



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:13 AM UTC

PDB ID : 1E4D / pdb_00001e4d
Title : Structure of OXA10 beta-lactamase at pH 8.3
Authors : Maveyraud, L.; Golemi, D.; Kotra, L.P.; Tranier, S.; Vakulenko, S.; Mobashery, S.; Samama, J.P.
Deposited on : 2000-07-03
Resolution : 1.80 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

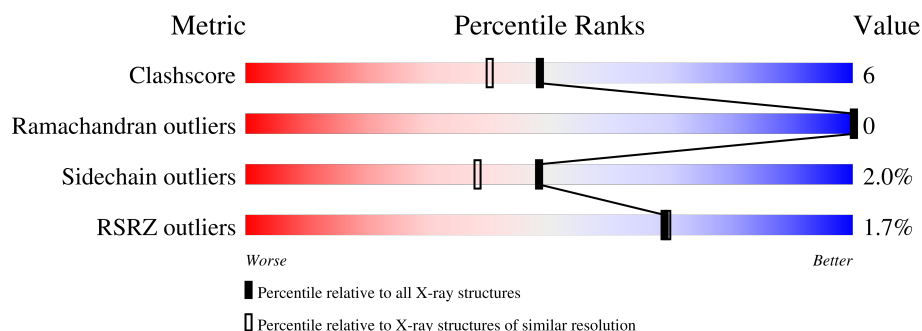
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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))
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	246	% 85% 14% .
1	B	246	87% 11% ..
1	C	246	4% 79% 18% .
1	D	246	% 84% 15% .

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
2	SO4	C	902	-	-	X	-
3	EDO	B	912	-	X	-	-
3	EDO	C	911	-	X	-	-
3	EDO	D	911	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8883 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 BETA-LACTAMASE OXA-10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	243	Total	C	N	O	S	0	3	0
			1953	1248	329	369	7			
1	B	244	Total	C	N	O	S	0	3	0
			1952	1246	327	373	6			
1	C	246	Total	C	N	O	S	0	3	0
			1961	1250	329	376	6			
1	D	246	Total	C	N	O	S	0	4	0
			1970	1255	331	378	6			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	70	KCX	LYS	modified residue	UNP P14489
B	70	KCX	LYS	modified residue	UNP P14489
C	70	KCX	LYS	modified residue	UNP P14489
D	70	KCX	LYS	modified residue	UNP P14489

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	225	Total	O	0	0
			225	225		

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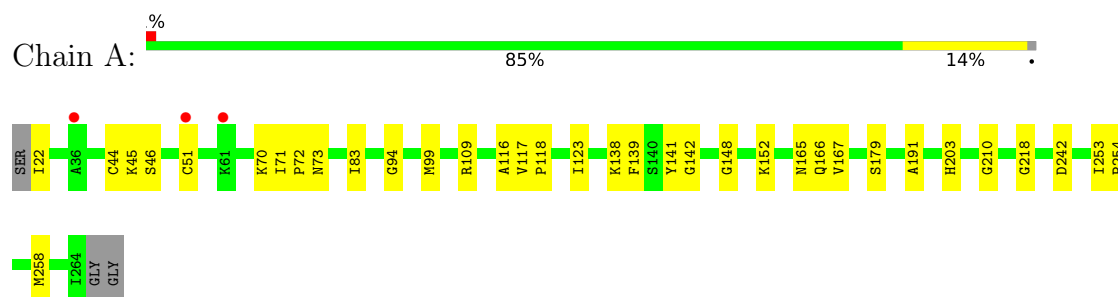
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	250	Total 250	O 250	0	0
4	C	219	Total 219	O 219	0	0
4	D	249	Total 249	O 249	0	0

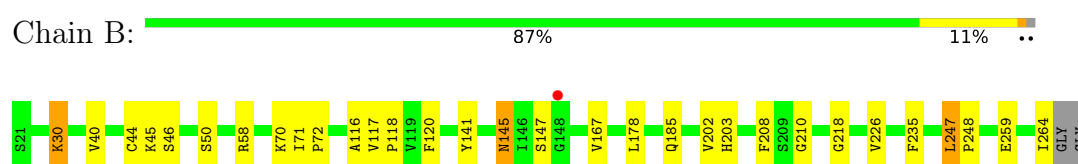
3 Residue-property plots [i](#)

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.

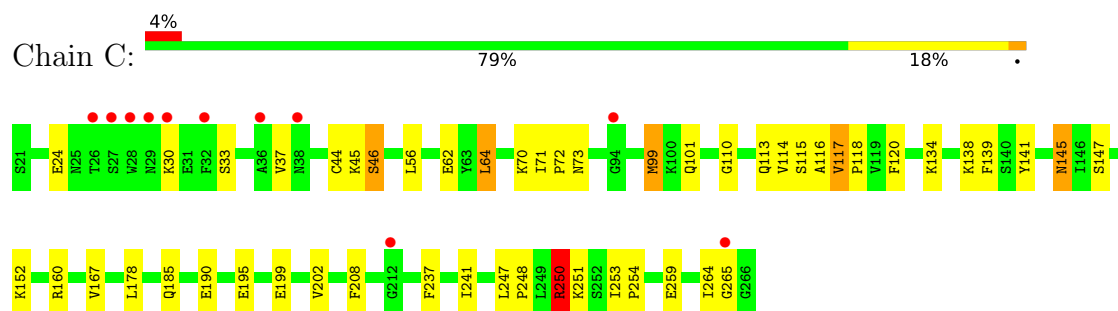
• Molecule 1: BETA-LACTAMASE OXA-10



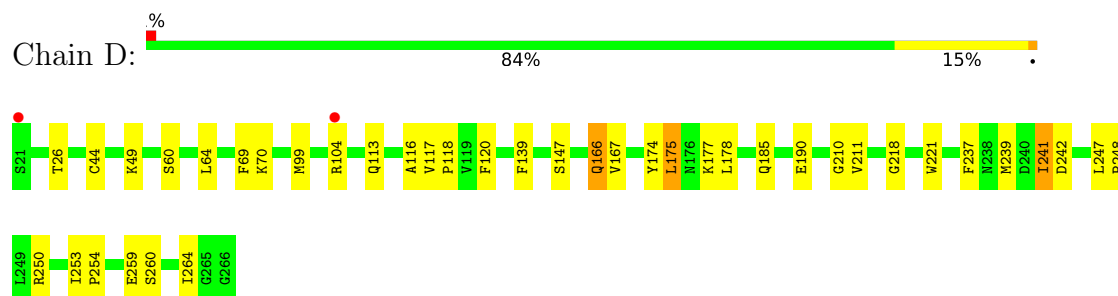
• Molecule 1: BETA-LACTAMASE OXA-10



• Molecule 1: BETA-LACTAMASE OXA-10



• Molecule 1: BETA-LACTAMASE OXA-10



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.28Å 82.44Å 101.17Å 90.00° 95.89° 90.00°	Depositor
Resolution (Å)	29.54 – 1.80 29.54 – 1.80	Depositor EDS
% Data completeness (in resolution range)	84.3 (29.54-1.80) 84.4 (29.54-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 1.80Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.198 , 0.243 0.193 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	20.3	Xtriage
Anisotropy	0.402	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 46.2	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	0.95	EDS
Total number of atoms	8883	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.38% 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, KCX, SO4

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	0.35	0/1981	0.95	4/2675 (0.1%)
1	B	0.38	0/1980	0.98	5/2675 (0.2%)
1	C	0.39	0/1989	1.04	7/2685 (0.3%)
1	D	0.40	0/1998	1.04	7/2697 (0.3%)
All	All	0.38	0/7948	1.01	23/10732 (0.2%)

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
1	C	0	1
1	D	0	3
All	All	0	5

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	250	ARG	CD-NE-CZ	7.93	135.50	124.40
1	C	139	PHE	CA-CB-CG	7.35	121.15	113.80
1	D	104	ARG	CA-CB-CG	6.47	127.03	114.10
1	B	235	PHE	CA-CB-CG	6.41	120.21	113.80
1	D	139	PHE	CA-CB-CG	6.18	119.98	113.80
1	D	120	PHE	CA-CB-CG	-6.05	107.75	113.80
1	D	242	ASP	CA-CB-CG	5.87	118.47	112.60
1	C	250	ARG	NE-CZ-NH1	5.66	127.16	121.50
1	C	116	ALA	CA-C-N	5.57	125.78	120.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	116	ALA	C-N-CA	5.57	125.78	120.43
1	B	208	PHE	CA-CB-CG	5.49	119.29	113.80
1	D	69	PHE	CA-CB-CG	-5.44	108.36	113.80
1	B	116	ALA	CA-C-N	5.41	125.62	120.43
1	B	116	ALA	C-N-CA	5.41	125.62	120.43
1	A	139	PHE	CA-CB-CG	5.34	119.14	113.80
1	D	116	ALA	CA-C-N	5.33	125.55	120.43
1	D	116	ALA	C-N-CA	5.33	125.55	120.43
1	A	116	ALA	CA-C-N	5.33	125.55	120.43
1	A	116	ALA	C-N-CA	5.33	125.55	120.43
1	A	242	ASP	CA-CB-CG	5.32	117.92	112.60
1	C	117	VAL	CB-CA-C	-5.24	108.72	113.70
1	B	120	PHE	CA-CB-CG	-5.04	108.76	113.80
1	C	120	PHE	CA-CB-CG	-5.01	108.79	113.80

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	50[B]	SER	Mainchain
1	C	46[B]	SER	Mainchain
1	D	147[B]	SER	Mainchain
1	D	260[B]	SER	Mainchain
1	D	60[B]	SER	Mainchain

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	1953	0	1944	18	0
1	B	1952	0	1937	19	0
1	C	1961	0	1943	33	0
1	D	1970	0	1950	20	0
2	A	20	0	0	1	0
2	B	25	0	0	0	0
2	C	20	0	0	3	0
2	D	15	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	4	0	6	1	0
3	B	8	0	12	0	0
3	C	4	0	6	1	0
3	D	8	0	12	1	0
4	A	225	0	0	1	0
4	B	250	0	0	1	0
4	C	219	0	0	1	0
4	D	249	0	0	3	0
All	All	8883	0	7810	87	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 (87) 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:259:GLU:HB2	1:D:264:ILE:HD12	1.65	0.79
1:D:113:GLN:HA	3:D:911:EDO:H21	1.73	0.70
1:C:259:GLU:HB2	1:C:264:ILE:HD12	1.78	0.65
1:B:30:LYS:HD3	1:B:30:LYS:N	2.13	0.64
1:B:44:CYS:SG	1:B:167:VAL:HG11	2.39	0.63
1:C:44:CYS:SG	1:C:167:VAL:HG11	2.41	0.60
1:C:145:ASN:C	1:C:145:ASN:HD22	2.11	0.58
1:C:24:GLU:HB2	1:C:56:LEU:HG	1.86	0.58
1:C:248:PRO:HA	2:C:902:SO4:O2	2.04	0.57
1:D:44:CYS:SG	1:D:167:VAL:HG11	2.46	0.56
1:A:99[A]:MET:HE3	2:A:903:SO4:O2	2.08	0.54
1:A:94:GLY:HA3	1:D:99:MET:HE1	1.90	0.54
1:C:145:ASN:HD22	1:C:147:SER:H	1.55	0.53
1:B:145:ASN:HD22	1:B:147[A]:SER:H	1.55	0.53
1:C:253:ILE:HB	1:C:254:PRO:HD3	1.89	0.53
1:D:239:MET:HE2	4:D:2219:HOH:O	2.08	0.53
1:C:71:ILE:HB	1:C:72:PRO:CD	2.40	0.52
1:B:145:ASN:HD22	1:B:147[B]:SER:H	1.55	0.52
1:C:178:LEU:H	1:C:185:GLN:HE22	1.58	0.52
1:B:145:ASN:HD22	1:B:145:ASN:C	2.18	0.52
1:A:44:CYS:SG	1:A:167:VAL:HG11	2.49	0.51
1:B:259:GLU:HG2	1:B:264:ILE:HD12	1.92	0.51
1:A:258:MET:HE3	3:A:911:EDO:H11	1.93	0.51
1:C:99:MET:HB3	1:C:101:GLN:HG2	1.93	0.50
1:C:145:ASN:ND2	1:C:147:SER:H	2.09	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:211:VAL:HG22	1:D:241:ILE:HD11	1.92	0.50
1:C:134:LYS:HE2	1:C:138:LYS:HE3	1.94	0.50
1:B:202:VAL:HG22	1:B:226:VAL:HG22	1.94	0.48
1:D:237:PHE:CE1	1:D:250:ARG:HA	2.49	0.48
1:C:199:GLU:HA	3:C:911:EDO:O2	2.13	0.48
1:D:117:VAL:N	1:D:118:PRO:HD2	2.28	0.48
1:A:138:LYS:HB3	1:A:179:SER:OG	2.13	0.48
1:B:203:HIS:HD2	4:B:2193:HOH:O	1.96	0.48
1:D:49:LYS:HA	1:D:49:LYS:HE2	1.95	0.48
1:D:210:GLY:O	1:D:218:GLY:HA3	2.14	0.47
1:C:250:ARG:HD3	2:C:902:SO4:O1	2.14	0.47
1:B:71:ILE:HB	1:B:72:PRO:CD	2.45	0.47
1:D:211:VAL:HG22	1:D:241:ILE:CD1	2.45	0.47
1:C:237:PHE:CE1	1:C:250:ARG:HA	2.50	0.47
1:B:40:VAL:CG1	1:B:58:ARG:HG2	2.46	0.46
1:C:72:PRO:HG3	1:C:141:TYR:CE2	2.50	0.46
1:C:30:LYS:HA	1:C:33:SER:OG	2.16	0.46
1:C:64:LEU:HD21	4:C:2130:HOH:O	2.15	0.46
1:C:117:VAL:N	1:C:118:PRO:HD2	2.30	0.46
1:D:64:LEU:HD11	4:D:2159:HOH:O	2.16	0.46
1:A:109[B]:ARG:HD2	1:A:191:ALA:HA	1.98	0.45
1:A:210:GLY:O	1:A:218:GLY:HA3	2.16	0.45
1:A:45:LYS:O	1:A:46:SER:HB2	2.17	0.45
1:C:62:GLU:HB3	1:C:160:ARG:HB3	2.00	0.44
1:C:71:ILE:HB	1:C:72:PRO:HD3	2.00	0.44
1:D:253:ILE:HB	1:D:254:PRO:HD3	2.00	0.44
1:A:142:GLY:H	1:A:165:ASN:ND2	2.15	0.44
1:B:247:LEU:N	1:B:248:PRO:HD2	2.32	0.44
1:A:71:ILE:HB	1:A:72:PRO:CD	2.48	0.43
1:B:72:PRO:HG3	1:B:141:TYR:CE2	2.53	0.43
1:A:83:ILE:HD11	1:A:123:ILE:HD13	1.99	0.43
1:A:148:GLY:O	1:A:152:LYS:HE3	2.17	0.43
1:D:26:THR:HG22	4:D:2001:HOH:O	2.17	0.43
1:A:117:VAL:N	1:A:118:PRO:HD2	2.33	0.43
1:A:203:HIS:HD2	4:A:2173:HOH:O	2.00	0.43
1:B:178:LEU:H	1:B:185:GLN:HE22	1.67	0.43
1:C:208:PHE:HE1	1:C:241:ILE:HG21	1.84	0.43
1:C:264:ILE:O	1:C:265:GLY:C	2.60	0.43
1:B:145:ASN:ND2	1:B:147[A]:SER:H	2.16	0.42
1:A:203:HIS:HE1	1:C:190:GLU:OE2	2.02	0.42
1:B:145:ASN:ND2	1:B:147[B]:SER:H	2.17	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:PRO:HG3	1:A:141:TYR:CE2	2.54	0.42
1:C:110:GLY:O	1:C:114:VAL:HG22	2.20	0.42
1:B:45:LYS:O	1:B:46[A]:SER:HB2	2.20	0.42
1:C:247:LEU:HB3	1:C:248:PRO:HD3	2.02	0.42
1:D:178:LEU:H	1:D:185:GLN:HE22	1.68	0.42
1:C:45:LYS:O	1:C:46[B]:SER:HB2	2.18	0.42
1:C:250:ARG:HG2	1:C:251:LYS:N	2.35	0.42
1:C:208:PHE:CE1	1:C:241:ILE:HG21	2.55	0.42
1:D:175:LEU:HB3	1:D:177:LYS:HG3	2.01	0.42
1:D:247:LEU:HB3	1:D:248:PRO:HD3	2.01	0.41
1:B:203:HIS:HE1	1:D:190:GLU:OE2	2.03	0.41
1:A:253:ILE:N	1:A:254:PRO:HD2	2.36	0.41
1:B:210:GLY:O	1:B:218:GLY:HA3	2.20	0.41
1:C:195:GLU:HB2	1:C:202:VAL:HB	2.03	0.41
1:C:208:PHE:CE1	1:C:241:ILE:HD13	2.56	0.41
1:C:237:PHE:CD2	1:C:254:PRO:HG3	2.56	0.41
1:D:166:GLN:HG3	1:D:221:TRP:CZ3	2.56	0.41
1:D:174:TYR:HD2	1:D:175:LEU:HD13	1.86	0.40
1:A:22:ILE:HA	1:A:51:CYS:O	2.21	0.40
1:C:250:ARG:NH2	2:C:904:SO4:O2	2.51	0.40
1:B:117:VAL:N	1:B:118:PRO:HD2	2.36	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	243/246 (99%)	236 (97%)	7 (3%)	0	100	100
1	B	244/246 (99%)	235 (96%)	9 (4%)	0	100	100
1	C	246/246 (100%)	235 (96%)	11 (4%)	0	100	100
1	D	247/246 (100%)	239 (97%)	8 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	980/984 (100%)	945 (96%)	35 (4%)	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	211/209 (101%)	209 (99%)	2 (1%)	70	67
1	B	212/209 (101%)	209 (99%)	3 (1%)	59	52
1	C	212/209 (101%)	203 (96%)	9 (4%)	26	14
1	D	213/209 (102%)	210 (99%)	3 (1%)	59	52
All	All	848/836 (101%)	831 (98%)	17 (2%)	48	38

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

Mol	Chain	Res	Type
1	A	73	ASN
1	A	166	GLN
1	B	30	LYS
1	B	145	ASN
1	B	247	LEU
1	C	37	VAL
1	C	64	LEU
1	C	73	ASN
1	C	99	MET
1	C	113	GLN
1	C	115	SER
1	C	145	ASN
1	C	152	LYS
1	C	250	ARG
1	D	166	GLN
1	D	175	LEU
1	D	241	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	ASN
1	A	165	ASN
1	A	203	HIS
1	B	73	ASN
1	B	122	GLN
1	B	145	ASN
1	B	185	GLN
1	B	203	HIS
1	C	73	ASN
1	C	113	GLN
1	C	133	GLN
1	C	145	ASN
1	C	185	GLN
1	D	113	GLN
1	D	121	GLN
1	D	185	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 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	$\# Z > 2$
1	KCX	D	70	1	10,11,12	2.33	2 (20%)	6,12,14	2.23	2 (33%)
1	KCX	A	70	1	10,11,12	2.73	2 (20%)	6,12,14	5.56	2 (33%)
1	KCX	B	70	1	10,11,12	2.73	2 (20%)	6,12,14	5.13	2 (33%)
1	KCX	C	70	1	10,11,12	2.59	2 (20%)	6,12,14	5.31	2 (33%)

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
1	KCX	D	70	1	-	1/9/10/12	-
1	KCX	A	70	1	-	0/9/10/12	-
1	KCX	B	70	1	-	0/9/10/12	-
1	KCX	C	70	1	-	0/9/10/12	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	70	KCX	CX-NZ	8.06	1.49	1.35
1	B	70	KCX	CX-NZ	8.05	1.49	1.35
1	C	70	KCX	CX-NZ	7.52	1.48	1.35
1	D	70	KCX	CX-NZ	6.73	1.47	1.35
1	C	70	KCX	OQ2-CX	-2.49	1.12	1.36
1	A	70	KCX	OQ2-CX	-2.49	1.12	1.36
1	B	70	KCX	OQ2-CX	-2.45	1.12	1.36
1	D	70	KCX	OQ2-CX	-2.38	1.13	1.36

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	70	KCX	CE-NZ-CX	-9.69	105.54	121.98
1	A	70	KCX	OQ1-CX-NZ	-9.49	110.50	124.92
1	B	70	KCX	CE-NZ-CX	-9.29	106.22	121.98
1	C	70	KCX	OQ1-CX-NZ	-9.21	110.92	124.92
1	C	70	KCX	CE-NZ-CX	-9.01	106.69	121.98
1	B	70	KCX	OQ1-CX-NZ	-8.31	112.30	124.92
1	D	70	KCX	CE-NZ-CX	-3.93	115.30	121.98
1	D	70	KCX	OQ1-CX-NZ	-3.03	120.31	124.92

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	D	70	KCX	C-CA-CB-CG

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

22 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	$\# Z > 2$
2	SO4	B	901	-	4,4,4	0.69	0	6,6,6	0.06	0
2	SO4	A	904	-	4,4,4	0.70	0	6,6,6	0.06	0
2	SO4	B	904	-	4,4,4	0.70	0	6,6,6	0.08	0
2	SO4	D	903	-	4,4,4	0.69	0	6,6,6	0.07	0
3	EDO	A	911	-	3,3,3	2.25	2 (66%)	2,2,2	0.45	0
2	SO4	C	904	-	4,4,4	0.66	0	6,6,6	0.25	0
2	SO4	D	902	-	4,4,4	0.68	0	6,6,6	0.16	0
2	SO4	C	901	-	4,4,4	0.67	0	6,6,6	0.12	0
3	EDO	D	911	-	3,3,3	2.13	2 (66%)	2,2,2	0.48	0
2	SO4	A	903	-	4,4,4	0.71	0	6,6,6	0.09	0
2	SO4	C	902	-	4,4,4	0.65	0	6,6,6	0.08	0
2	SO4	B	905	-	4,4,4	0.66	0	6,6,6	0.18	0
2	SO4	A	901	-	4,4,4	0.70	0	6,6,6	0.14	0
3	EDO	C	911	-	3,3,3	2.17	2 (66%)	2,2,2	0.39	0
2	SO4	B	903	-	4,4,4	0.60	0	6,6,6	0.11	0
3	EDO	D	912	-	3,3,3	2.23	2 (66%)	2,2,2	0.47	0
2	SO4	D	901	-	4,4,4	0.65	0	6,6,6	0.21	0
2	SO4	B	902	-	4,4,4	0.72	0	6,6,6	0.24	0
3	EDO	B	911	-	3,3,3	2.23	2 (66%)	2,2,2	0.45	0
3	EDO	B	912	-	3,3,3	2.23	2 (66%)	2,2,2	0.40	0
2	SO4	C	903	-	4,4,4	0.71	0	6,6,6	0.15	0
2	SO4	A	902	-	4,4,4	0.71	0	6,6,6	0.15	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	EDO	D	911	-	-	1/1/1/1	-
3	EDO	B	912	-	-	1/1/1/1	-
3	EDO	A	911	-	-	0/1/1/1	-
3	EDO	C	911	-	-	1/1/1/1	-
3	EDO	B	911	-	-	0/1/1/1	-
3	EDO	D	912	-	-	0/1/1/1	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	912	EDO	O2-C2	3.00	1.57	1.42
3	B	912	EDO	O2-C2	2.97	1.57	1.42
3	B	911	EDO	O2-C2	2.95	1.57	1.42
3	A	911	EDO	O2-C2	2.94	1.57	1.42
3	D	911	EDO	O2-C2	2.88	1.56	1.42
3	C	911	EDO	O2-C2	2.77	1.56	1.42
3	C	911	EDO	O1-C1	2.41	1.54	1.42
3	A	911	EDO	O1-C1	2.41	1.54	1.42
3	B	911	EDO	O1-C1	2.35	1.54	1.42
3	B	912	EDO	O1-C1	2.33	1.54	1.42
3	D	912	EDO	O1-C1	2.32	1.53	1.42
3	D	911	EDO	O1-C1	2.19	1.53	1.42

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	911	EDO	O1-C1-C2-O2
3	B	912	EDO	O1-C1-C2-O2
3	C	911	EDO	O1-C1-C2-O2

There are no ring outliers.

6 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	911	EDO	1	0
2	C	904	SO4	1	0
3	D	911	EDO	1	0
2	A	903	SO4	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	902	SO4	2	0
3	C	911	EDO	1	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	242/246 (98%)	0.21	3 (1%) 76 77	7, 22, 34, 44	3 (1%)
1	B	243/246 (98%)	0.08	1 (0%) 88 89	11, 19, 30, 41	3 (1%)
1	C	245/246 (99%)	0.23	11 (4%) 38 37	8, 19, 37, 47	3 (1%)
1	D	245/246 (99%)	-0.07	2 (0%) 82 83	8, 17, 32, 41	4 (1%)
All	All	975/984 (99%)	0.11	17 (1%) 69 69	7, 19, 34, 47	13 (1%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	94	GLY	4.0
1	C	26	THR	3.1
1	C	30	LYS	3.0
1	B	148	GLY	2.6
1	D	21	SER	2.6
1	C	265	GLY	2.5
1	D	104	ARG	2.4
1	C	212	GLY	2.4
1	A	36	ALA	2.3
1	C	28	TRP	2.3
1	C	36	ALA	2.2
1	C	27	SER	2.2
1	A	61	LYS	2.2
1	C	29	ASN	2.1
1	A	51	CYS	2.1
1	C	32	PHE	2.0
1	C	38	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	KCX	C	70	12/13	0.93	0.07	11,13,16,16	0
1	KCX	B	70	12/13	0.96	0.06	11,14,17,17	0
1	KCX	A	70	12/13	0.96	0.06	14,16,18,18	0
1	KCX	D	70	12/13	0.96	0.06	11,12,13,14	0

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(Å ²)	Q<0.9
2	SO4	C	901	5/5	0.65	0.25	41,41,41,42	0
2	SO4	B	902	5/5	0.67	0.20	44,44,44,45	0
2	SO4	A	903	5/5	0.74	0.14	55,55,55,55	0
2	SO4	B	905	5/5	0.79	0.11	44,44,44,44	0
2	SO4	C	903	5/5	0.79	0.13	47,47,48,48	0
2	SO4	D	903	5/5	0.80	0.12	52,52,52,52	0
2	SO4	A	902	5/5	0.81	0.12	46,46,47,47	0
2	SO4	B	904	5/5	0.82	0.10	55,56,56,56	0
2	SO4	C	902	5/5	0.84	0.12	47,47,47,47	0
3	EDO	C	911	4/4	0.84	0.17	33,33,33,34	0
3	EDO	D	912	4/4	0.84	0.18	30,30,30,31	0
2	SO4	D	902	5/5	0.85	0.31	43,43,43,43	0
2	SO4	A	904	5/5	0.86	0.13	49,49,49,49	0
2	SO4	B	901	5/5	0.86	0.12	50,50,50,50	0
3	EDO	D	911	4/4	0.87	0.20	25,27,28,29	0
2	SO4	D	901	5/5	0.92	0.12	30,30,31,31	0
3	EDO	B	912	4/4	0.93	0.08	27,27,27,27	0
2	SO4	B	903	5/5	0.93	0.12	29,29,30,30	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EDO	A	911	4/4	0.96	0.08	20,20,21,21	0
3	EDO	B	911	4/4	0.96	0.07	20,20,20,20	0
2	SO4	A	901	5/5	0.96	0.09	28,28,29,29	0
2	SO4	C	904	5/5	0.97	0.08	27,27,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.