	0
3	NA	A	502	1/1	0.99	0.06	16,16,16,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.