



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:22 PM UTC

PDB ID : 6S87 / pdb_00006s87
Title : Crystal structure of 2-methylcitrate synthase (PrpC) from *Pseudomonas aeruginosa* in complex with oxaloacetate.
Authors : Wijaya, A.J.; Brear, P.; Dolan, S.K.; Welch, M.
Deposited on : 2019-07-08
Resolution : 1.65 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

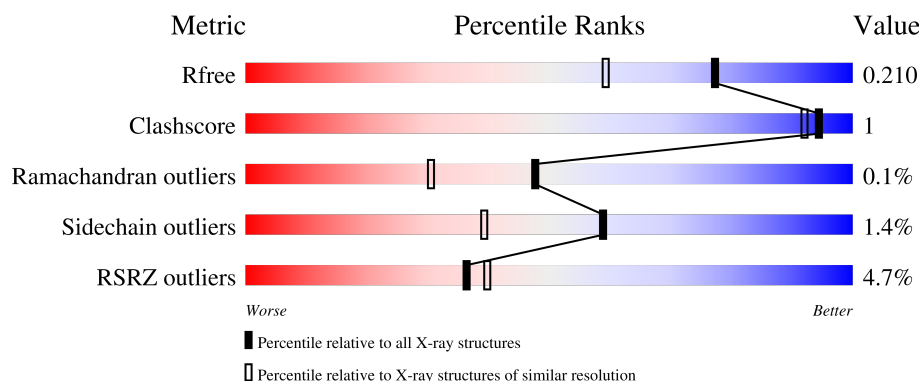
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	375	 4% 90% 5% •
1	B	375	 7% 86% 7% • 6%
1	C	375	 5% 90% • • 6%
1	D	375	 2% 90% 7% •

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 11684 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 Citrate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	359	Total	C	N	O	S	0	0	0
			2826	1796	489	528	13			
1	B	353	Total	C	N	O	S	0	0	0
			2790	1776	482	519	13			
1	C	354	Total	C	N	O	S	0	0	0
			2794	1778	483	520	13			
1	D	365	Total	C	N	O	S	0	1	0
			2877	1828	500	536	13			

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



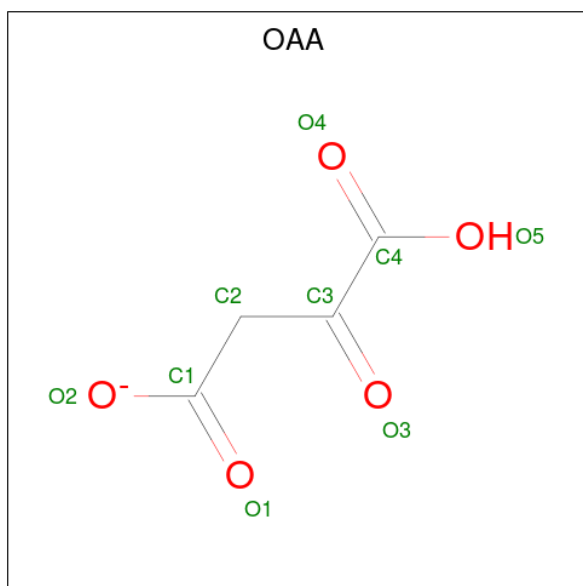
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is OXALOACETATE ION (CCD ID: OAA) (formula: $C_4H_3O_5$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	C	O	0	0
			9	4	5		

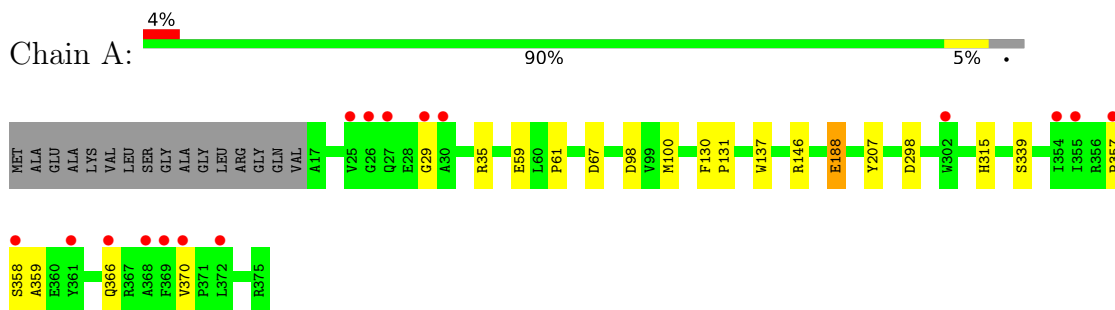
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	106	Total	O	0	0
			106	106		
4	B	44	Total	O	0	0
			44	44		
4	C	85	Total	O	0	1
			86	86		
4	D	116	Total	O	0	0
			116	116		

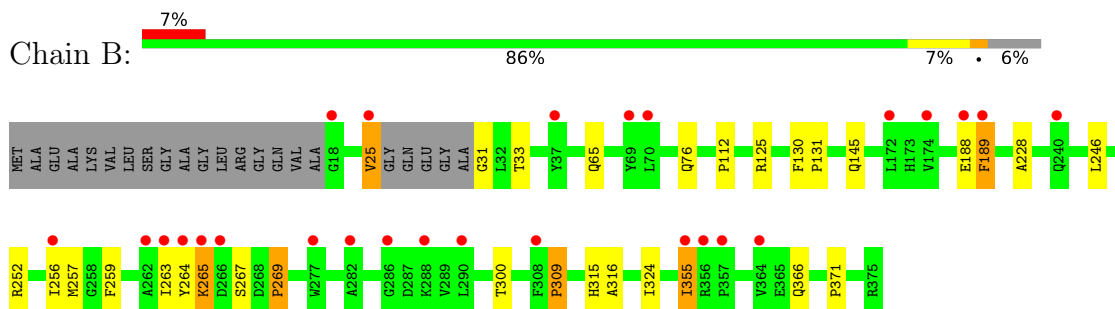
3 Residue-property plots [i](#)

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.

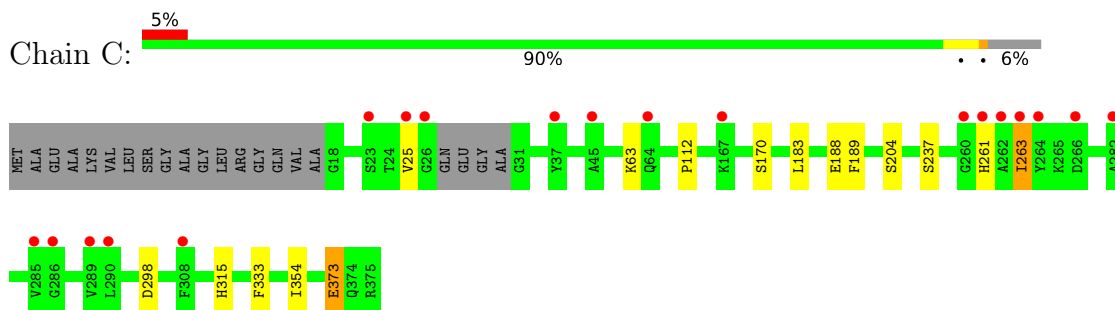
- Molecule 1: Citrate synthase



- Molecule 1: Citrate synthase

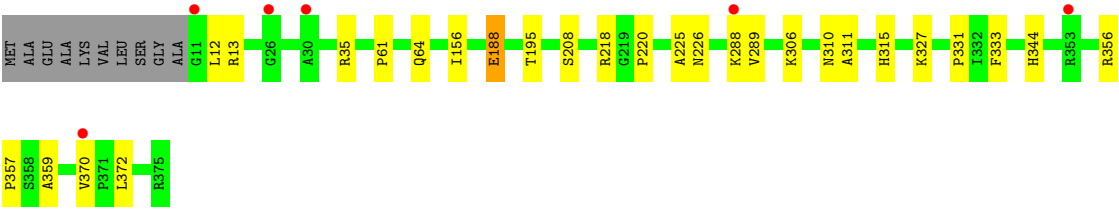


- Molecule 1: Citrate synthase



- Molecule 1: Citrate synthase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.64Å 86.76Å 161.84Å 90.00° 95.89° 90.00°	Depositor
Resolution (Å)	27.23 – 1.65 27.23 – 1.65	Depositor EDS
% Data completeness (in resolution range)	99.8 (27.23-1.65) 99.8 (27.23-1.65)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 1.65Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.197 , 0.215 0.210 , 0.210	Depositor DCC
R_{free} test set	8371 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	22.0	Xtriage
Anisotropy	0.484	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 41.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11684	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 68.25 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.5037e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	1.33	8/2890 (0.3%)	1.33	4/3911 (0.1%)
1	B	1.22	4/2853 (0.1%)	1.42	6/3860 (0.2%)
1	C	1.25	3/2857 (0.1%)	1.32	3/3865 (0.1%)
1	D	1.35	14/2941 (0.5%)	1.30	4/3980 (0.1%)
All	All	1.29	29/11541 (0.3%)	1.34	17/15616 (0.1%)

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	61	PRO	C-O	-7.13	1.15	1.23
1	D	331	PRO	C-O	-7.00	1.15	1.24
1	D	311	ALA	C-O	-6.84	1.15	1.24
1	A	61	PRO	C-O	-6.79	1.16	1.23
1	A	131	PRO	C-O	-6.59	1.15	1.24
1	D	188	GLU	C-O	6.17	1.30	1.24
1	C	112	PRO	C-O	-5.98	1.17	1.23
1	A	100	MET	C-O	-5.96	1.17	1.24
1	D	156	ILE	C-O	5.91	1.30	1.24
1	D	370	VAL	N-CA	5.89	1.50	1.46
1	D	220	PRO	C-O	-5.85	1.16	1.24
1	D	35	ARG	C-O	5.67	1.30	1.23
1	D	310	ASN	C-O	5.62	1.30	1.23
1	A	188	GLU	C-O	-5.54	1.19	1.24
1	D	208	SER	C-O	5.53	1.30	1.24
1	B	309	PRO	C-O	-5.50	1.17	1.23
1	B	324	ILE	N-CA	5.46	1.50	1.46
1	A	98	ASP	C-O	5.45	1.31	1.24
1	D	344	HIS	CE1-NE2	5.38	1.38	1.32
1	B	188	GLU	C-O	-5.38	1.19	1.24
1	A	35	ARG	C-O	5.34	1.30	1.23
1	C	204	SER	C-O	5.22	1.31	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	225	ALA	C-O	-5.22	1.17	1.24
1	D	357	PRO	C-O	-5.19	1.18	1.23
1	C	183	LEU	C-O	5.19	1.30	1.24
1	D	195	THR	C-O	-5.10	1.18	1.24
1	A	137	TRP	C-O	5.10	1.29	1.24
1	B	371	PRO	C-O	-5.10	1.17	1.23
1	A	339	SER	C-O	-5.04	1.18	1.24

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	289	VAL	CA-C-O	-5.89	115.56	120.80
1	D	333	PHE	CA-CB-CG	5.73	119.53	113.80
1	B	265	LYS	CB-CA-C	5.72	119.34	109.15
1	C	188	GLU	N-CA-C	5.71	116.99	108.31
1	B	189	PHE	CA-C-N	5.53	131.28	122.10
1	B	189	PHE	C-N-CA	5.53	131.28	122.10
1	D	218	ARG	CG-CD-NE	5.53	124.16	112.00
1	A	298	ASP	CA-CB-CG	5.51	118.11	112.60
1	A	207	TYR	CA-C-O	-5.43	114.79	120.55
1	C	263	ILE	CA-C-O	-5.41	115.42	120.34
1	B	269	PRO	CB-CA-C	-5.39	103.26	112.26
1	A	146	ARG	CA-C-O	-5.38	114.82	120.58
1	B	112	PRO	N-CA-C	5.37	119.74	111.21
1	A	59	GLU	CB-CG-CD	5.25	121.53	112.60
1	B	316	ALA	CA-C-O	-5.16	115.40	120.82
1	D	226	ASN	CB-CA-C	5.10	119.25	110.79
1	C	298	ASP	CA-CB-CG	5.07	117.67	112.60

There are no chirality outliers.

There are no planarity outliers.

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	2826	0	2787	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2790	0	2756	17	0
1	C	2794	0	2759	7	0
1	D	2877	0	2845	9	0
2	A	6	0	8	0	0
2	B	12	0	16	0	0
2	C	6	0	8	1	0
2	D	12	0	16	0	0
3	D	9	0	2	0	0
4	A	106	0	0	0	0
4	B	44	0	0	0	0
4	C	86	0	0	0	0
4	D	116	0	0	2	0
All	All	11684	0	11197	30	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 1.

All (30) 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:359:ALA:HB2	1:B:189:PHE:CE1	2.26	0.70
1:A:359:ALA:HB2	1:B:189:PHE:CZ	2.28	0.67
1:B:252:ARG:O	1:B:252:ARG:HG2	2.04	0.57
1:B:31:GLY:HA3	1:B:269:PRO:HG3	1.87	0.55
1:C:25:VAL:HG21	1:C:263:ILE:HG12	1.90	0.54
1:B:25:VAL:HG11	1:B:263:ILE:HG22	1.88	0.54
1:C:373:GLU:H	1:C:373:GLU:CD	2.17	0.52
1:B:25:VAL:HG11	1:B:263:ILE:CG2	2.40	0.52
1:D:12:LEU:HD22	1:D:12:LEU:N	2.25	0.52
1:A:357:PRO:HG2	1:B:355:ILE:HG21	1.93	0.51
1:B:76:GLN:O	1:B:125:ARG:NH1	2.44	0.49
1:B:264:TYR:HB3	1:B:267:SER:O	2.13	0.48
1:C:333:PHE:CE2	2:C:401:GOL:H12	2.49	0.47
1:D:64:GLN:NE2	4:D:501:HOH:O	2.48	0.46
1:C:189:PHE:CZ	1:D:359:ALA:HB2	2.51	0.46
1:A:366:GLN:HG2	1:B:33:THR:HG21	1.98	0.45
1:B:130:PHE:N	1:B:131:PRO:CD	2.79	0.45
1:D:356:ARG:O	1:D:356:ARG:HG3	2.17	0.45
1:C:63:LYS:HD2	1:D:372:LEU:HD21	1.98	0.45
1:B:259:PHE:CE2	1:B:309:PRO:HB3	2.52	0.44
1:D:306:LYS:HD2	1:D:306:LYS:N	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:359:ALA:CB	1:B:189:PHE:CE1	2.99	0.43
1:A:370:VAL:O	1:A:370:VAL:HG23	2.19	0.43
1:B:228:ALA:HB1	1:B:257:MET:HG3	2.01	0.42
1:C:261:HIS:CE1	1:C:263:ILE:HG22	2.55	0.42
1:D:327:LYS:NZ	4:D:504:HOH:O	2.53	0.41
1:A:29:GLY:HA3	1:B:366:GLN:HE22	1.85	0.40
1:B:65:GLN:NE2	1:D:288:LYS:HE2	2.36	0.40
1:B:246:LEU:HD23	1:B:300:THR:CG2	2.51	0.40
1:C:354:ILE:HD11	1:D:13:ARG:NH1	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	357/375 (95%)	352 (99%)	4 (1%)	1 (0%)	36	21
1	B	349/375 (93%)	344 (99%)	5 (1%)	0	100	100
1	C	350/375 (93%)	343 (98%)	7 (2%)	0	100	100
1	D	364/375 (97%)	358 (98%)	6 (2%)	0	100	100
All	All	1420/1500 (95%)	1397 (98%)	22 (2%)	1 (0%)	48	30

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	358	SER

5.3.2 Protein sidechains [i](#)

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

resolution.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	296/306 (97%)	292 (99%)	4 (1%)	59	39
1	B	294/306 (96%)	288 (98%)	6 (2%)	48	26
1	C	294/306 (96%)	290 (99%)	4 (1%)	59	39
1	D	301/306 (98%)	299 (99%)	2 (1%)	76	64
All	All	1185/1224 (97%)	1169 (99%)	16 (1%)	59	39

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	ASP
1	A	130	PHE
1	A	188	GLU
1	A	315	HIS
1	B	25	VAL
1	B	145	GLN
1	B	256	ILE
1	B	265	LYS
1	B	315	HIS
1	B	355	ILE
1	C	170	SER
1	C	237	SER
1	C	315	HIS
1	C	373	GLU
1	D	188	GLU
1	D	315	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	GLN
1	A	374	GLN
1	B	76	GLN
1	B	150	ASN
1	B	374	GLN
1	C	64	GLN
1	C	76	GLN

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Mol	Chain	Res	Type
1	C	165	HIS
1	D	64	GLN
1	D	150	ASN
1	D	165	HIS
1	D	190	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 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	GOL	C	401	-	5,5,5	0.21	0	5,5,5	0.38	0
3	OAA	D	403	-	8,8,8	5.15	3 (37%)	8,10,10	1.92	3 (37%)
2	GOL	D	402	-	5,5,5	0.22	0	5,5,5	1.04	0
2	GOL	D	401	-	5,5,5	0.17	0	5,5,5	0.36	0
2	GOL	B	401	-	5,5,5	0.15	0	5,5,5	0.44	0
2	GOL	A	401	-	5,5,5	0.28	0	5,5,5	0.44	0
2	GOL	B	402	-	5,5,5	0.15	0	5,5,5	0.53	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	GOL	C	401	-	-	0/4/4/4	-
3	OAA	D	403	-	-	3/8/8/8	-
2	GOL	D	402	-	-	3/4/4/4	-
2	GOL	D	401	-	-	0/4/4/4	-
2	GOL	B	401	-	-	2/4/4/4	-
2	GOL	A	401	-	-	0/4/4/4	-
2	GOL	B	402	-	-	2/4/4/4	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	403	OAA	C3-C4	-13.36	1.33	1.53
3	D	403	OAA	O2-C1	-4.81	1.14	1.30
3	D	403	OAA	O5-C4	-2.40	1.24	1.30

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	403	OAA	O4-C4-C3	-3.73	117.17	121.81
3	D	403	OAA	O5-C4-C3	2.82	121.42	113.59
3	D	403	OAA	O3-C3-C2	-2.11	117.16	120.41

There are no chirality outliers.

All (10) torsion outliers are listed below:

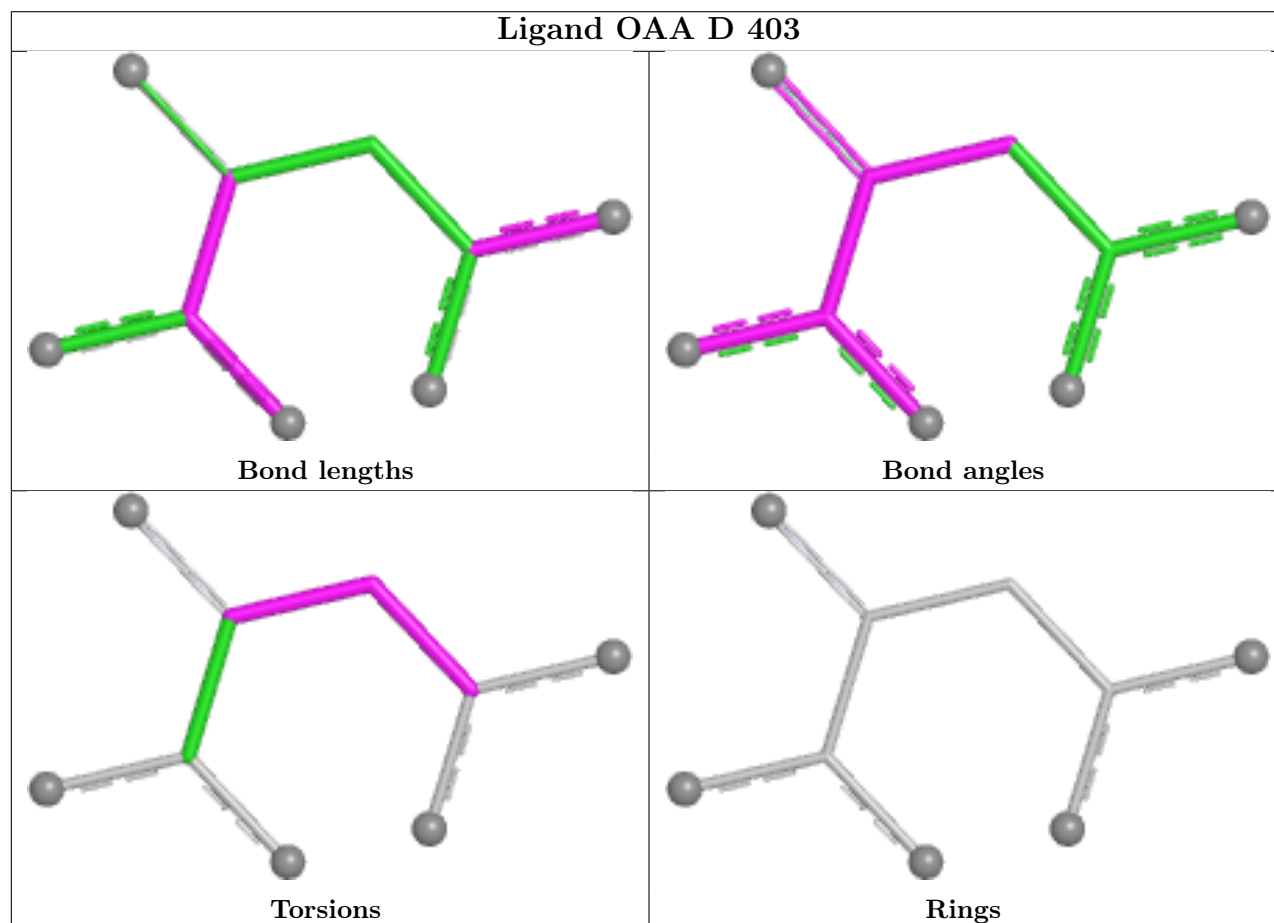
Mol	Chain	Res	Type	Atoms
2	D	402	GOL	O1-C1-C2-C3
2	B	401	GOL	O1-C1-C2-C3
2	B	402	GOL	O1-C1-C2-C3
2	D	402	GOL	C1-C2-C3-O3
2	B	402	GOL	O1-C1-C2-O2
2	D	402	GOL	O1-C1-C2-O2
2	B	401	GOL	O1-C1-C2-O2
3	D	403	OAA	O1-C1-C2-C3
3	D	403	OAA	C1-C2-C3-O3
3	D	403	OAA	C1-C2-C3-C4

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	GOL	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	359/375 (95%)	0.33	16 (4%) 38 42	18, 25, 55, 73	0
1	B	353/375 (94%)	0.87	26 (7%) 20 23	21, 38, 59, 93	0
1	C	354/375 (94%)	0.50	19 (5%) 31 35	19, 30, 49, 77	0
1	D	365/375 (97%)	0.12	6 (1%) 70 75	10, 23, 46, 75	1 (0%)
All	All	1431/1500 (95%)	0.45	67 (4%) 36 40	10, 29, 53, 93	1 (0%)

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	263	ILE	4.9
1	C	263	ILE	4.6
1	C	289	VAL	4.2
1	B	18	GLY	4.1
1	B	308	PHE	4.0
1	A	27	GLN	4.0
1	C	26	GLY	3.9
1	A	366	GLN	3.9
1	B	25	VAL	3.8
1	C	262	ALA	3.8
1	C	23	SER	3.5
1	C	260	GLY	3.5
1	A	30	ALA	3.5
1	A	26	GLY	3.4
1	A	355	ILE	3.3
1	A	357	PRO	3.3
1	B	355	ILE	3.2
1	B	356	ARG	3.2
1	A	25	VAL	3.2
1	B	262	ALA	3.1
1	C	266	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	264	TYR	3.0
1	B	364	VAL	3.0
1	A	370	VAL	2.9
1	C	64	GLN	2.9
1	B	282	ALA	2.8
1	B	189	PHE	2.8
1	D	11	GLY	2.8
1	B	290	LEU	2.7
1	C	286	GLY	2.7
1	C	261	HIS	2.6
1	A	302	TRP	2.6
1	C	45	ALA	2.6
1	A	372	LEU	2.5
1	B	288	LYS	2.5
1	D	26	GLY	2.5
1	B	70	LEU	2.4
1	B	172	LEU	2.4
1	B	357	PRO	2.4
1	C	37	TYR	2.4
1	B	174	VAL	2.4
1	A	358	SER	2.4
1	B	286	GLY	2.3
1	B	265	LYS	2.3
1	A	361	TYR	2.3
1	D	370	VAL	2.3
1	C	290	LEU	2.3
1	B	37	TYR	2.3
1	B	264	TYR	2.3
1	A	368	ALA	2.2
1	B	277	TRP	2.2
1	B	240	GLN	2.2
1	C	282	ALA	2.2
1	D	30	ALA	2.2
1	D	288	LYS	2.2
1	B	256	ILE	2.2
1	B	266	ASP	2.1
1	B	69	TYR	2.1
1	C	285	VAL	2.1
1	C	25	VAL	2.1
1	D	353	ARG	2.1
1	B	188	GLU	2.0
1	A	29	GLY	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	354	ILE	2.0
1	A	369	PHE	2.0
1	C	308	PHE	2.0
1	C	167	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

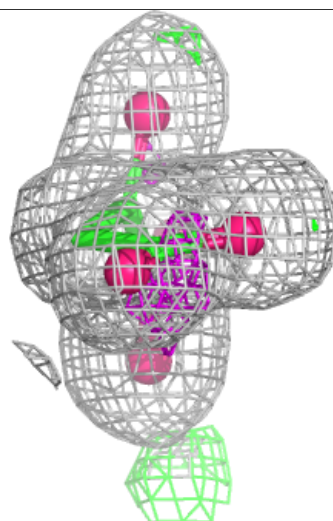
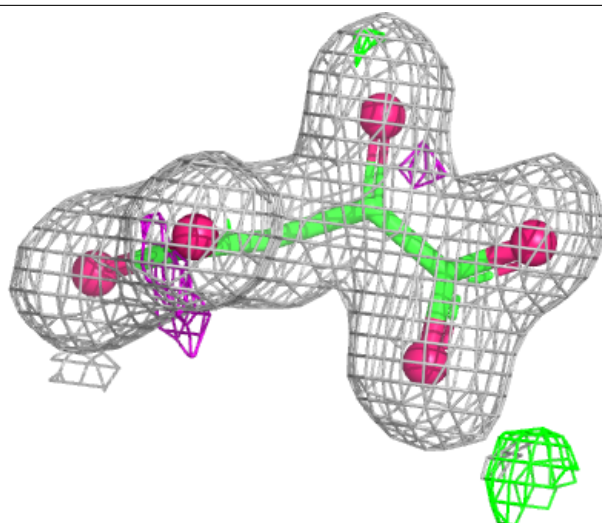
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	GOL	C	401	6/6	0.80	0.13	43,45,49,49	0
2	GOL	B	402	6/6	0.82	0.14	51,52,53,54	0
2	GOL	A	401	6/6	0.88	0.10	26,32,34,36	0
2	GOL	B	401	6/6	0.88	0.12	44,50,53,54	0
2	GOL	D	401	6/6	0.89	0.11	28,34,38,38	0
3	OAA	D	403	9/9	0.92	0.08	20,23,24,24	0
2	GOL	D	402	6/6	0.94	0.09	21,30,33,34	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around OAA D 403:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.