



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 01:09 PM UTC

PDB ID : 4Q2B / pdb_00004q2b
Title : The crystal structure of an endo-1,4-D-glucanase from *Pseudomonas putida* KT2440
Authors : Tan, K.; Joachimiak, G.; Endres, M.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2014-04-07
Resolution : 2.12 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

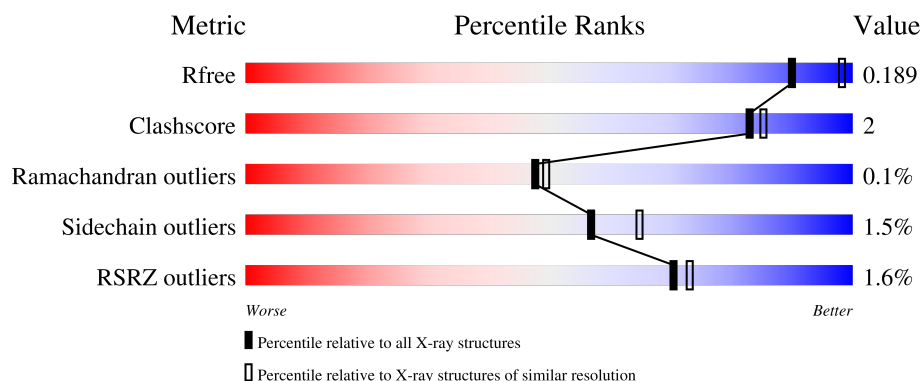
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.12 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	8290 (2.14-2.10)
Clashscore	190562	8817 (2.14-2.10)
Ramachandran outliers	187476	8738 (2.14-2.10)
Sidechain outliers	187428	8739 (2.14-2.10)
RSRZ outliers	180081	8294 (2.14-2.10)

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	349	0% 93% 5% .
1	B	349	2% 90% 8% .
1	C	349	0% 91% 7% .
1	D	349	2% 94% . .
1	E	349	2% 92% 5% .

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Mol	Chain	Length	Quality of chain
1	F	349	% 90% 8%

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
3	TRS	F	402	-	X	-	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 17169 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 Endo-1,4-beta-D-glucanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	342	Total	C	N	O	Se	0	2	0
			2685	1721	477	484	3			
1	B	344	Total	C	N	O	Se	0	3	0
			2711	1736	488	484	3			
1	C	343	Total	C	N	O	Se	0	2	0
			2709	1735	483	488	3			
1	D	344	Total	C	N	O	Se	0	1	0
			2688	1725	480	480	3			
1	E	342	Total	C	N	O	Se	0	0	0
			2673	1710	483	477	3			
1	F	343	Total	C	N	O	Se	0	0	0
			2679	1715	480	481	3			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	SER	-	expression tag	UNP Q88JL2
A	23	ASN	-	expression tag	UNP Q88JL2
A	24	ALA	-	expression tag	UNP Q88JL2
B	22	SER	-	expression tag	UNP Q88JL2
B	23	ASN	-	expression tag	UNP Q88JL2
B	24	ALA	-	expression tag	UNP Q88JL2
C	22	SER	-	expression tag	UNP Q88JL2
C	23	ASN	-	expression tag	UNP Q88JL2
C	24	ALA	-	expression tag	UNP Q88JL2
D	22	SER	-	expression tag	UNP Q88JL2
D	23	ASN	-	expression tag	UNP Q88JL2
D	24	ALA	-	expression tag	UNP Q88JL2
E	22	SER	-	expression tag	UNP Q88JL2
E	23	ASN	-	expression tag	UNP Q88JL2
E	24	ALA	-	expression tag	UNP Q88JL2
F	22	SER	-	expression tag	UNP Q88JL2
F	23	ASN	-	expression tag	UNP Q88JL2

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Chain	Residue	Modelled	Actual	Comment	Reference
F	24	ALA	-	expression tag	UNP Q88JL2

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



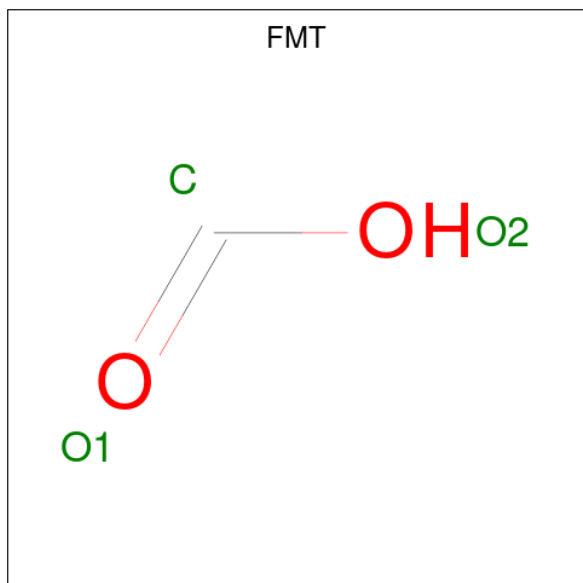
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: C₄H₁₂NO₃).



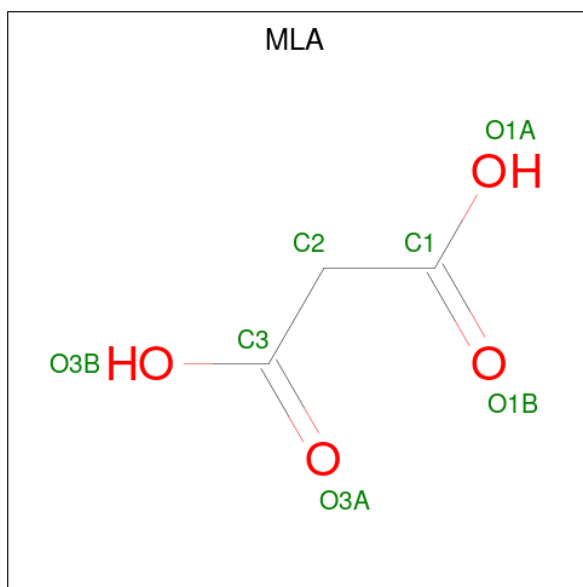
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			8	4	1	3		
3	D	1	Total	C	N	O	0	0
			8	4	1	3		
3	E	1	Total	C	N	O	0	0
			8	4	1	3		
3	F	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 4 is FORMIC ACID (CCD ID: FMT) (formula: CH₂O₂).



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

- Molecule 5 is MALONIC ACID (CCD ID: MLA) (formula: $C_3H_4O_4$).



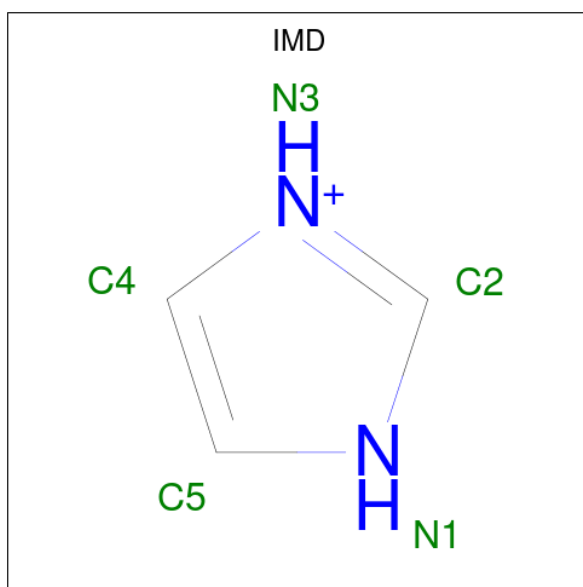
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			7	3	4		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 7 is IMIDAZOLE (CCD ID: IMD) (formula: $C_3H_5N_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	C	1	Total	C	N	0	0
			5	3	2		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	185	Total 185	O 185	0	0
8	B	123	Total 123	O 123	0	0
8	C	232	Total 232	O 232	0	0
8	D	189	Total 189	O 189	0	0
8	E	80	Total 80	O 80	0	0
8	F	91	Total 91	O 91	0	0

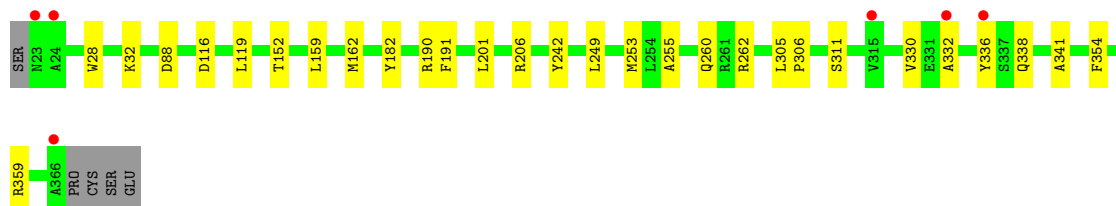
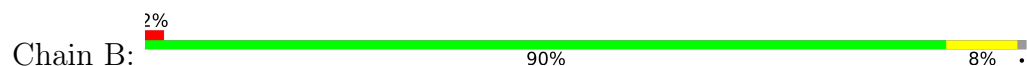
3 Residue-property plots [i](#)

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.

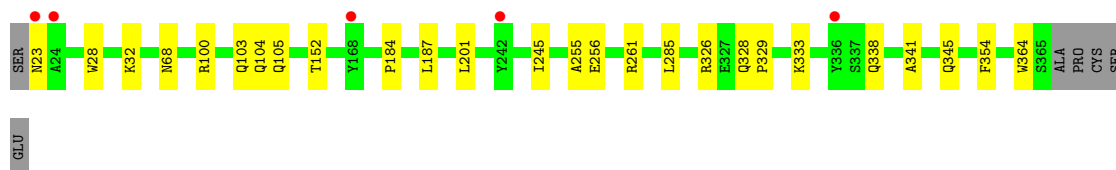
- Molecule 1: Endo-1,4-beta-D-glucanase



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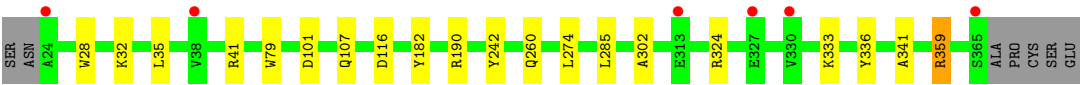


- Molecule 1: Endo-1,4-beta-D-glucanase

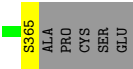
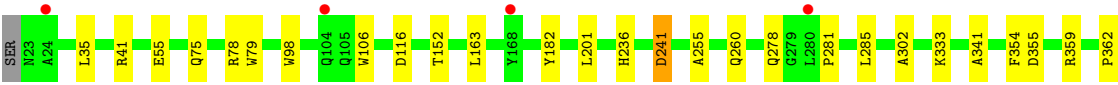
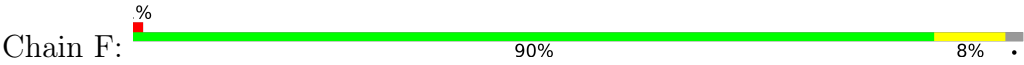


- Molecule 1: Endo-1,4-beta-D-glucanase





● Molecule 1: Endo-1,4-beta-D-glucanase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	126.94Å 129.19Å 165.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.78 – 2.12 40.78 – 2.12	Depositor EDS
% Data completeness (in resolution range)	99.5 (40.78-2.12) 99.8 (40.78-2.12)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.74 (at 2.12Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.163 , 0.198 (Not available) , 0.189	Depositor DCC
R_{free} test set	7719 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	38.7	Xtriage
Anisotropy	0.286	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.005 for k,h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	17169	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.86% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, GOL, IMD, FMT, TRS, MLA

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	0.49	0/2770	0.74	0/3783
1	B	0.45	0/2796	0.74	1/3817 (0.0%)
1	C	0.54	0/2794	0.77	0/3812
1	D	0.50	0/2770	0.74	0/3781
1	E	0.40	0/2750	0.72	0/3751
1	F	0.42	0/2756	0.75	1/3759 (0.0%)
All	All	0.47	0/16636	0.74	2/22703 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	88	ASP	N-CA-C	5.51	116.56	107.20
1	F	281	PRO	O-C-N	5.45	123.71	121.15

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	2685	0	2569	10	0
1	B	2711	0	2609	14	0
1	C	2709	0	2605	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2688	0	2582	9	0
1	E	2673	0	2581	9	0
1	F	2679	0	2580	12	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	10	0	0	1	0
2	D	5	0	0	1	0
2	E	5	0	0	0	0
2	F	5	0	0	1	0
3	A	8	0	12	3	0
3	D	8	0	12	3	0
3	E	8	0	12	3	0
3	F	8	0	12	4	0
4	A	9	0	3	0	0
4	B	6	0	2	0	0
4	C	9	0	3	0	0
4	D	3	0	1	0	0
5	C	7	0	2	0	0
6	C	12	0	16	0	0
6	D	6	0	8	0	0
7	C	5	0	5	1	0
8	A	185	0	0	1	0
8	B	123	0	0	0	0
8	C	232	0	0	1	0
8	D	189	0	0	3	0
8	E	80	0	0	2	0
8	F	91	0	0	2	0
All	All	17169	0	15614	75	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 2.

All (75) 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:B:332:ALA:O	1:B:338:GLN:NE2	2.17	0.78
1:C:326:ARG:NH1	8:C:644:HOH:O	2.27	0.68
1:D:362:PRO:HG2	1:D:365:SER:HB2	1.76	0.67
1:F:236:HIS:NE2	2:F:401:SO4:O2	2.29	0.65
1:A:28:TRP:CE2	1:A:32:LYS:HD2	2.35	0.61
1:F:75:GLN:OE1	1:F:78:ARG:NH1	2.33	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:402:SO4:O3	7:C:409:IMD:H4	2.00	0.61
1:B:119:LEU:HD22	1:B:191:PHE:HE2	1.68	0.59
1:E:116:ASP:HB3	1:E:182:TYR:CG	2.40	0.57
1:F:55:GLU:OE2	3:F:402:TRS:N	2.32	0.56
1:A:35:LEU:HD21	1:A:333:LYS:HA	1.86	0.56
1:E:101:ASP:OD2	1:E:107:GLN:NE2	2.40	0.55
1:A:116:ASP:HB3	1:A:182:TYR:CG	2.42	0.54
1:D:242:TYR:HD2	1:D:336:TYR:CE2	2.26	0.53
3:F:402:TRS:H32	8:F:587:HOH:O	2.07	0.53
1:B:28:TRP:CE2	1:B:32:LYS:HD2	2.43	0.53
1:B:242:TYR:O	1:B:336[B]:TYR:OH	2.23	0.52
1:F:116:ASP:HB3	1:F:182:TYR:CG	2.45	0.52
3:F:402:TRS:H11	8:F:591:HOH:O	2.09	0.52
1:D:48:GLU:O	1:D:100:ARG:NH1	2.43	0.52
3:A:402:TRS:H11	8:A:677:HOH:O	2.10	0.51
1:F:35:LEU:HD21	1:F:333:LYS:HA	1.93	0.51
1:B:116:ASP:HB3	1:B:182:TYR:CG	2.46	0.50
1:C:28:TRP:CE2	1:C:32:LYS:HD3	2.47	0.50
1:D:116:ASP:HB3	1:D:182:TYR:CG	2.47	0.50
1:E:28:TRP:CE2	1:E:32:LYS:HD2	2.46	0.49
1:F:41:ARG:HB3	1:F:79:TRP:CE2	2.48	0.48
1:A:341:ALA:O	1:A:345:GLN:HB2	2.14	0.48
1:E:190:ARG:HD3	1:E:359:ARG:HG3	1.95	0.48
1:D:236:HIS:NE2	2:D:401:SO4:O4	2.46	0.47
1:E:35:LEU:HD21	1:E:333:LYS:HA	1.97	0.47
1:F:355:ASP:OD2	1:F:359:ARG:NH1	2.41	0.47
1:C:103:GLN:HG3	1:C:105:GLN:HG2	1.95	0.47
1:A:28:TRP:CZ2	1:A:32:LYS:HD2	2.50	0.47
1:C:333:LYS:HA	1:C:338:GLN:HE21	1.80	0.47
3:D:403:TRS:H11	8:D:557:HOH:O	2.15	0.46
1:B:262:ARG:HH22	1:B:311:SER:HB3	1.80	0.46
1:B:249:LEU:HG	1:B:253:MSE:HE2	1.98	0.45
1:E:41:ARG:HB3	1:E:79:TRP:CE2	2.51	0.45
3:D:403:TRS:H12	8:D:685:HOH:O	2.16	0.45
1:B:159:LEU:HD21	1:B:162:MSE:HE3	1.99	0.45
3:E:402:TRS:H11	8:E:575:HOH:O	2.18	0.44
1:B:190:ARG:HD3	1:B:359:ARG:HG3	1.98	0.44
1:A:302:ALA:HB1	1:A:341:ALA:HA	2.00	0.44
1:E:116:ASP:HB3	1:E:182:TYR:CD1	2.53	0.44
1:F:98:TRP:CH2	1:F:106:TRP:HB3	2.53	0.44
1:B:242:TYR:HD2	1:B:336[B]:TYR:CE2	2.36	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:302:ALA:HB1	1:E:341:ALA:HA	2.00	0.43
1:C:184:PRO:HG2	1:C:187:LEU:HD22	1.99	0.43
3:E:402:TRS:O3	3:E:402:TRS:O1	2.37	0.43
1:B:330:VAL:HG11	1:B:341:ALA:HB2	2.00	0.43
1:F:362:PRO:HD2	1:F:365:SER:HB3	2.00	0.43
1:A:98:TRP:CH2	1:A:106:TRP:HB3	2.53	0.43
1:D:326:ARG:NH1	8:D:519:HOH:O	2.39	0.43
1:F:182:TYR:OH	3:F:402:TRS:H21	2.19	0.43
1:A:55:GLU:OE2	3:A:402:TRS:O1	2.30	0.43
1:C:341:ALA:O	1:C:345:GLN:HB2	2.19	0.43
3:E:402:TRS:H12	8:E:513:HOH:O	2.19	0.43
1:A:116:ASP:OD2	3:A:402:TRS:O1	2.37	0.42
1:F:255:ALA:HB2	1:F:354:PHE:O	2.20	0.42
1:F:302:ALA:HB1	1:F:341:ALA:HA	2.00	0.42
1:C:68:ASN:HB2	1:C:364:TRP:CD1	2.54	0.42
1:A:68:ASN:HB2	1:A:364:TRP:CD1	2.54	0.42
1:B:206:ARG:HD3	1:B:260:GLN:OE1	2.20	0.42
1:D:115:SER:HB3	1:D:152:THR:HG21	2.02	0.42
1:D:241:ASP:OD1	1:D:284:HIS:ND1	2.43	0.42
1:B:255:ALA:HB2	1:B:354:PHE:O	2.20	0.41
1:D:55:GLU:OE2	3:D:403:TRS:O1	2.27	0.41
1:B:305:LEU:HB2	1:B:306:PRO:HD3	2.02	0.41
1:C:328:GLN:HA	1:C:329:PRO:HD3	1.85	0.41
1:C:256:GLU:HG2	1:C:261:ARG:NH1	2.35	0.41
1:E:242:TYR:HD2	1:E:336:TYR:CE2	2.39	0.41
1:C:255:ALA:HB2	1:C:354:PHE:O	2.21	0.41
1:C:100:ARG:HG2	1:C:104:GLN:HA	2.02	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	71/349 (20%)	70 (99%)	1 (1%)	0	100	100
1	B	283/349 (81%)	279 (99%)	4 (1%)	0	100	100
1	C	343/349 (98%)	338 (98%)	5 (2%)	0	100	100
1	D	343/349 (98%)	339 (99%)	4 (1%)	0	100	100
1	E	340/349 (97%)	332 (98%)	8 (2%)	0	100	100
1	F	341/349 (98%)	336 (98%)	4 (1%)	1 (0%)	36	36
All	All	1721/2094 (82%)	1694 (98%)	26 (2%)	1 (0%)	48	49

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	241	ASP

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	259/267 (97%)	255 (98%)	4 (2%)	57	65
1	B	262/267 (98%)	260 (99%)	2 (1%)	73	80
1	C	263/267 (98%)	259 (98%)	4 (2%)	57	65
1	D	257/267 (96%)	255 (99%)	2 (1%)	73	80
1	E	258/267 (97%)	253 (98%)	5 (2%)	50	57
1	F	259/267 (97%)	252 (97%)	7 (3%)	39	44
All	All	1558/1602 (97%)	1534 (98%)	24 (2%)	57	65

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	119	LEU
1	A	201	LEU
1	A	241	ASP
1	A	285	LEU

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Mol	Chain	Res	Type
1	B	152	THR
1	B	201	LEU
1	C	23	ASN
1	C	152	THR
1	C	201	LEU
1	C	285	LEU
1	D	152	THR
1	D	201	LEU
1	E	260	GLN
1	E	274	LEU
1	E	285	LEU
1	E	324	ARG
1	E	359	ARG
1	F	152	THR
1	F	163	LEU
1	F	201	LEU
1	F	241	ASP
1	F	260	GLN
1	F	278	GLN
1	F	285	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	295	HIS
1	B	92	HIS
1	B	236	HIS
1	B	238	ASN
1	B	338	GLN
1	C	23	ASN
1	C	83	ASN
1	C	174	GLN
1	C	238	ASN
1	C	278	GLN
1	E	92	HIS
1	E	314	HIS
1	F	284	HIS
1	F	314	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

25 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	$\# Z > 2$
3	TRS	E	402	-	7,7,7	0.53	0	9,9,9	0.81	0
4	FMT	A	403	-	2,2,2	0.63	0	1,1,1	0.04	0
6	GOL	C	405	-	5,5,5	0.36	0	5,5,5	0.31	0
4	FMT	D	404	-	2,2,2	0.65	0	1,1,1	0.14	0
4	FMT	C	407	-	2,2,2	0.69	0	1,1,1	0.16	0
2	SO4	E	401	-	4,4,4	0.23	0	6,6,6	0.12	0
5	MLA	C	403	-	6,6,6	1.17	0	7,7,7	1.30	0
4	FMT	B	403	-	2,2,2	0.69	0	1,1,1	0.16	0
2	SO4	B	401	-	4,4,4	0.22	0	6,6,6	0.14	0
2	SO4	C	402	-	4,4,4	0.22	0	6,6,6	0.14	0
6	GOL	D	402	-	5,5,5	0.40	0	5,5,5	0.63	0
4	FMT	A	404	-	2,2,2	0.70	0	1,1,1	0.20	0
3	TRS	D	403	-	7,7,7	0.42	0	9,9,9	1.02	1 (11%)
2	SO4	F	401	-	4,4,4	0.23	0	6,6,6	0.12	0
4	FMT	C	406	-	2,2,2	0.68	0	1,1,1	0.13	0
2	SO4	D	401	-	4,4,4	0.27	0	6,6,6	0.10	0
3	TRS	F	402	-	7,7,7	0.35	0	9,9,9	1.96	2 (22%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	FMT	B	402	-	2,2,2	0.71	0	1,1,1	0.15	0
7	IMD	C	409	-	5,5,5	0.68	0	5,5,5	0.48	0
3	TRS	A	402	-	7,7,7	0.65	0	9,9,9	1.27	1 (11%)
4	FMT	A	405	-	2,2,2	0.63	0	1,1,1	0.09	0
4	FMT	C	408	-	2,2,2	0.65	0	1,1,1	0.09	0
2	SO4	A	401	-	4,4,4	0.25	0	6,6,6	0.07	0
6	GOL	C	404	-	5,5,5	0.32	0	5,5,5	0.51	0
2	SO4	C	401	-	4,4,4	0.32	0	6,6,6	0.18	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	TRS	F	402	-	-	8/9/9/9	-
3	TRS	E	402	-	-	4/9/9/9	-
7	IMD	C	409	-	-	-	0/1/1/1
3	TRS	A	402	-	-	3/9/9/9	-
3	TRS	D	403	-	-	3/9/9/9	-
5	MLA	C	403	-	-	4/4/4/4	-
6	GOL	C	405	-	-	2/4/4/4	-
6	GOL	C	404	-	-	0/4/4/4	-
6	GOL	D	402	-	-	4/4/4/4	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	402	TRS	O1-C1-C	3.66	121.09	110.88
3	F	402	TRS	C2-C-C1	3.31	119.48	110.66
3	A	402	TRS	O1-C1-C	2.71	118.43	110.88
3	D	403	TRS	O1-C1-C	2.07	116.66	110.88

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	403	TRS	C1-C-C3-O3

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Mol	Chain	Res	Type	Atoms
3	D	403	TRS	C2-C-C3-O3
3	F	402	TRS	C2-C-C1-O1
3	F	402	TRS	C3-C-C1-O1
3	F	402	TRS	N-C-C1-O1
3	F	402	TRS	N-C-C3-O3
6	C	405	GOL	O1-C1-C2-C3
6	D	402	GOL	O1-C1-C2-C3
6	D	402	GOL	C1-C2-C3-O3
6	C	405	GOL	O1-C1-C2-O2
6	D	402	GOL	O1-C1-C2-O2
3	E	402	TRS	C3-C-C1-O1
3	F	402	TRS	C2-C-C3-O3
6	D	402	GOL	O2-C2-C3-O3
3	A	402	TRS	C3-C-C2-O2
3	D	403	TRS	N-C-C3-O3
3	E	402	TRS	N-C-C1-O1
3	F	402	TRS	C1-C-C2-O2
3	F	402	TRS	N-C-C2-O2
5	C	403	MLA	O1A-C1-C2-C3
3	A	402	TRS	C1-C-C2-O2
3	E	402	TRS	C2-C-C1-O1
3	F	402	TRS	C3-C-C2-O2
5	C	403	MLA	O1B-C1-C2-C3
5	C	403	MLA	C1-C2-C3-O3A
3	A	402	TRS	N-C-C2-O2
5	C	403	MLA	C1-C2-C3-O3B
3	E	402	TRS	C3-C-C2-O2

There are no ring outliers.

8 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	402	TRS	3	0
2	C	402	SO4	1	0
3	D	403	TRS	3	0
2	F	401	SO4	1	0
2	D	401	SO4	1	0
3	F	402	TRS	4	0
7	C	409	IMD	1	0
3	A	402	TRS	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 ⓘ

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	339/349 (97%)	-0.49	5 (1%) 72 74	21, 34, 59, 85	2 (0%)
1	B	341/349 (97%)	-0.21	6 (1%) 67 70	17, 39, 77, 105	3 (0%)
1	C	340/349 (97%)	-0.58	5 (1%) 72 74	18, 27, 56, 94	2 (0%)
1	D	341/349 (97%)	-0.49	6 (1%) 67 70	21, 32, 64, 105	1 (0%)
1	E	339/349 (97%)	-0.02	6 (1%) 67 70	31, 51, 86, 107	0
1	F	340/349 (97%)	-0.11	4 (1%) 76 79	30, 47, 77, 114	0
All	All	2040/2094 (97%)	-0.32	32 (1%) 70 73	17, 38, 74, 114	8 (0%)

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	336[A]	TYR	4.2
1	D	103	GLN	3.9
1	B	24	ALA	3.8
1	D	104	GLN	3.8
1	D	366	ALA	3.8
1	A	333	LYS	3.5
1	A	24	ALA	3.4
1	A	332	ALA	3.3
1	C	242[A]	TYR	3.0
1	E	365	SER	3.0
1	B	366	ALA	2.9
1	F	104	GLN	2.8
1	F	280	LEU	2.7
1	A	168[A]	TYR	2.7
1	C	23	ASN	2.7
1	E	327	GLU	2.7
1	D	168[A]	TYR	2.6
1	E	313	GLU	2.5
1	B	23	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	168	TYR	2.5
1	D	24	ALA	2.5
1	D	332	ALA	2.4
1	E	330	VAL	2.3
1	C	336	TYR	2.2
1	E	24	ALA	2.2
1	C	24	ALA	2.2
1	B	332	ALA	2.1
1	B	315	VAL	2.1
1	E	38	VAL	2.1
1	A	331	GLU	2.0
1	F	24	ALA	2.1
1	F	168	TYR	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(Å ²)	Q<0.9
3	TRS	F	402	8/8	0.74	0.16	37,44,51,53	0
2	SO4	C	401	5/5	0.76	0.20	36,56,63,65	5
2	SO4	E	401	5/5	0.78	0.15	82,87,91,92	5
4	FMT	A	405	3/3	0.78	0.16	54,54,55,62	0
4	FMT	A	404	3/3	0.79	0.16	68,68,71,73	0
2	SO4	B	401	5/5	0.81	0.16	77,80,82,83	5
4	FMT	C	408	3/3	0.81	0.14	55,55,57,62	0
6	GOL	D	402	6/6	0.81	0.18	47,50,52,56	0
7	IMD	C	409	5/5	0.82	0.15	62,65,67,68	0
4	FMT	B	403	3/3	0.83	0.13	65,65,66,70	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	FMT	D	404	3/3	0.84	0.13	48,48,56,59	0
2	SO4	D	401	5/5	0.85	0.14	72,73,75,76	5
3	TRS	D	403	8/8	0.86	0.11	30,45,52,68	0
4	FMT	C	407	3/3	0.86	0.11	70,70,73,73	0
4	FMT	B	402	3/3	0.86	0.21	63,63,67,67	0
2	SO4	F	401	5/5	0.87	0.11	81,84,85,85	5
5	MLA	C	403	7/7	0.87	0.12	53,68,70,77	0
6	GOL	C	405	6/6	0.88	0.14	62,65,71,71	0
3	TRS	E	402	8/8	0.88	0.09	33,38,41,53	0
2	SO4	A	401	5/5	0.88	0.09	82,82,84,84	5
3	TRS	A	402	8/8	0.89	0.12	25,36,43,58	0
4	FMT	A	403	3/3	0.92	0.17	40,40,41,45	0
4	FMT	C	406	3/3	0.92	0.08	40,40,54,57	0
6	GOL	C	404	6/6	0.92	0.12	30,43,53,57	0
2	SO4	C	402	5/5	0.95	0.09	39,47,53,61	5

6.5 Other polymers [i](#)

There are no such residues in this entry.