



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 8E0R / pdb_00008e0r
Title : Crystal structure of mouse APCDD1 in P21 space group
Authors : Hsieh, F.L.; Chang, T.H.; Gabelli, S.B.; Nathans, J.
Deposited on : 2022-08-09
Resolution : 1.95 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

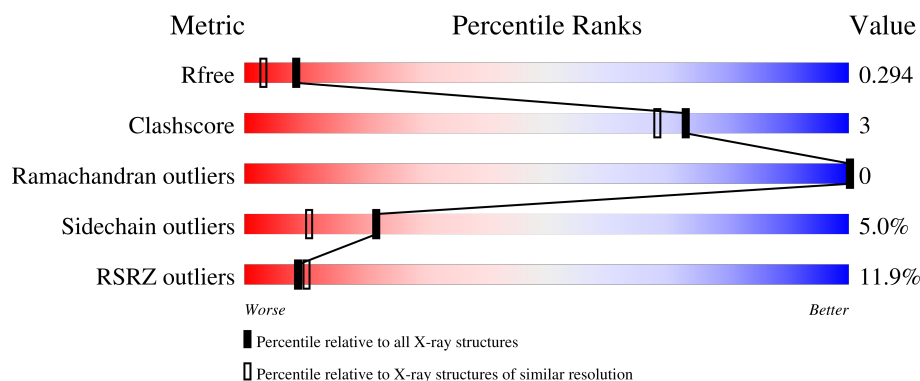
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	468	12% 80% 8% 11%
1	B	468	9% 74% 12% 13%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7140 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 Protein APCDD1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	418	Total	C	N	O	S	0	3	0
			3389	2140	618	609	22			
1	B	405	Total	C	N	O	S	0	2	0
			3307	2096	604	587	20			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	GLU	-	expression tag	UNP Q3U128
A	25	THR	-	expression tag	UNP Q3U128
A	26	GLY	-	expression tag	UNP Q3U128
A	483	THR	-	expression tag	UNP Q3U128
A	484	HIS	-	expression tag	UNP Q3U128
A	485	HIS	-	expression tag	UNP Q3U128
A	486	HIS	-	expression tag	UNP Q3U128
A	487	HIS	-	expression tag	UNP Q3U128
A	488	HIS	-	expression tag	UNP Q3U128
A	489	HIS	-	expression tag	UNP Q3U128
A	490	HIS	-	expression tag	UNP Q3U128
A	491	HIS	-	expression tag	UNP Q3U128
B	24	GLU	-	expression tag	UNP Q3U128
B	25	THR	-	expression tag	UNP Q3U128
B	26	GLY	-	expression tag	UNP Q3U128
B	483	THR	-	expression tag	UNP Q3U128
B	484	HIS	-	expression tag	UNP Q3U128
B	485	HIS	-	expression tag	UNP Q3U128
B	486	HIS	-	expression tag	UNP Q3U128
B	487	HIS	-	expression tag	UNP Q3U128
B	488	HIS	-	expression tag	UNP Q3U128
B	489	HIS	-	expression tag	UNP Q3U128
B	490	HIS	-	expression tag	UNP Q3U128
B	491	HIS	-	expression tag	UNP Q3U128

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	194	Total	O	0	0
			194	194		
3	B	180	Total	O	0	0
			180	180		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.86Å 144.13Å 64.38Å 90.00° 99.52° 90.00°	Depositor
Resolution (Å)	34.00 – 1.95 34.00 – 1.95	Depositor EDS
% Data completeness (in resolution range)	76.6 (34.00-1.95) 76.6 (34.00-1.95)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.11 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.209 , 0.247 0.264 , 0.294	Depositor DCC
R_{free} test set	2407 reflections (3.68%)	wwPDB-VP
Wilson B-factor (Å ²)	32.0	Xtriage
Anisotropy	0.021	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 30.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7140	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.50% 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: NAG

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	0.97	0/3498	1.24	2/4748 (0.0%)
1	B	1.08	3/3416 (0.1%)	1.29	10/4637 (0.2%)
All	All	1.03	3/6914 (0.0%)	1.27	12/9385 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	368[A]	HIS	C-O	15.76	1.44	1.23
1	B	368[B]	HIS	C-O	15.76	1.44	1.23
1	B	368[C]	HIS	C-O	15.76	1.44	1.23

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	368[A]	HIS	CA-C-O	11.05	134.04	121.11
1	B	368[B]	HIS	CA-C-O	11.05	134.04	121.11
1	B	368[C]	HIS	CA-C-O	11.05	134.04	121.11
1	B	368[A]	HIS	O-C-N	-8.40	112.31	123.23
1	B	368[B]	HIS	O-C-N	-8.40	112.31	123.23
1	B	368[C]	HIS	O-C-N	-8.40	112.31	123.23
1	A	244	PHE	CB-CA-C	-6.73	98.79	110.17
1	A	244	PHE	CA-CB-CG	5.65	119.45	113.80
1	B	368[A]	HIS	N-CA-C	5.32	118.07	109.40
1	B	368[B]	HIS	N-CA-C	5.32	118.07	109.40
1	B	368[C]	HIS	N-CA-C	5.32	118.07	109.40
1	B	333	VAL	N-CA-C	-5.03	107.39	112.17

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	3389	0	3244	19	0
1	B	3307	0	3169	25	3
2	A	42	0	39	0	0
2	B	28	0	26	0	0
3	A	194	0	0	2	0
3	B	180	0	0	2	0
All	All	7140	0	6478	42	3

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 3.

All (42) 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:403:ASP:OD1	1:A:405:THR:HG23	1.91	0.70
1:A:298:GLN:OE1	3:A:601:HOH:O	2.09	0.70
1:A:368[A]:HIS:CD2	1:A:405:THR:HG21	2.30	0.66
1:B:339:PHE:HB3	1:B:378:THR:HG21	1.83	0.61
1:B:224:ASP:OD2	3:B:601:HOH:O	2.15	0.60
1:B:369:MET:H	1:B:405:THR:HG22	1.68	0.59
1:A:339:PHE:HB3	1:A:378:THR:HG21	1.85	0.58
1:B:219:PRO:HD2	1:B:225:HIS:O	2.06	0.55
1:B:157:GLU:CD	1:B:157:GLU:H	2.14	0.55
1:A:101:ASN:HB2	1:A:127:LYS:HE2	1.90	0.51
1:A:225[B]:HIS:CE1	1:A:262:HIS:H	2.29	0.51
1:B:129:ARG:NH2	3:B:604:HOH:O	2.32	0.50
1:B:157:GLU:CD	1:B:157:GLU:N	2.71	0.49
1:A:219:PRO:HB2	1:A:222:SER:HB3	1.94	0.49
1:B:451:ARG:NH1	1:B:454:LYS:HD2	2.27	0.48
1:B:56:LEU:HD11	1:B:312:HIS:HB2	1.94	0.48
1:A:263:ALA:HA	1:B:431:THR:O	2.13	0.48
1:B:403:ASP:OD1	1:B:405:THR:HG23	2.14	0.48
1:B:322:TRP:HZ2	1:B:366:VAL:HG22	1.80	0.47
1:B:425:PHE:HD2	1:B:437:LEU:HD11	1.81	0.46
1:B:76:HIS:HE2	1:B:94:SER:HG	1.62	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:THR:HB	1:A:204:PHE:HB3	1.98	0.46
1:A:427[B]:MET:HE3	1:A:427[B]:MET:HB2	1.76	0.46
1:A:403:ASP:CG	1:A:405:THR:HG23	2.40	0.46
1:B:369:MET:HE3	1:B:369:MET:HB2	1.85	0.44
1:A:114:ARG:HD3	1:A:284:LYS:O	2.17	0.44
1:A:285:ALA:HB2	1:B:225:HIS:NE2	2.32	0.44
1:A:405:THR:HA	1:A:417:LEU:HD12	2.01	0.43
1:B:162:GLN:O	1:B:165:ARG:HG2	2.19	0.43
1:A:225[B]:HIS:CE1	1:A:261:ASN:HA	2.55	0.42
1:B:164:SER:HB3	1:B:181:TRP:HB2	2.01	0.42
1:B:199:THR:HB	1:B:204:PHE:HB3	2.02	0.42
1:B:427:MET:HE3	1:B:427:MET:HB2	1.88	0.41
1:A:219:PRO:HA	3:A:672:HOH:O	2.20	0.41
1:B:178:GLY:O	1:B:180:PRO:HD3	2.21	0.41
1:A:176:ALA:HB1	1:A:177:PRO:CD	2.50	0.41
1:A:368[A]:HIS:CE1	1:A:370:LYS:HG3	2.56	0.41
1:B:112:SER:OG	1:B:114:ARG:HG2	2.21	0.41
1:B:226:LEU:O	1:B:259:ASN:HB3	2.21	0.41
1:A:159:VAL:O	1:A:163:LEU:HB2	2.20	0.40
1:B:291:LEU:O	1:B:292:HIS:C	2.65	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:191:GLN:NE2	1:B:406:HIS:CD2[1_556]	1.95	0.25
1:B:224:ASP:OD2	1:B:354:SER:OG[1_455]	2.11	0.09
1:B:191:GLN:NE2	1:B:406:HIS:NE2[1_556]	2.13	0.07

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	415/468 (89%)	391 (94%)	24 (6%)	0	100	100
1	B	400/468 (86%)	383 (96%)	17 (4%)	0	100	100
All	All	815/936 (87%)	774 (95%)	41 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	369/411 (90%)	351 (95%)	18 (5%)	22	11
1	B	360/411 (88%)	342 (95%)	18 (5%)	22	11
All	All	729/822 (89%)	693 (95%)	36 (5%)	22	11

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

Mol	Chain	Res	Type
1	A	86	SER
1	A	129	ARG
1	A	163	LEU
1	A	184	ASP
1	A	194	SER
1	A	197	GLU
1	A	200	LYS
1	A	243	VAL
1	A	244	PHE
1	A	306	VAL
1	A	335	LYS
1	A	405	THR
1	A	410	CYS
1	A	411	VAL
1	A	413	LEU
1	A	424	ILE
1	A	451	ARG
1	A	469	SER

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Mol	Chain	Res	Type
1	B	60	HIS
1	B	64	ARG
1	B	65	ILE
1	B	86	SER
1	B	104	LYS
1	B	114	ARG
1	B	157	GLU
1	B	163	LEU
1	B	175	LEU
1	B	184	ASP
1	B	217	GLN
1	B	243	VAL
1	B	299	ARG
1	B	355	LYS
1	B	405	THR
1	B	411	VAL
1	B	417	LEU
1	B	424	ILE

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	58	HIS
1	A	107	GLN
1	A	191	GLN
1	A	261	ASN
1	A	397	GLN
1	B	107	GLN
1	B	151	GLN
1	B	210	GLN
1	B	217	GLN
1	B	241	GLN
1	B	276	HIS
1	B	328	HIS

5.3.3 RNA ⓘ

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](#)

5 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	NAG	A	503	1	14,14,15	0.46	0	17,19,21	0.98	2 (11%)
2	NAG	B	502	1	14,14,15	0.55	0	17,19,21	1.34	4 (23%)
2	NAG	A	502	1	14,14,15	0.35	0	17,19,21	0.88	1 (5%)
2	NAG	B	501	1	14,14,15	0.33	0	17,19,21	1.30	1 (5%)
2	NAG	A	501	1	14,14,15	0.28	0	17,19,21	1.34	1 (5%)

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	NAG	A	503	1	-	0/6/23/26	0/1/1/1
2	NAG	B	502	1	-	0/6/23/26	0/1/1/1
2	NAG	A	502	1	-	0/6/23/26	0/1/1/1
2	NAG	B	501	1	-	1/6/23/26	0/1/1/1
2	NAG	A	501	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NAG	C1-O5-C5	4.84	118.68	112.19
2	B	501	NAG	C1-O5-C5	3.78	117.25	112.19
2	B	502	NAG	C4-C3-C2	-2.59	107.22	111.02
2	B	502	NAG	C3-C4-C5	-2.48	105.73	110.23
2	A	503	NAG	O4-C4-C3	-2.30	104.97	110.38
2	A	503	NAG	C1-C2-N2	2.26	114.00	110.43
2	B	502	NAG	O5-C5-C4	-2.15	105.59	110.83
2	A	502	NAG	C2-N2-C7	2.12	125.74	122.90
2	B	502	NAG	O5-C5-C6	2.03	111.62	107.66

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	NAG	C4-C5-C6-O6
2	A	501	NAG	O5-C5-C6-O6
2	B	501	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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	418/468 (89%)	0.84	54 (12%) 7 9	2, 8, 19, 25	3 (0%)
1	B	405/468 (86%)	0.76	44 (10%) 10 12	2, 9, 25, 42	2 (0%)
All	All	823/936 (87%)	0.80	98 (11%) 9 10	2, 9, 22, 42	5 (0%)

All (98) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	413	LEU	6.9
1	A	135	TRP	6.1
1	B	135	TRP	5.7
1	B	406	HIS	5.5
1	B	178	GLY	5.5
1	B	177	PRO	5.4
1	A	409	GLY	5.4
1	B	468	SER	5.1
1	A	220	HIS	4.7
1	A	136	ILE	4.6
1	A	221	HIS	4.4
1	B	60	HIS	4.3
1	B	50	SER	4.2
1	B	170	THR	4.0
1	B	160	ALA	4.0
1	A	271	PHE	4.0
1	A	178	GLY	3.9
1	B	53	HIS	3.9
1	B	467	ALA	3.8
1	A	174	PHE	3.8
1	A	388	ASN	3.8
1	A	61	ASN	3.8
1	B	171	CYS	3.8
1	B	65	ILE	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	246	ARG	3.7
1	A	414	GLY	3.7
1	A	472	ARG	3.7
1	A	68	GLN	3.7
1	A	411	VAL	3.6
1	B	64	ARG	3.5
1	A	50	SER	3.5
1	B	196	HIS	3.4
1	B	174	PHE	3.4
1	A	393	GLU	3.4
1	A	243	VAL	3.4
1	A	392	ALA	3.4
1	B	385	PHE	3.4
1	A	194	SER	3.3
1	B	176	ALA	3.2
1	B	198	CYS	3.2
1	B	394	GLY	3.2
1	A	410	CYS	3.2
1	A	240	THR	3.1
1	A	191	GLN	3.1
1	B	175	LEU	3.1
1	A	177	PRO	3.1
1	A	471	PRO	3.1
1	B	351	VAL	3.1
1	A	176	ALA	3.0
1	A	60	HIS	2.9
1	A	192	GLU	2.9
1	B	414	GLY	2.9
1	A	203	ASN	2.8
1	B	58	HIS	2.8
1	A	405	THR	2.8
1	B	225	HIS	2.8
1	B	197	GLU	2.7
1	B	62	GLY	2.6
1	B	182	VAL	2.6
1	B	165	ARG	2.5
1	B	172	PRO	2.5
1	B	192	GLU	2.5
1	A	244	PHE	2.4
1	B	179	GLY	2.4
1	A	386	SER	2.4
1	B	240	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	62	GLY	2.4
1	A	351	VAL	2.4
1	B	136	ILE	2.3
1	B	220	HIS	2.3
1	A	219	PRO	2.3
1	A	377	ALA	2.3
1	B	61	ASN	2.2
1	A	394	GLY	2.2
1	A	201	ALA	2.2
1	A	376	ALA	2.2
1	A	169	ARG	2.2
1	B	169	ARG	2.2
1	B	199	THR	2.2
1	B	173	GLY	2.2
1	A	283	PRO	2.2
1	B	349	ARG	2.2
1	A	57	LYS	2.1
1	A	172	PRO	2.1
1	A	470	SER	2.1
1	A	129	ARG	2.1
1	A	173	GLY	2.1
1	B	129	ARG	2.1
1	A	217	GLN	2.1
1	A	412	ALA	2.1
1	B	162	GLN	2.1
1	B	224	ASP	2.1
1	A	166	LEU	2.1
1	A	130	LEU	2.0
1	A	314	ILE	2.0
1	A	132	GLN	2.0
1	B	111	GLY	2.0
1	A	70	PRO	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	A	501	14/15	0.64	0.15	18,20,22,23	0
2	NAG	B	501	14/15	0.79	0.14	23,24,27,28	0
2	NAG	A	502	14/15	0.83	0.08	21,22,24,25	0
2	NAG	A	503	14/15	0.86	0.09	10,10,12,13	0
2	NAG	B	502	14/15	0.87	0.09	11,11,12,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.