



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:49 PM UTC

PDB ID : 6NXD / pdb_00006nxd
Title : TYPE I L-ASPARAGINASE FROM ESCHERICHIA COLI IN COMPLEX
WITH CITRATE AT PH 4
Authors : Lubkowski, J.; Wlodawer, A.
Deposited on : 2019-02-08
Resolution : 1.90 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

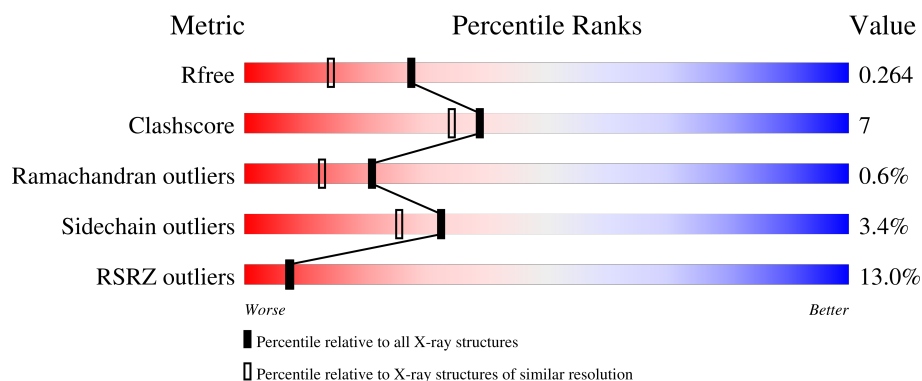
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	358	10% 72% 12% • 15%
1	B	358	12% 78% 11% • 9%
1	C	358	13% 72% 14% • 13%
1	D	358	11% 76% 12% • 9%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 10411 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 L-asparaginase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	305	Total	C	N	O	S	0	0	0
			2336	1482	398	446	10			
1	B	325	Total	C	N	O	S	0	0	0
			2491	1579	428	472	12			
1	C	311	Total	C	N	O	S	0	1	0
			2394	1522	406	454	12			
1	D	324	Total	C	N	O	S	0	1	0
			2480	1574	423	472	11			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP P0A962
A	-18	GLY	-	expression tag	UNP P0A962
A	-17	SER	-	expression tag	UNP P0A962
A	-16	SER	-	expression tag	UNP P0A962
A	-15	HIS	-	expression tag	UNP P0A962
A	-14	HIS	-	expression tag	UNP P0A962
A	-13	HIS	-	expression tag	UNP P0A962
A	-12	HIS	-	expression tag	UNP P0A962
A	-11	HIS	-	expression tag	UNP P0A962
A	-10	HIS	-	expression tag	UNP P0A962
A	-9	SER	-	expression tag	UNP P0A962
A	-8	SER	-	expression tag	UNP P0A962
A	-7	GLY	-	expression tag	UNP P0A962
A	-6	LEU	-	expression tag	UNP P0A962
A	-5	VAL	-	expression tag	UNP P0A962
A	-4	PRO	-	expression tag	UNP P0A962
A	-3	ARG	-	expression tag	UNP P0A962
A	-2	GLY	-	expression tag	UNP P0A962
A	-1	SER	-	expression tag	UNP P0A962
A	0	HIS	-	expression tag	UNP P0A962
B	-19	MET	-	initiating methionine	UNP P0A962

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	expression tag	UNP P0A962
B	-17	SER	-	expression tag	UNP P0A962
B	-16	SER	-	expression tag	UNP P0A962
B	-15	HIS	-	expression tag	UNP P0A962
B	-14	HIS	-	expression tag	UNP P0A962
B	-13	HIS	-	expression tag	UNP P0A962
B	-12	HIS	-	expression tag	UNP P0A962
B	-11	HIS	-	expression tag	UNP P0A962
B	-10	HIS	-	expression tag	UNP P0A962
B	-9	SER	-	expression tag	UNP P0A962
B	-8	SER	-	expression tag	UNP P0A962
B	-7	GLY	-	expression tag	UNP P0A962
B	-6	LEU	-	expression tag	UNP P0A962
B	-5	VAL	-	expression tag	UNP P0A962
B	-4	PRO	-	expression tag	UNP P0A962
B	-3	ARG	-	expression tag	UNP P0A962
B	-2	GLY	-	expression tag	UNP P0A962
B	-1	SER	-	expression tag	UNP P0A962
B	0	HIS	-	expression tag	UNP P0A962
C	-19	MET	-	initiating methionine	UNP P0A962
C	-18	GLY	-	expression tag	UNP P0A962
C	-17	SER	-	expression tag	UNP P0A962
C	-16	SER	-	expression tag	UNP P0A962
C	-15	HIS	-	expression tag	UNP P0A962
C	-14	HIS	-	expression tag	UNP P0A962
C	-13	HIS	-	expression tag	UNP P0A962
C	-12	HIS	-	expression tag	UNP P0A962
C	-11	HIS	-	expression tag	UNP P0A962
C	-10	HIS	-	expression tag	UNP P0A962
C	-9	SER	-	expression tag	UNP P0A962
C	-8	SER	-	expression tag	UNP P0A962
C	-7	GLY	-	expression tag	UNP P0A962
C	-6	LEU	-	expression tag	UNP P0A962
C	-5	VAL	-	expression tag	UNP P0A962
C	-4	PRO	-	expression tag	UNP P0A962
C	-3	ARG	-	expression tag	UNP P0A962
C	-2	GLY	-	expression tag	UNP P0A962
C	-1	SER	-	expression tag	UNP P0A962
C	0	HIS	-	expression tag	UNP P0A962
D	-19	MET	-	initiating methionine	UNP P0A962
D	-18	GLY	-	expression tag	UNP P0A962
D	-17	SER	-	expression tag	UNP P0A962

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	expression tag	UNP P0A962
D	-15	HIS	-	expression tag	UNP P0A962
D	-14	HIS	-	expression tag	UNP P0A962
D	-13	HIS	-	expression tag	UNP P0A962
D	-12	HIS	-	expression tag	UNP P0A962
D	-11	HIS	-	expression tag	UNP P0A962
D	-10	HIS	-	expression tag	UNP P0A962
D	-9	SER	-	expression tag	UNP P0A962
D	-8	SER	-	expression tag	UNP P0A962
D	-7	GLY	-	expression tag	UNP P0A962
D	-6	LEU	-	expression tag	UNP P0A962
D	-5	VAL	-	expression tag	UNP P0A962
D	-4	PRO	-	expression tag	UNP P0A962
D	-3	ARG	-	expression tag	UNP P0A962
D	-2	GLY	-	expression tag	UNP P0A962
D	-1	SER	-	expression tag	UNP P0A962
D	0	HIS	-	expression tag	UNP P0A962

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

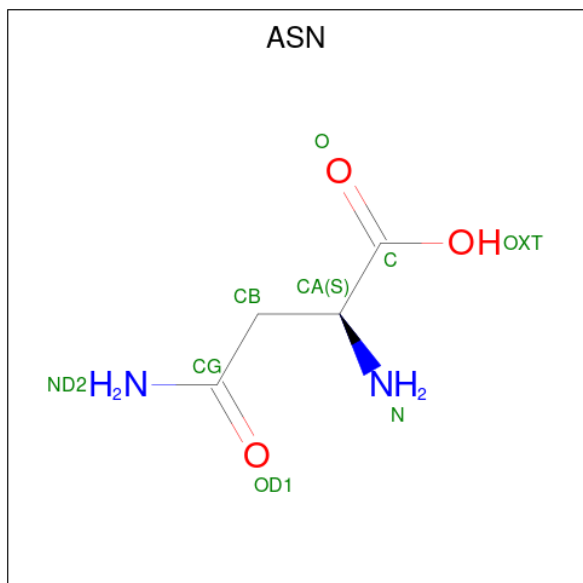
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	4	Total Cl 4 4	0	0
2	B	4	Total Cl 4 4	0	0
2	C	1	Total Cl 1 1	0	0
2	D	1	Total Cl 1 1	0	0

- Molecule 3 is CITRIC ACID (CCD ID: CIT) (formula: C₆H₈O₇).



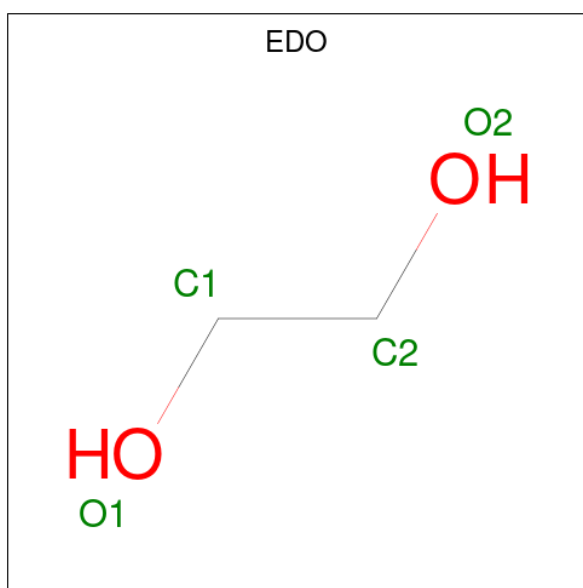
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		
3	C	1	Total	C	O	0	0
			12	6	6		
3	D	1	Total	C	O	0	0
			12	6	6		

- Molecule 4 is ASPARAGINE (CCD ID: ASN) (formula: $C_4H_8N_2O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N O 9 4 2 3	0	0
4	B	1	Total C N O 9 4 2 3	0	0
4	C	1	Total C N O 9 4 2 3	0	0
4	D	1	Total C N O 9 4 2 3	0	0

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

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

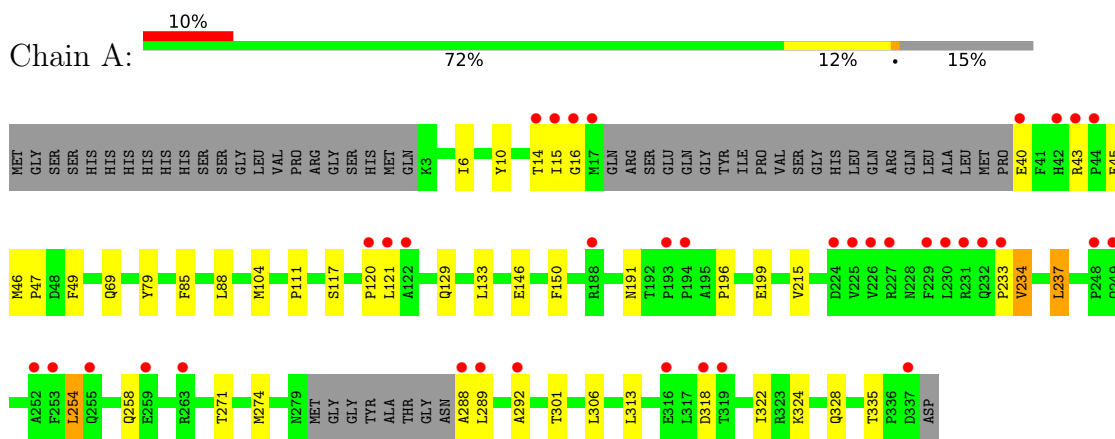
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	129	Total O 129 129	0	0
6	B	150	Total O 150 150	0	0
6	C	124	Total O 124 124	0	0
6	D	161	Total O 161 161	0	0

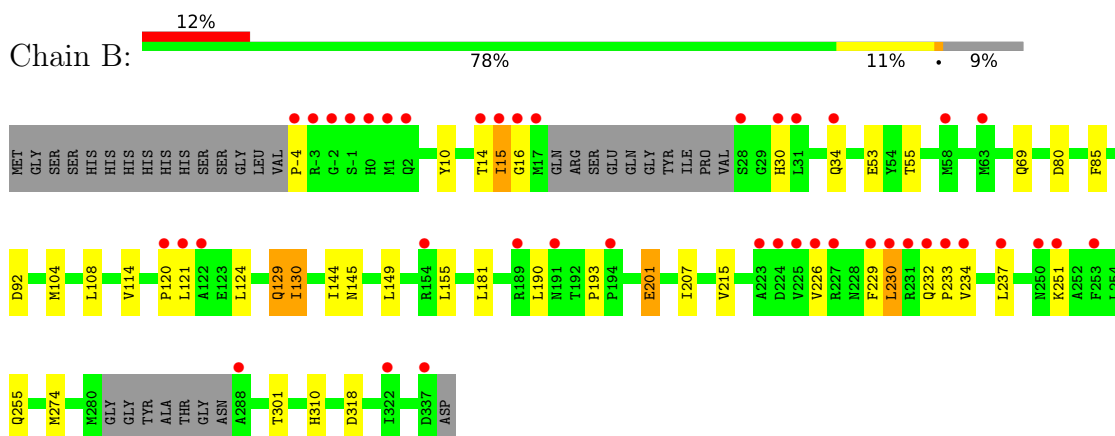
3 Residue-property plots [i](#)

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.

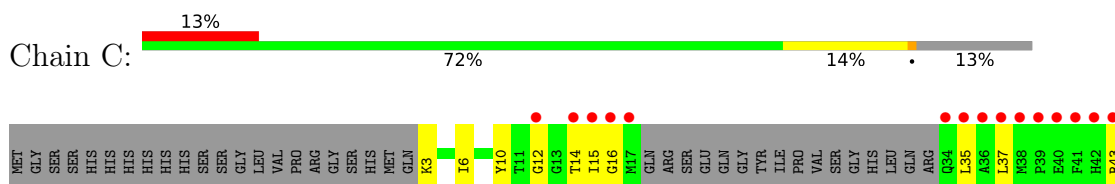
• Molecule 1: L-asparaginase 1

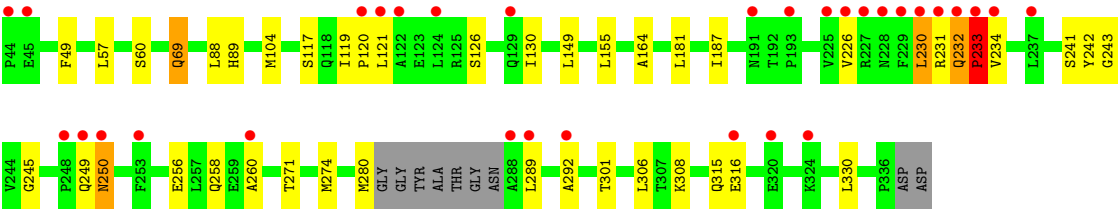


• Molecule 1: L-asparaginase 1

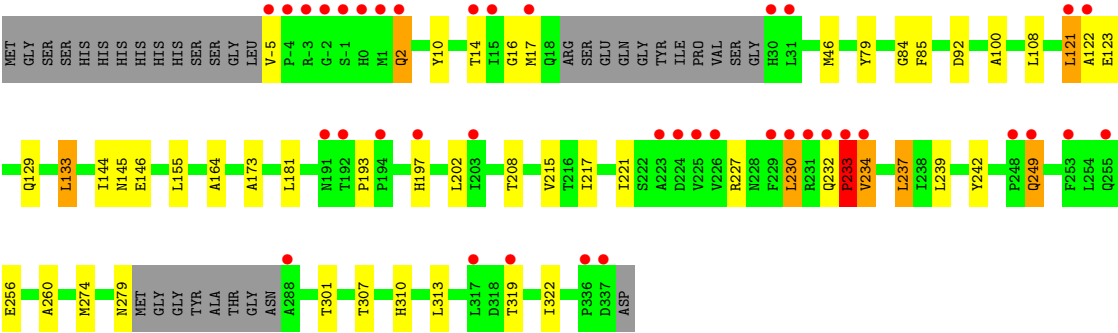
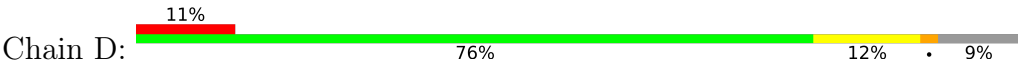


• Molecule 1: L-asparaginase 1





● Molecule 1: L-asparaginase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	90.31Å 89.76Å 93.08Å 90.00° 117.03° 90.00°	Depositor
Resolution (Å)	45.17 – 1.90 45.17 – 1.90	Depositor EDS
% Data completeness (in resolution range)	90.7 (45.17-1.90) 90.7 (45.17-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.46 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.8.0232	Depositor
R, R_{free}	0.203 , 0.244 (Not available) , 0.264	Depositor DCC
R_{free} test set	4749 reflections (4.56%)	wwPDB-VP
Wilson B-factor (Å ²)	18.3	Xtriage
Anisotropy	0.121	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 44.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.148 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10411	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% 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: CL, CIT, EDO

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	1.13	0/2387	1.43	5/3251 (0.2%)
1	B	1.12	0/2546	1.42	3/3465 (0.1%)
1	C	1.10	0/2447	1.40	2/3332 (0.1%)
1	D	1.12	2/2537 (0.1%)	1.42	6/3457 (0.2%)
All	All	1.12	2/9917 (0.0%)	1.42	16/13505 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	146	GLU	C-O	6.98	1.32	1.23
1	D	310	HIS	CE1-NE2	5.78	1.38	1.32

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	271	THR	CA-CB-OG1	-8.23	97.26	109.60
1	C	271	THR	CA-CB-OG1	-7.98	97.63	109.60
1	C	233	PRO	N-CA-CB	-6.60	96.32	103.25
1	A	335	THR	CB-CA-C	6.22	116.33	110.17
1	D	208	THR	CB-CA-C	6.16	119.86	109.32
1	A	196	PRO	N-CA-C	-6.12	101.61	111.03
1	D	173	ALA	CA-C-N	5.42	129.71	123.15
1	D	173	ALA	C-N-CA	5.42	129.71	123.15
1	D	233	PRO	N-CA-CB	-5.31	97.68	103.25
1	B	193	PRO	N-CA-CB	5.28	105.97	103.22
1	A	150	PHE	CA-CB-CG	-5.26	108.54	113.80
1	B	318	ASP	CA-CB-CG	5.15	117.75	112.60
1	A	111	PRO	CB-CA-C	-5.11	104.41	111.21
1	D	121	LEU	CA-C-N	5.07	127.03	120.44
1	D	121	LEU	C-N-CA	5.07	127.03	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	130	ILE	CA-C-O	-5.05	115.50	120.85

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	2336	0	2286	28	0
1	B	2491	0	2453	32	0
1	C	2394	0	2369	34	0
1	D	2480	0	2435	42	0
2	A	4	0	0	1	0
2	B	4	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	12	0	5	1	0
3	B	12	0	5	1	0
3	C	12	0	5	0	0
3	D	12	0	5	3	0
4	A	9	0	5	0	0
4	B	9	0	5	0	0
4	C	9	0	5	0	0
4	D	9	0	5	0	0
5	A	12	0	18	3	0
5	B	12	0	18	1	0
5	C	8	0	12	1	0
5	D	20	0	30	1	0
6	A	129	0	0	2	0
6	B	150	0	0	4	0
6	C	124	0	0	0	0
6	D	161	0	0	2	0
All	All	10411	0	9661	133	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 7.

All (133) 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:10:TYR:OH	1:B:16:GLY:HA3	1.80	0.82
1:B:274:MET:HE3	1:D:274:MET:HB2	1.61	0.81
1:C:232:GLN:H	1:C:233:PRO:HD3	1.50	0.76
1:D:164:ALA:HB2	1:D:274:MET:HE1	1.73	0.70
1:D:-5:VAL:HA	1:D:144:ILE:HG12	1.73	0.69
1:A:318:ASP:OD1	6:A:1101:HOH:O	2.11	0.69
1:B:201:GLU:CD	6:B:1101:HOH:O	2.36	0.67
1:D:14:THR:HA	1:D:17:MET:CE	2.24	0.67
1:B:229:PHE:HB3	1:B:234:VAL:HG21	1.77	0.66
1:B:215:VAL:CG2	1:B:237:LEU:HD21	2.26	0.66
1:C:181:LEU:HD21	1:C:187:ILE:HG23	1.77	0.65
1:B:274:MET:HE1	6:D:1236:HOH:O	1.97	0.65
1:B:207:ILE:HG23	1:B:310:HIS:HB3	1.80	0.63
1:D:14:THR:HA	1:D:17:MET:HE3	1.79	0.63
1:B:274:MET:CE	1:D:274:MET:HB2	2.28	0.63
1:D:46:MET:HE2	1:D:133:LEU:HD22	1.82	0.62
1:C:10:TYR:OH	1:C:16:GLY:HA3	2.00	0.62
1:D:234:VAL:HG11	1:D:237:LEU:HD13	1.80	0.62
1:D:227:ARG:N	1:D:256:GLU:OE2	2.33	0.61
1:A:40:GLU:O	1:A:43:ARG:HG2	2.01	0.61
1:D:319:THR:HG22	6:D:1244:HOH:O	2.01	0.60
1:B:92:ASP:OD2	3:B:1005:CIT:H21	2.01	0.59
1:D:-5:VAL:CB	1:D:144:ILE:HD11	2.32	0.59
1:C:241[A]:SER:OG	1:C:245:GLY:HA2	2.04	0.58
1:B:215:VAL:HG21	1:B:237:LEU:HD21	1.85	0.58
1:B:129:GLN:HB2	6:B:1156:HOH:O	2.04	0.57
1:D:17:MET:HE1	3:D:1002:CIT:O7	2.05	0.57
1:D:10:TYR:OH	1:D:16:GLY:HA3	2.04	0.56
1:B:-4:PRO:HD2	1:B:144:ILE:HG12	1.88	0.56
1:A:10:TYR:OH	1:A:16:GLY:HA3	2.04	0.56
1:D:215[A]:VAL:HG12	1:D:239:LEU:HD23	1.87	0.56
1:A:301:THR:HA	5:A:1008:EDO:H12	1.88	0.55
1:A:15:ILE:HD13	1:A:88:LEU:HD22	1.88	0.55
1:A:215:VAL:HG23	1:A:237:LEU:CD1	2.37	0.55
1:C:232:GLN:H	1:C:233:PRO:CD	2.19	0.55
1:C:230:LEU:HD21	1:C:260:ALA:HB2	1.89	0.55
1:A:199:GLU:OE2	6:A:1102:HOH:O	2.19	0.54
1:A:15:ILE:O	1:A:121:LEU:CB	2.56	0.53
1:C:15:ILE:HD13	1:C:88:LEU:HD22	1.91	0.53
1:A:274:MET:HE3	1:C:274:MET:HB2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:215:VAL:HG23	1:B:237:LEU:HD21	1.92	0.52
1:D:230:LEU:HD11	1:D:260:ALA:HA	1.92	0.52
1:A:121:LEU:CB	1:A:129:GLN:HE22	2.23	0.52
1:C:14:THR:HG23	1:C:120:PRO:HD3	1.92	0.52
1:B:226:VAL:O	1:B:230:LEU:HD23	2.09	0.51
1:D:215[A]:VAL:HG11	1:D:239:LEU:HD22	1.93	0.51
1:B:233:PRO:O	1:B:234:VAL:HG23	2.10	0.51
1:C:230:LEU:HD21	1:C:260:ALA:CB	2.41	0.51
1:A:45:GLU:HB2	1:B:124:LEU:HD22	1.93	0.50
1:A:215:VAL:CG2	1:A:237:LEU:HD11	2.41	0.50
1:A:14:THR:HG22	1:A:120:PRO:N	2.26	0.50
1:C:15:ILE:HA	1:C:117:SER:OG	2.12	0.50
1:A:234:VAL:HG11	1:A:237:LEU:HD23	1.92	0.50
1:C:258:GLN:NE2	1:C:292:ALA:HA	2.27	0.49
1:C:15:ILE:O	1:C:121:LEU:HB2	2.12	0.49
1:C:6:ILE:O	1:C:49:PHE:HA	2.12	0.49
1:B:121:LEU:HD11	1:B:129:GLN:HG3	1.95	0.49
1:D:145:ASN:HB2	1:D:197:HIS:HE1	1.76	0.49
1:B:80:ASP:OD1	6:B:1101:HOH:O	2.20	0.49
1:C:60:SER:HB2	1:C:89:HIS:CE1	2.47	0.49
1:C:232:GLN:N	1:C:233:PRO:CD	2.76	0.49
1:C:230:LEU:HD11	1:C:260:ALA:HA	1.95	0.48
1:B:14:THR:HG23	1:B:120:PRO:HD3	1.95	0.48
1:C:301:THR:HA	5:C:1004:EDO:H12	1.95	0.48
1:B:155:LEU:HB3	1:B:181:LEU:HB3	1.95	0.48
1:C:250:ASN:OD1	1:C:250:ASN:C	2.56	0.48
1:D:215[A]:VAL:CG1	1:D:239:LEU:CD2	2.91	0.48
1:C:280:MET:CE	1:C:289:LEU:HB2	2.44	0.48
1:D:215[A]:VAL:CG1	1:D:239:LEU:HD22	2.43	0.48
1:A:14:THR:CG2	1:A:120:PRO:HD3	2.44	0.47
1:A:15:ILE:HA	1:A:117:SER:OG	2.13	0.47
1:C:14:THR:CG2	1:C:120:PRO:HD3	2.45	0.47
1:A:215:VAL:CG2	1:A:237:LEU:CD1	2.92	0.47
1:D:2:GLN:HE21	1:D:2:GLN:HB2	1.54	0.47
1:C:15:ILE:CA	1:C:117:SER:OG	2.63	0.47
1:A:46:MET:HE3	1:A:47:PRO:HD2	1.97	0.47
1:D:279:ASN:C	1:D:279:ASN:HD22	2.22	0.47
1:D:46:MET:CE	1:D:133:LEU:HD22	2.44	0.47
1:A:79:TYR:HA	1:A:85:PHE:CZ	2.50	0.46
1:D:14:THR:HA	1:D:17:MET:HE2	1.93	0.46
1:A:191:ASN:HD22	1:D:193:PRO:HD3	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:84:GLY:C	1:D:85:PHE:CD1	2.94	0.46
1:A:258:GLN:OE1	1:A:292:ALA:HA	2.16	0.46
1:C:149:LEU:C	1:C:149:LEU:HD23	2.41	0.45
1:D:121:LEU:HD11	1:D:129:GLN:HG3	1.98	0.45
1:D:233:PRO:O	1:D:234:VAL:HG23	2.17	0.45
1:D:85:PHE:CD1	1:D:85:PHE:N	2.85	0.45
1:D:108:LEU:O	1:D:145:ASN:HB3	2.17	0.44
1:C:43:ARG:NH2	1:D:122:ALA:O	2.51	0.44
1:D:145:ASN:HB2	1:D:197:HIS:CE1	2.52	0.44
1:D:234:VAL:HG11	1:D:237:LEU:CD1	2.44	0.44
2:A:1003:CL:CL	5:A:1007:EDO:O1	2.73	0.43
1:A:324:LYS:O	1:A:328:GLN:HG3	2.18	0.43
1:D:100:ALA:HA	1:D:307:THR:HB	2.00	0.43
1:B:30:HIS:O	1:B:34:GLN:HG2	2.17	0.43
1:C:226:VAL:CG1	1:C:256:GLU:HB3	2.48	0.43
1:D:301:THR:HA	5:D:1004:EDO:H21	2.00	0.43
1:A:146:GLU:HG2	5:A:1007:EDO:H22	2.00	0.43
1:D:164:ALA:CB	1:D:274:MET:HE1	2.46	0.43
1:D:92:ASP:OD2	3:D:1002:CIT:H21	2.19	0.43
1:D:217:ILE:HG22	1:D:242:TYR:CZ	2.54	0.43
1:A:313:LEU:HD23	1:A:322:ILE:HD13	2.00	0.43
1:B:14:THR:O	1:B:120:PRO:HA	2.18	0.43
1:D:155:LEU:HB3	1:D:181:LEU:HB3	2.01	0.43
1:D:215[A]:VAL:HG12	1:D:239:LEU:CD2	2.50	0.42
1:A:6:ILE:O	1:A:49:PHE:HA	2.19	0.42
1:B:69:GLN:HA	1:B:104:MET:HE1	2.01	0.42
1:B:14:THR:CG2	1:B:120:PRO:HD3	2.48	0.42
1:C:308:LYS:HE3	1:C:330:LEU:HD12	2.01	0.42
1:D:249:GLN:HE21	1:D:249:GLN:CA	2.32	0.42
1:B:85:PHE:CD1	1:B:85:PHE:N	2.88	0.42
1:C:12:GLY:HA2	1:C:57:LEU:HD22	2.00	0.42
1:C:164:ALA:CB	1:C:274:MET:HE1	2.50	0.42
1:B:15:ILE:HG13	6:B:1170:HOH:O	2.19	0.41
1:B:121:LEU:CD1	1:B:129:GLN:HG3	2.51	0.41
1:D:313:LEU:HD23	1:D:322:ILE:HD13	2.02	0.41
1:B:114:VAL:O	1:B:149:LEU:HA	2.21	0.41
1:B:108:LEU:O	1:B:145:ASN:HB3	2.21	0.41
1:B:130:ILE:HD12	1:B:130:ILE:HA	1.91	0.41
1:B:10:TYR:HB3	1:B:53:GLU:HA	2.02	0.41
1:C:119:ILE:HB	1:C:126:SER:HA	2.02	0.41
1:B:301:THR:HA	5:B:1007:EDO:H21	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:14:THR:HG22	1:C:120:PRO:N	2.36	0.41
1:C:242:TYR:O	1:C:243:GLY:C	2.63	0.41
1:A:254:LEU:HD21	1:A:288:ALA:HB1	2.03	0.41
1:C:130:ILE:HD12	1:C:130:ILE:HA	1.92	0.40
3:D:1002:CIT:O2	3:D:1002:CIT:C6	2.69	0.40
3:A:1005:CIT:C6	3:A:1005:CIT:O2	2.69	0.40
1:C:155:LEU:HB3	1:C:181:LEU:HB3	2.03	0.40
1:A:14:THR:CG2	1:A:120:PRO:CD	3.00	0.40
1:A:69:GLN:HA	1:A:104:MET:HE1	2.03	0.40
1:D:79:TYR:HA	1:D:85:PHE:CZ	2.56	0.40
1:C:69:GLN:HA	1:C:104:MET:HE1	2.02	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	299/358 (84%)	287 (96%)	11 (4%)	1 (0%)	36	29
1	B	319/358 (89%)	308 (97%)	10 (3%)	1 (0%)	36	29
1	C	306/358 (86%)	294 (96%)	9 (3%)	3 (1%)	12	5
1	D	319/358 (89%)	309 (97%)	8 (2%)	2 (1%)	21	13
All	All	1243/1432 (87%)	1198 (96%)	38 (3%)	7 (1%)	21	13

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	231	ARG
1	D	232	GLN
1	B	232	GLN
1	A	233	PRO

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Mol	Chain	Res	Type
1	C	233	PRO
1	D	233	PRO
1	C	232	GLN

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	250/298 (84%)	244 (98%)	6 (2%)	43	38
1	B	267/298 (90%)	259 (97%)	8 (3%)	36	30
1	C	258/298 (87%)	246 (95%)	12 (5%)	23	15
1	D	265/298 (89%)	256 (97%)	9 (3%)	32	25
All	All	1040/1192 (87%)	1005 (97%)	35 (3%)	32	25

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	133	LEU
1	A	234	VAL
1	A	237	LEU
1	A	254	LEU
1	A	289	LEU
1	A	306	LEU
1	B	15	ILE
1	B	55	THR
1	B	129	GLN
1	B	190	LEU
1	B	201	GLU
1	B	230	LEU
1	B	251	LYS
1	B	255	GLN
1	C	3	LYS
1	C	35	LEU
1	C	37	LEU
1	C	69	GLN

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Mol	Chain	Res	Type
1	C	230	LEU
1	C	233	PRO
1	C	234	VAL
1	C	249	GLN
1	C	250	ASN
1	C	306	LEU
1	C	315	GLN
1	C	316	GLU
1	D	2	GLN
1	D	123	GLU
1	D	133	LEU
1	D	202	LEU
1	D	221	ILE
1	D	230	LEU
1	D	234	VAL
1	D	237	LEU
1	D	249	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	129	GLN
1	A	145	ASN
1	A	153	ASN
1	A	255	GLN
1	B	129	GLN
1	B	152	ASN
1	B	153	ASN
1	B	246	ASN
1	B	258	GLN
1	B	328	GLN
1	C	34	GLN
1	C	145	ASN
1	C	186	HIS
1	C	258	GLN
1	D	2	GLN
1	D	152	ASN
1	D	153	ASN
1	D	249	GLN
1	D	250	ASN
1	D	255	GLN
1	D	272	GLN

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Mol	Chain	Res	Type
1	D	279	ASN
1	D	315	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 31 ligands modelled in this entry, 10 are monoatomic - leaving 21 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	EDO	A	1007	-	3,3,3	0.10	0	2,2,2	0.13	0
3	CIT	B	1005	1	11,11,12	1.49	3 (27%)	12,15,17	1.23	2 (16%)
4	ASN	B	1006	-	7,8,8	1.16	1 (14%)	6,10,10	0.47	0
5	EDO	A	1008	-	3,3,3	0.19	0	2,2,2	0.35	0
5	EDO	D	1006	-	3,3,3	0.56	0	2,2,2	0.43	0
5	EDO	D	1007	-	3,3,3	0.45	0	2,2,2	0.31	0
4	ASN	A	1006	-	7,8,8	0.75	0	6,10,10	0.76	0
3	CIT	A	1005	1	11,11,12	1.09	1 (9%)	14,15,17	1.28	3 (21%)
5	EDO	D	1004	-	3,3,3	0.46	0	2,2,2	0.52	0
5	EDO	C	1004	-	3,3,3	0.27	0	2,2,2	0.40	0
5	EDO	A	1009	-	3,3,3	0.34	0	2,2,2	0.51	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	B	1009	-	3,3,3	0.32	0	2,2,2	0.48	0
3	CIT	C	1002	1	11,11,12	1.16	0	14,15,17	1.21	2 (14%)
4	ASN	D	1003	-	7,8,8	0.82	0	6,10,10	0.46	0
5	EDO	D	1008	-	3,3,3	0.55	0	2,2,2	0.33	0
3	CIT	D	1002	1	11,11,12	1.08	0	14,15,17	1.51	4 (28%)
5	EDO	B	1008	-	3,3,3	0.24	0	2,2,2	0.39	0
5	EDO	B	1007	-	3,3,3	0.33	0	2,2,2	0.37	0
5	EDO	D	1005	-	3,3,3	0.55	0	2,2,2	0.72	0
4	ASN	C	1003	-	7,8,8	0.58	0	6,10,10	0.71	0
5	EDO	C	1005	-	3,3,3	0.34	0	2,2,2	0.46	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	A	1007	-	-	1/1/1/1	-
3	CIT	B	1005	1	-	0/15/15/16	-
4	ASN	B	1006	-	-	1/8/8/8	-
5	EDO	A	1008	-	-	0/1/1/1	-
5	EDO	D	1006	-	-	0/1/1/1	-
5	EDO	D	1007	-	-	1/1/1/1	-
4	ASN	A	1006	-	-	1/8/8/8	-
3	CIT	A	1005	1	-	3/15/15/16	-
5	EDO	D	1004	-	-	0/1/1/1	-
5	EDO	C	1004	-	-	0/1/1/1	-
5	EDO	A	1009	-	-	0/1/1/1	-
5	EDO	B	1009	-	-	0/1/1/1	-
3	CIT	C	1002	1	-	1/15/15/16	-
4	ASN	D	1003	-	-	2/8/8/8	-
5	EDO	D	1008	-	-	0/1/1/1	-
3	CIT	D	1002	1	-	2/15/15/16	-
5	EDO	B	1008	-	-	1/1/1/1	-
5	EDO	B	1007	-	-	0/1/1/1	-
5	EDO	D	1005	-	-	1/1/1/1	-
4	ASN	C	1003	-	-	2/8/8/8	-
5	EDO	C	1005	-	-	0/1/1/1	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1005	CIT	C3-C6	2.67	1.56	1.53
3	B	1005	CIT	O4-C5	-2.50	1.29	1.42
4	B	1006	ASN	O-C	2.48	1.29	1.22
3	A	1005	CIT	C3-C6	2.05	1.55	1.53
3	B	1005	CIT	O5-C6	2.00	1.28	1.22

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1002	CIT	O5-C6-C3	-3.21	115.88	122.09
3	C	1002	CIT	O5-C6-C3	-3.04	116.20	122.09
3	D	1002	CIT	O6-C6-C3	2.57	118.06	113.14
3	A	1005	CIT	O5-C6-C3	-2.49	117.27	122.09
3	B	1005	CIT	O4-C5-C4	2.31	117.82	111.33
3	D	1002	CIT	O1-C1-C2	-2.24	116.62	122.95
3	C	1002	CIT	O6-C6-C3	2.23	117.42	113.14
3	A	1005	CIT	O6-C6-C3	2.13	117.23	113.14
3	A	1005	CIT	O1-C1-C2	-2.05	117.15	122.95
3	D	1002	CIT	O3-C5-C4	-2.04	120.34	126.38
3	B	1005	CIT	O5-C6-C3	-2.04	118.14	122.09

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	1008	EDO	O1-C1-C2-O2
4	D	1003	ASN	OXT-C-CA-CB
4	D	1003	ASN	O-C-CA-CB
4	C	1003	ASN	OXT-C-CA-CB
4	C	1003	ASN	O-C-CA-CB
3	A	1005	CIT	C3-C4-C5-O3
3	C	1002	CIT	C3-C4-C5-O3
3	D	1002	CIT	C3-C4-C5-O3
4	B	1006	ASN	OXT-C-CA-CB
5	D	1007	EDO	O1-C1-C2-O2
5	D	1005	EDO	O1-C1-C2-O2
3	A	1005	CIT	O1-C1-C2-C3
3	A	1005	CIT	O2-C1-C2-C3
4	A	1006	ASN	OXT-C-CA-CB
5	A	1007	EDO	O1-C1-C2-O2
3	D	1002	CIT	O1-C1-C2-C3

There are no ring outliers.

8 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1007	EDO	2	0
3	B	1005	CIT	1	0
5	A	1008	EDO	1	0
3	A	1005	CIT	1	0
5	D	1004	EDO	1	0
5	C	1004	EDO	1	0
3	D	1002	CIT	3	0
5	B	1007	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 [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	305/358 (85%)	0.69	37 (12%) 8 9	10, 19, 46, 84	0
1	B	325/358 (90%)	0.76	42 (12%) 7 7	10, 18, 46, 74	0
1	C	311/358 (86%)	0.83	46 (14%) 5 6	7, 20, 53, 79	1 (0%)
1	D	324/358 (90%)	0.72	39 (12%) 9 9	10, 18, 44, 74	1 (0%)
All	All	1265/1432 (88%)	0.75	164 (12%) 7 7	7, 19, 48, 84	2 (0%)

All (164) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	233	PRO	10.8
1	D	233	PRO	10.6
1	C	39	PRO	8.1
1	B	15	ILE	6.6
1	B	232	GLN	6.2
1	C	37	LEU	6.2
1	D	232	GLN	5.9
1	A	14	THR	5.6
1	A	232	GLN	5.4
1	B	231	ARG	5.3
1	A	233	PRO	5.3
1	C	233	PRO	5.2
1	B	234	VAL	5.2
1	C	288	ALA	5.1
1	C	229	PHE	4.7
1	A	229	PHE	4.5
1	D	230	LEU	4.5
1	B	230	LEU	4.4
1	C	14	THR	4.4
1	B	225	VAL	4.3
1	C	232	GLN	4.3

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Mol	Chain	Res	Type	RSRZ
1	B	288	ALA	4.2
1	C	36	ALA	4.2
1	C	225	VAL	4.0
1	C	35	LEU	3.9
1	B	1	MET	3.9
1	A	225	VAL	3.9
1	C	230	LEU	3.8
1	B	194	PRO	3.8
1	D	288	ALA	3.8
1	D	225	VAL	3.8
1	D	14	THR	3.7
1	D	122	ALA	3.7
1	C	42	HIS	3.7
1	A	316	GLU	3.7
1	D	0	HIS	3.7
1	B	122	ALA	3.7
1	C	231	ARG	3.7
1	D	234	VAL	3.6
1	B	-4	PRO	3.6
1	B	17	MET	3.6
1	D	2	GLN	3.5
1	A	252	ALA	3.5
1	B	31	LEU	3.5
1	B	223	ALA	3.5
1	C	121	LEU	3.5
1	D	194	PRO	3.5
1	D	-1	SER	3.5
1	D	1	MET	3.5
1	C	43	ARG	3.5
1	C	15	ILE	3.4
1	B	121	LEU	3.4
1	B	-2	GLY	3.4
1	A	42	HIS	3.3
1	A	15	ILE	3.2
1	A	121	LEU	3.2
1	D	-3	ARG	3.2
1	B	14	THR	3.1
1	C	226	VAL	3.1
1	C	38	MET	3.1
1	A	43	ARG	3.1
1	D	231	ARG	3.1
1	C	16	GLY	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	253	PHE	3.0
1	B	16	GLY	3.0
1	C	129	GLN	3.0
1	D	17	MET	3.0
1	D	317	LEU	3.0
1	B	-1	SER	3.0
1	C	234	VAL	3.0
1	D	-5	VAL	3.0
1	B	253	PHE	3.0
1	A	16	GLY	3.0
1	A	226	VAL	2.9
1	D	337	ASP	2.9
1	D	229	PHE	2.9
1	D	30	HIS	2.9
1	A	120	PRO	2.8
1	A	249	GLN	2.8
1	B	226	VAL	2.8
1	C	44	PRO	2.8
1	C	289	LEU	2.8
1	C	250	ASN	2.7
1	A	230	LEU	2.7
1	B	34	GLN	2.7
1	A	288	ALA	2.7
1	D	319	THR	2.7
1	D	-4	PRO	2.7
1	B	2	GLN	2.7
1	B	189	ARG	2.7
1	A	248	PRO	2.7
1	A	318	ASP	2.6
1	B	120	PRO	2.6
1	A	289	LEU	2.6
1	B	0	HIS	2.6
1	B	30	HIS	2.6
1	A	44	PRO	2.6
1	B	-3	ARG	2.6
1	C	320	GLU	2.6
1	D	-2	GLY	2.6
1	D	31	LEU	2.5
1	C	191	ASN	2.5
1	C	249	GLN	2.5
1	A	194	PRO	2.5
1	D	253	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	15	ILE	2.5
1	A	259	GLU	2.5
1	B	337	ASP	2.5
1	D	223	ALA	2.5
1	C	324	LYS	2.5
1	A	263	ARG	2.4
1	B	28	SER	2.4
1	C	34	GLN	2.4
1	C	17	MET	2.4
1	A	40	GLU	2.4
1	C	253	PHE	2.3
1	D	192	THR	2.3
1	D	255	GLN	2.3
1	B	251	LYS	2.3
1	A	224	ASP	2.3
1	C	227	ARG	2.3
1	C	45	GLU	2.3
1	A	122	ALA	2.3
1	C	292	ALA	2.3
1	D	191	ASN	2.3
1	D	121	LEU	2.3
1	D	336	PRO	2.2
1	B	63	MET	2.2
1	B	227	ARG	2.2
1	C	40	GLU	2.2
1	C	122	ALA	2.2
1	D	203	ILE	2.2
1	B	191	ASN	2.2
1	B	250	ASN	2.2
1	C	228	ASN	2.2
1	D	226	VAL	2.2
1	C	316	GLU	2.2
1	A	337	ASP	2.2
1	B	224	ASP	2.2
1	B	58	MET	2.2
1	D	249	GLN	2.2
1	D	248	PRO	2.2
1	B	229	PHE	2.1
1	D	224	ASP	2.1
1	A	193	PRO	2.1
1	A	231	ARG	2.1
1	B	154	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	237	LEU	2.1
1	C	248	PRO	2.1
1	A	17	MET	2.1
1	D	197	HIS	2.1
1	C	260	ALA	2.1
1	A	227	ARG	2.1
1	C	124	LEU	2.1
1	A	292	ALA	2.1
1	B	322	ILE	2.1
1	C	12	GLY	2.1
1	A	255	GLN	2.1
1	C	193	PRO	2.1
1	B	237	LEU	2.1
1	C	120	PRO	2.0
1	A	188	ARG	2.0
1	C	41	PHE	2.0
1	A	319	THR	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
5	EDO	D	1007	4/4	0.82	0.17	31,37,37,37	0
5	EDO	D	1008	4/4	0.82	0.12	15,15,16,17	0
5	EDO	B	1008	4/4	0.83	0.14	28,37,37,42	0
5	EDO	D	1006	4/4	0.84	0.15	31,33,35,37	0
5	EDO	A	1007	4/4	0.84	0.15	37,37,37,40	0
3	CIT	B	1005	12/13	0.84	0.11	22,27,32,35	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	EDO	A	1009	4/4	0.88	0.10	17,18,18,19	0
5	EDO	D	1005	4/4	0.89	0.11	18,20,20,22	0
5	EDO	B	1009	4/4	0.89	0.09	13,13,13,14	0
3	CIT	C	1002	12/13	0.90	0.09	17,22,28,33	0
5	EDO	C	1005	4/4	0.90	0.09	16,17,17,20	0
3	CIT	D	1002	12/13	0.91	0.08	18,22,27,30	0
3	CIT	A	1005	12/13	0.92	0.09	19,23,28,31	0
4	ASN	A	1006	9/9	0.93	0.08	9,10,11,11	0
5	EDO	B	1007	4/4	0.93	0.08	18,19,20,22	0
5	EDO	A	1008	4/4	0.93	0.07	14,15,16,17	0
4	ASN	D	1003	9/9	0.95	0.07	10,12,14,15	0
2	CL	C	1001	1/1	0.95	0.06	19,19,19,19	0
2	CL	A	1003	1/1	0.95	0.08	21,21,21,21	0
4	ASN	C	1003	9/9	0.95	0.07	11,11,13,14	0
4	ASN	B	1006	9/9	0.96	0.06	10,12,14,14	0
5	EDO	D	1004	4/4	0.96	0.09	17,17,19,19	0
5	EDO	C	1004	4/4	0.96	0.06	14,15,16,16	0
2	CL	B	1003	1/1	0.97	0.07	19,19,19,19	0
2	CL	A	1001	1/1	0.98	0.03	13,13,13,13	0
2	CL	A	1004	1/1	0.98	0.03	25,25,25,25	0
2	CL	D	1001	1/1	0.98	0.05	19,19,19,19	0
2	CL	B	1002	1/1	0.98	0.04	13,13,13,13	0
2	CL	B	1004	1/1	0.99	0.03	21,21,21,21	0
2	CL	A	1002	1/1	0.99	0.03	14,14,14,14	0
2	CL	B	1001	1/1	0.99	0.04	13,13,13,13	0

6.5 Other polymers

There are no such residues in this entry.