



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:52 AM UTC

PDB ID : 5KIL / pdb_00005kil
Title : CmlA beta-hydroxylase E377D mutant
Authors : Knoot, C.J.; Lipscomb, J.D.
Deposited on : 2016-06-16
Resolution : 2.72 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

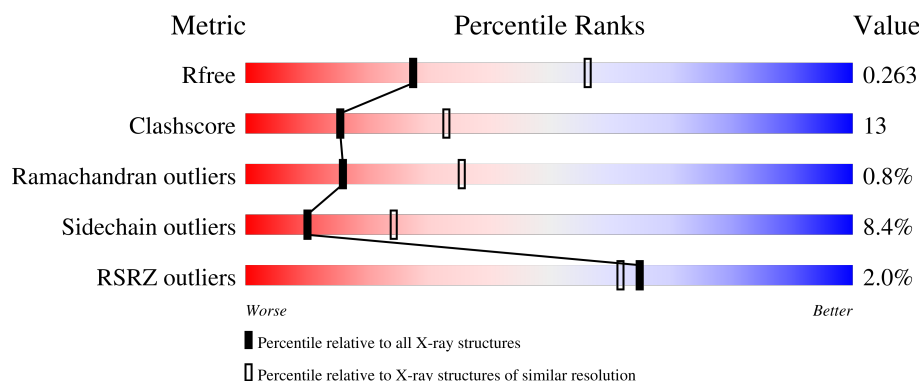
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

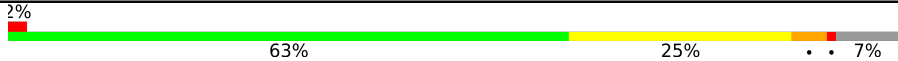
The reported resolution of this entry is 2.72 Å.

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	4348 (2.74-2.70)
Clashscore	190562	4665 (2.74-2.70)
Ramachandran outliers	187476	4584 (2.74-2.70)
Sidechain outliers	187428	4585 (2.74-2.70)
RSRZ outliers	180081	4348 (2.74-2.70)

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	551	

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
4	ACT	A	603	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4137 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.

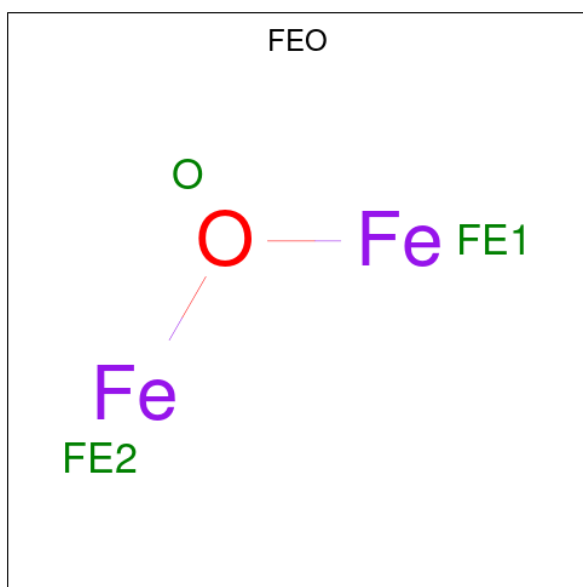
- Molecule 1 is a protein called CmlA protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	510	Total	C	N	O	S	0	1	0
			4110	2610	736	752	12			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP F2RB80
A	-18	GLY	-	expression tag	UNP F2RB80
A	-17	SER	-	expression tag	UNP F2RB80
A	-16	SER	-	expression tag	UNP F2RB80
A	-15	HIS	-	expression tag	UNP F2RB80
A	-14	HIS	-	expression tag	UNP F2RB80
A	-13	HIS	-	expression tag	UNP F2RB80
A	-12	HIS	-	expression tag	UNP F2RB80
A	-11	HIS	-	expression tag	UNP F2RB80
A	-10	SER	-	expression tag	UNP F2RB80
A	-9	SER	-	expression tag	UNP F2RB80
A	-8	GLY	-	expression tag	UNP F2RB80
A	-7	LEU	-	expression tag	UNP F2RB80
A	-6	VAL	-	expression tag	UNP F2RB80
A	-5	PRO	-	expression tag	UNP F2RB80
A	-4	ARG	-	expression tag	UNP F2RB80
A	-3	GLY	-	expression tag	UNP F2RB80
A	-2	SER	-	expression tag	UNP F2RB80
A	-1	HIS	-	expression tag	UNP F2RB80
A	189	ASN	ASP	see remark 999	UNP F2RB80
A	377	ASP	GLU	engineered mutation	UNP F2RB80

- Molecule 2 is MU-OXO-DIIRON (CCD ID: FEO) (formula: Fe₂O).

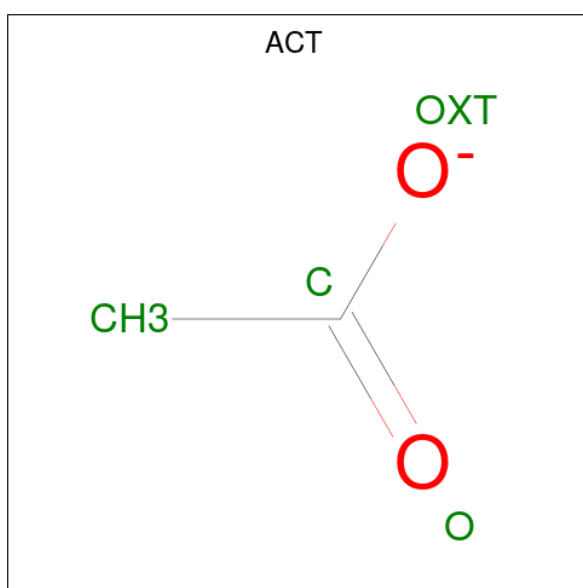


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	O	0	0
			3	2	1		

- Molecule 3 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	K	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		

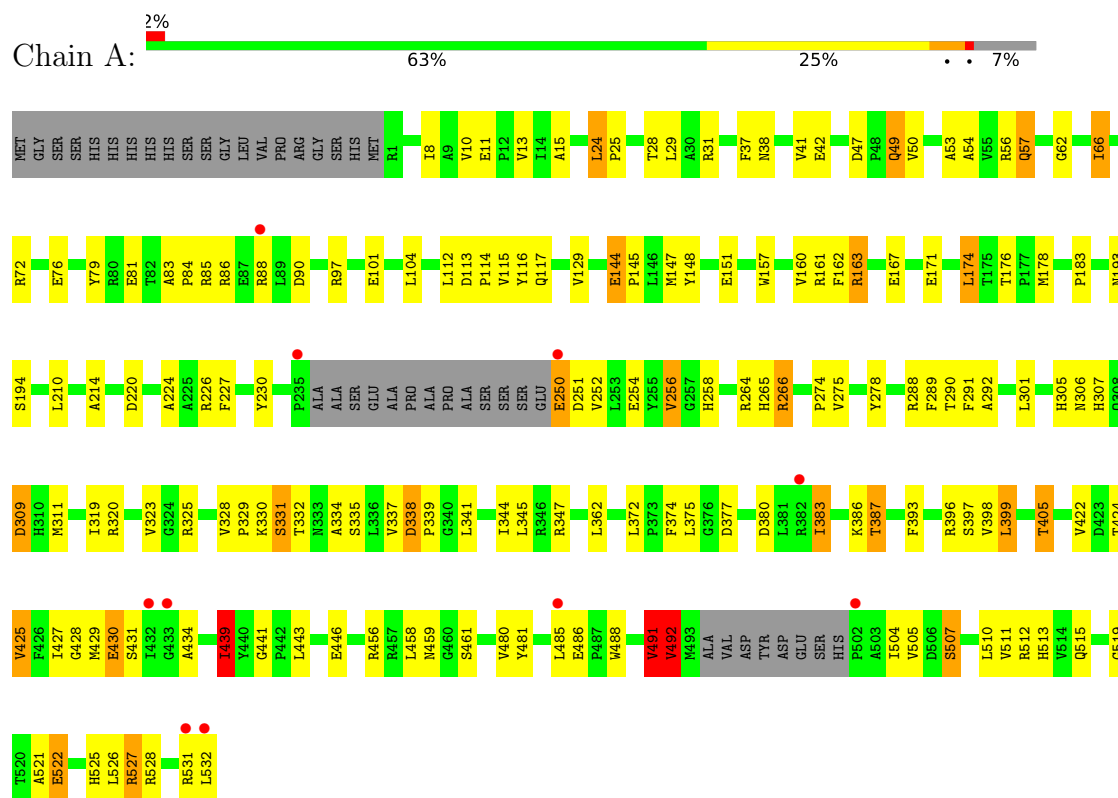
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	19	Total	O	0	0
			19	19		

3 Residue-property plots

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: CmlA protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	153.95Å 153.95Å 92.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.49 – 2.72 38.49 – 2.72	Depositor EDS
% Data completeness (in resolution range)	99.7 (38.49-2.72) 99.7 (38.49-2.72)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.83 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.202 , 0.262 0.202 , 0.263	Depositor DCC
R_{free} test set	1536 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	59.5	Xtriage
Anisotropy	0.202	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 37.2	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	4137	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.45% 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: FEO, ACT, K

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	1.20	11/4217 (0.3%)	1.29	31/5736 (0.5%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	338	ASP	CA-CB	12.71	1.66	1.53
1	A	383	ILE	C-O	11.57	1.35	1.24
1	A	250	GLU	C-O	10.31	1.44	1.23
1	A	334	ALA	C-O	9.14	1.35	1.23
1	A	250	GLU	CD-OE2	8.93	1.42	1.25
1	A	250	GLU	CD-OE1	7.78	1.40	1.25
1	A	335	SER	C-O	7.28	1.34	1.23
1	A	331	SER	N-CA	6.38	1.53	1.45
1	A	334	ALA	N-CA	-6.27	1.37	1.46
1	A	54	ALA	C-O	5.33	1.30	1.24
1	A	492	VAL	N-CA	-5.05	1.40	1.46

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	104	LEU	CA-C-N	-8.54	110.81	120.04
1	A	104	LEU	C-N-CA	-8.54	110.81	120.04
1	A	434	ALA	N-CA-C	7.77	120.61	110.43
1	A	337	VAL	CA-C-N	-6.87	114.58	122.00
1	A	337	VAL	C-N-CA	-6.87	114.58	122.00
1	A	148	TYR	N-CA-C	-6.68	104.60	112.89
1	A	491	VAL	CB-CA-C	-6.54	102.50	112.05
1	A	446	GLU	CA-C-N	-6.49	113.26	120.14
1	A	446	GLU	C-N-CA	-6.49	113.26	120.14
1	A	176	THR	CB-CA-C	-6.15	100.96	110.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	24	LEU	CA-C-N	6.02	126.58	120.38
1	A	24	LEU	C-N-CA	6.02	126.58	120.38
1	A	338	ASP	O-C-N	6.00	125.43	121.23
1	A	37	PHE	N-CA-C	5.94	118.52	111.33
1	A	194	SER	N-CA-C	5.92	116.41	109.60
1	A	439	ILE	CB-CA-C	-5.91	104.18	111.92
1	A	305	HIS	N-CA-C	5.75	116.97	108.86
1	A	62	GLY	N-CA-C	-5.73	107.87	114.69
1	A	383	ILE	N-CA-C	5.67	116.45	108.17
1	A	56	ARG	N-CA-C	5.66	117.44	111.28
1	A	330	LYS	N-CA-CB	-5.58	101.63	110.06
1	A	383	ILE	CB-CA-C	-5.57	101.48	110.28
1	A	174	LEU	N-CA-C	5.56	119.91	113.18
1	A	66	ILE	N-CA-C	5.53	116.14	108.84
1	A	441	GLY	CA-C-N	-5.51	114.01	120.12
1	A	441	GLY	C-N-CA	-5.51	114.01	120.12
1	A	331	SER	N-CA-C	-5.46	100.41	108.99
1	A	323	VAL	N-CA-C	5.39	117.06	108.87
1	A	163	ARG	CA-C-N	-5.30	114.78	120.03
1	A	163	ARG	C-N-CA	-5.30	114.78	120.03
1	A	306	ASN	N-CA-C	5.05	117.19	109.62

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	4110	0	3994	109	0
2	A	3	0	0	0	0
3	A	1	0	0	0	0
4	A	4	0	3	3	0
5	A	19	0	0	0	0
All	All	4137	0	3997	109	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 13.

All (109) 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:250:GLU:HG3	1:A:251:ASP:H	1.08	1.14
1:A:15:ALA:HA	1:A:157[B]:TRP:CE3	2.00	0.96
1:A:250:GLU:HG3	1:A:251:ASP:N	1.84	0.93
1:A:430:GLU:HG2	4:A:603:ACT:H2	1.54	0.90
1:A:183:PRO:O	1:A:226:ARG:HD2	1.74	0.88
1:A:250:GLU:CG	1:A:251:ASP:H	1.86	0.88
1:A:383:ILE:O	1:A:386:LYS:NZ	2.06	0.88
1:A:113:ASP:N	1:A:114:PRO:HD2	1.92	0.84
1:A:250:GLU:OE2	1:A:250:GLU:HA	1.77	0.83
1:A:41:VAL:HG11	1:A:79:TYR:CD1	2.13	0.81
1:A:250:GLU:HB3	1:A:264:ARG:NH1	2.03	0.72
1:A:13:VAL:HG22	1:A:178:MET:HG3	1.71	0.72
1:A:220:ASP:HA	1:A:224:ALA:HB2	1.73	0.70
1:A:488:TRP:O	1:A:491:VAL:HG23	1.93	0.69
1:A:372:LEU:HD11	1:A:399:LEU:HD13	1.75	0.68
1:A:85:ARG:CZ	1:A:157[A]:TRP:CD1	2.76	0.68
1:A:398:VAL:HG22	1:A:424:THR:HB	1.76	0.67
1:A:53:ALA:O	1:A:57:GLN:NE2	2.26	0.67
1:A:113:ASP:N	1:A:114:PRO:CD	2.58	0.66
1:A:430:GLU:CG	4:A:603:ACT:H2	2.25	0.66
1:A:372:LEU:CD1	1:A:399:LEU:HD13	2.26	0.65
1:A:430:GLU:O	1:A:430:GLU:HG3	1.97	0.65
1:A:15:ALA:HA	1:A:157[B]:TRP:CD2	2.33	0.63
1:A:278:TYR:O	1:A:288:ARG:NH2	2.31	0.63
1:A:374:PHE:CD1	1:A:386:LYS:HD2	2.36	0.61
1:A:88:ARG:HH11	1:A:88:ARG:HB3	1.65	0.60
1:A:396:ARG:NH2	1:A:531:ARG:O	2.34	0.60
1:A:88:ARG:HB3	1:A:88:ARG:NH1	2.15	0.60
1:A:329:PRO:HB3	1:A:387:THR:HB	1.84	0.59
1:A:274:PRO:HG3	1:A:301:LEU:HD11	1.84	0.59
1:A:112:LEU:C	1:A:114:PRO:HD2	2.27	0.59
1:A:38:ASN:O	1:A:42:GLU:HG3	2.03	0.59
1:A:331:SER:OG	1:A:338:ASP:OD2	2.20	0.58
1:A:399:LEU:HD23	1:A:425:VAL:HG22	1.86	0.56
1:A:512:ARG:O	1:A:513:HIS:C	2.50	0.54
1:A:522:GLU:OE1	1:A:528:ARG:NH1	2.40	0.54
1:A:24:LEU:CB	1:A:25:PRO:HD3	2.38	0.54
1:A:290:THR:HG22	1:A:291:PHE:N	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:ASP:HA	1:A:224:ALA:CB	2.40	0.52
1:A:328:VAL:HG11	1:A:345:LEU:CD1	2.40	0.52
1:A:396:ARG:HH21	1:A:532:LEU:HA	1.75	0.52
1:A:15:ALA:HA	1:A:157[B]:TRP:CZ3	2.45	0.52
1:A:41:VAL:HG11	1:A:79:TYR:CE1	2.45	0.51
1:A:10:VAL:HG22	1:A:162:PHE:CD2	2.45	0.51
1:A:338:ASP:HB2	1:A:383:ILE:HG12	1.92	0.51
1:A:250:GLU:OE2	1:A:250:GLU:CA	2.52	0.51
1:A:429:MET:O	1:A:431:SER:N	2.43	0.51
1:A:79:TYR:O	1:A:86:ARG:NH1	2.45	0.50
1:A:83:ALA:N	1:A:84:PRO:HD2	2.26	0.50
1:A:256:VAL:O	1:A:256:VAL:HG23	2.11	0.50
1:A:254:GLU:OE2	1:A:527:ARG:NH1	2.47	0.48
1:A:114:PRO:O	1:A:117:GLN:HB2	2.13	0.48
1:A:372:LEU:HD12	1:A:399:LEU:CD1	2.43	0.48
1:A:214:ALA:HB2	1:A:227:PHE:CD2	2.49	0.48
1:A:256:VAL:HB	1:A:289:PHE:HB3	1.96	0.47
1:A:507:SER:O	1:A:511:VAL:HG23	2.14	0.47
1:A:525:HIS:CE1	1:A:526:LEU:HD23	2.49	0.47
1:A:430:GLU:CB	4:A:603:ACT:H2	2.45	0.47
1:A:174:LEU:HD11	1:A:492:VAL:HG23	1.97	0.47
1:A:250:GLU:HB3	1:A:264:ARG:HH11	1.77	0.46
1:A:311:MET:O	1:A:311:MET:HG3	2.14	0.46
1:A:439:ILE:HG13	1:A:458:LEU:HD11	1.97	0.46
1:A:362:LEU:HA	1:A:362:LEU:HD23	1.67	0.46
1:A:254:GLU:CD	1:A:527:ARG:NH1	2.73	0.46
1:A:47:ASP:O	1:A:50:VAL:HG23	2.16	0.46
1:A:112:LEU:O	1:A:115:VAL:HG23	2.16	0.46
1:A:521:ALA:O	1:A:522:GLU:HB2	2.16	0.45
1:A:24:LEU:HD12	1:A:344:ILE:HG23	1.98	0.45
1:A:144:GLU:N	1:A:145:PRO:HD2	2.31	0.45
1:A:266:ARG:HH11	1:A:266:ARG:HB2	1.82	0.45
1:A:347:ARG:HD3	1:A:347:ARG:HA	1.70	0.45
1:A:8:ILE:HD12	1:A:230:TYR:CD1	2.51	0.44
1:A:214:ALA:HB2	1:A:227:PHE:CE2	2.52	0.44
1:A:289:PHE:C	1:A:289:PHE:CD1	2.95	0.44
1:A:372:LEU:CD1	1:A:399:LEU:CD1	2.95	0.44
1:A:31:ARG:HB2	1:A:443:LEU:HD23	2.00	0.44
1:A:374:PHE:CG	1:A:386:LYS:HB2	2.52	0.44
1:A:429:MET:C	1:A:431:SER:N	2.76	0.44
1:A:85:ARG:NE	1:A:157[A]:TRP:CD1	2.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:LEU:HD23	1:A:210:LEU:HA	1.84	0.43
1:A:47:ASP:C	1:A:49:GLN:N	2.75	0.43
1:A:329:PRO:HD2	1:A:341:LEU:HD23	2.01	0.43
1:A:405:THR:HA	1:A:461:SER:OG	2.18	0.43
1:A:116:TYR:CZ	1:A:129:VAL:HG22	2.54	0.43
1:A:290:THR:CG2	1:A:291:PHE:N	2.81	0.43
1:A:57:GLN:HE21	1:A:57:GLN:H	1.67	0.43
1:A:515:GLN:HA	1:A:519:GLY:O	2.19	0.43
1:A:28:THR:O	1:A:29:LEU:C	2.61	0.42
1:A:427:ILE:HD12	1:A:428:GLY:H	1.82	0.42
1:A:254:GLU:CD	1:A:527:ARG:HH12	2.27	0.42
1:A:307:HIS:HE1	1:A:377:ASP:OD2	1.99	0.42
1:A:193:ASN:HB3	1:A:347:ARG:O	2.20	0.42
1:A:319:ILE:O	1:A:320:ARG:C	2.62	0.42
1:A:11:GLU:OE1	1:A:163:ARG:HD3	2.20	0.42
1:A:486:GLU:H	1:A:486:GLU:CD	2.27	0.42
1:A:24:LEU:HB2	1:A:25:PRO:HD3	2.02	0.42
1:A:338:ASP:CG	1:A:383:ILE:HG23	2.46	0.41
1:A:250:GLU:CG	1:A:251:ASP:N	2.54	0.41
1:A:47:ASP:C	1:A:49:GLN:H	2.27	0.41
1:A:72:ARG:NH2	1:A:76:GLU:OE1	2.54	0.41
1:A:265:HIS:CE1	1:A:393:PHE:HB3	2.55	0.41
1:A:290:THR:HG22	1:A:292:ALA:H	1.85	0.41
1:A:24:LEU:N	1:A:25:PRO:CD	2.84	0.40
1:A:309:ASP:OD1	1:A:309:ASP:N	2.53	0.40
1:A:485:LEU:HB3	1:A:525:HIS:CD2	2.56	0.40
1:A:504:ILE:O	1:A:505:VAL:C	2.64	0.40
1:A:97:ARG:O	1:A:101:GLU:HB2	2.22	0.40
1:A:147:MET:HE3	1:A:147:MET:HB3	1.95	0.40
1:A:183:PRO:O	1:A:226:ARG:CD	2.58	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	505/551 (92%)	460 (91%)	41 (8%)	4 (1%)	16	35

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	430	GLU
1	A	258	HIS
1	A	522	GLU
1	A	339	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	431/462 (93%)	395 (92%)	36 (8%)	10	24

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	GLN
1	A	57	GLN
1	A	66	ILE
1	A	81	GLU
1	A	90	ASP
1	A	144	GLU
1	A	151	GLU
1	A	160	VAL
1	A	161	ARG
1	A	167	GLU
1	A	171	GLU
1	A	252	VAL
1	A	256	VAL
1	A	266	ARG
1	A	275	VAL

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Mol	Chain	Res	Type
1	A	309	ASP
1	A	325	ARG
1	A	332	THR
1	A	375	LEU
1	A	380	ASP
1	A	387	THR
1	A	397	SER
1	A	399	LEU
1	A	405	THR
1	A	422	VAL
1	A	425	VAL
1	A	439	ILE
1	A	456	ARG
1	A	459	ASN
1	A	480	VAL
1	A	481	TYR
1	A	491	VAL
1	A	492	VAL
1	A	507	SER
1	A	510	LEU
1	A	527	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	GLN
1	A	71	HIS
1	A	265	HIS
1	A	459	ASN
1	A	525	HIS

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	$\# Z > 2$
4	ACT	A	603	2	3,3,3	0.82	0	3,3,3	1.14	0
2	FEO	A	601	4,1	0,2,2	-	-	-		

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	603	ACT	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	510/551 (92%)	-0.14	10 (1%) 65 62	20, 60, 95, 128	1 (0%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	502	PRO	7.4
1	A	531	ARG	3.5
1	A	485	LEU	3.2
1	A	88	ARG	2.8
1	A	250	GLU	2.7
1	A	432	ILE	2.5
1	A	382	ARG	2.2
1	A	235	PRO	2.2
1	A	433	GLY	2.2
1	A	532	LEU	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	A	603	4/4	0.85	0.28	64,69,74,74	4
3	K	A	602	1/1	0.97	0.09	78,78,78,78	0
2	FEO	A	601	3/3	0.99	0.03	48,48,61,67	1

6.5 Other polymers [i](#)

There are no such residues in this entry.