



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:55 AM UTC

PDB ID : 4X0U / pdb_00004x0u
Title : Structure ALDH7A1 inactivated by 4-diethylaminobenzaldehyde
Authors : Luo, M.; Tanner, J.J.
Deposited on : 2014-11-23
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

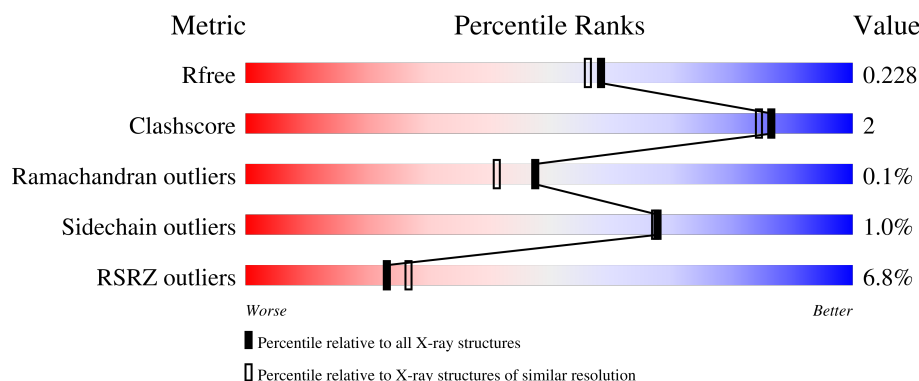
1 Overall quality at a glance

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	513	4% 88% 5% 7%
1	B	513	5% 89% 5% 6%
1	C	513	10% 88% 7% 5%
1	D	513	7% 88% 7% 5%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15034 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.

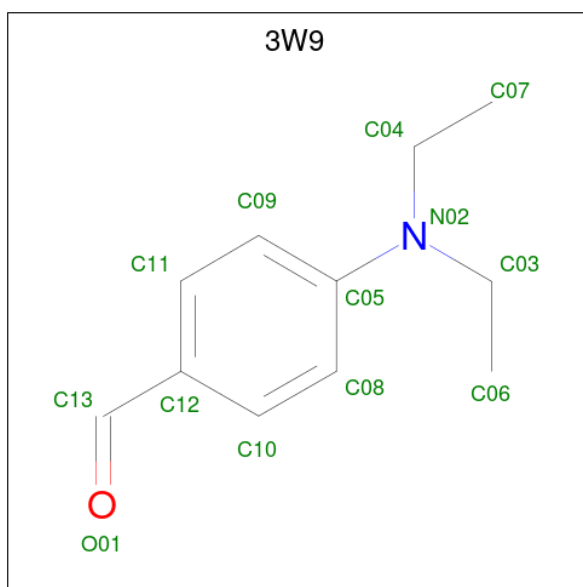
- Molecule 1 is a protein called Alpha-aminoadipic semialdehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	477	Total	C	N	O	S	0	0	0
			3588	2297	618	656	17			
1	B	484	Total	C	N	O	S	0	0	0
			3629	2316	626	671	16			
1	C	485	Total	C	N	O	S	0	0	0
			3613	2307	622	667	17			
1	D	487	Total	C	N	O	S	0	0	0
			3635	2318	632	668	17			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P49419
A	0	HIS	-	expression tag	UNP P49419
B	-1	GLY	-	expression tag	UNP P49419
B	0	HIS	-	expression tag	UNP P49419
C	-1	GLY	-	expression tag	UNP P49419
C	0	HIS	-	expression tag	UNP P49419
D	-1	GLY	-	expression tag	UNP P49419
D	0	HIS	-	expression tag	UNP P49419

- Molecule 2 is 4-(diethylamino)benzaldehyde (CCD ID: 3W9) (formula: C₁₁H₁₅NO).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			13	11	1	1		
2	B	1	Total	C	N	O	0	0
			13	11	1	1		
2	C	1	Total	C	N	O	0	0
			13	11	1	1		
2	D	1	Total	C	N	O	0	0
			13	11	1	1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total	Mg	0	0
			1	1		

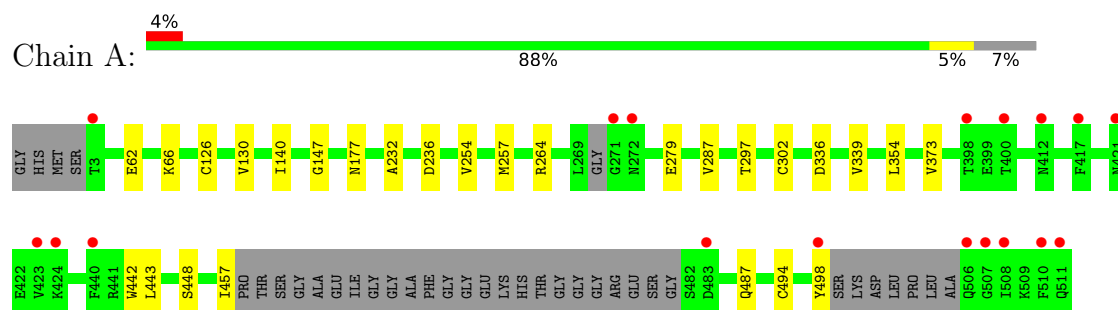
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	143	Total	O	0	0
			143	143		
4	B	125	Total	O	0	0
			125	125		
4	C	101	Total	O	0	0
			101	101		
4	D	147	Total	O	0	0
			147	147		

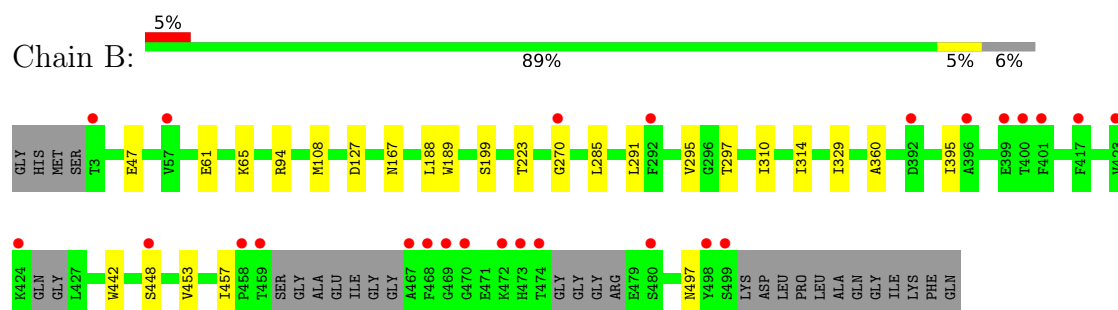
3 Residue-property plots [i](#)

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.

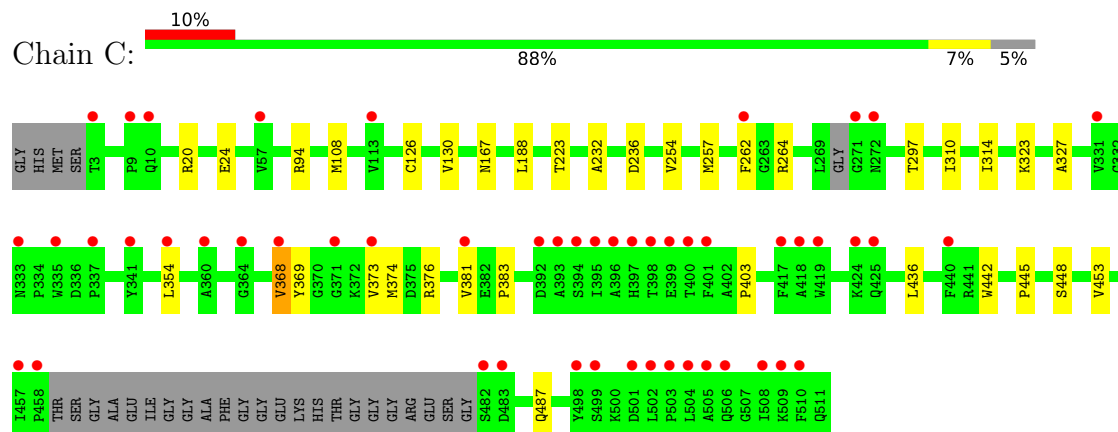
- Molecule 1: Alpha-aminoadipic semialdehyde dehydrogenase



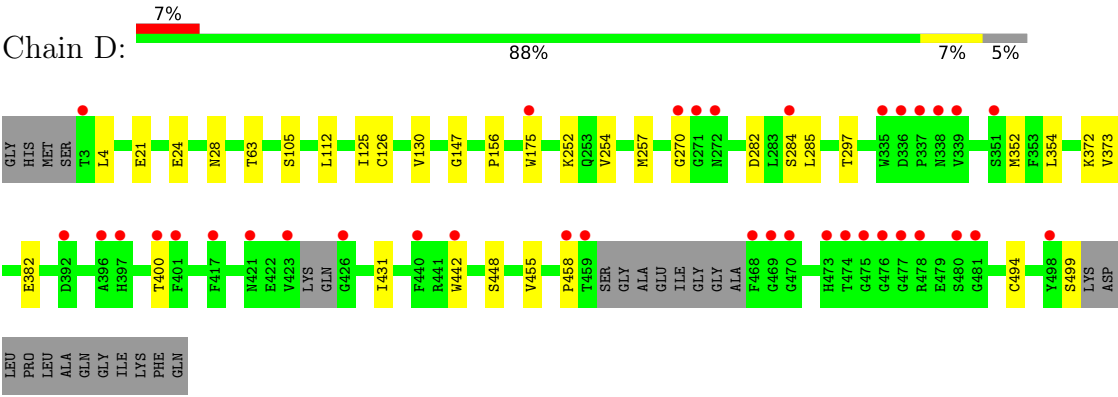
- Molecule 1: Alpha-aminoadipic semialdehyde dehydrogenase



- Molecule 1: Alpha-aminoadipic semialdehyde dehydrogenase



- Molecule 1: Alpha-aminoadipic semialdehyde dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	F 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	161.54Å 219.41Å 233.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	51.49 – 1.95 51.49 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.7 (51.49-1.95) 99.7 (51.49-1.95)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 1.95Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.188 , 0.227 0.190 , 0.228	Depositor DCC
R_{free} test set	7464 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	25.6	Xtriage
Anisotropy	0.393	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	15034	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.36% 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: 3W9, MG

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	0.46	0/3660	0.75	3/4974 (0.1%)
1	B	0.45	0/3703	0.76	0/5036
1	C	0.44	0/3687	0.76	0/5022
1	D	0.47	0/3708	0.76	0/5038
All	All	0.46	0/14758	0.76	3/20070 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	287	VAL	CA-C-N	-5.29	113.46	119.28
1	A	287	VAL	C-N-CA	-5.29	113.46	119.28
1	A	302	CYS	N-CA-C	-5.28	106.35	112.89

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	3588	0	3549	14	0
1	B	3629	0	3566	12	0
1	C	3613	0	3525	20	0
1	D	3635	0	3595	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	13	0	14	0	0
2	B	13	0	14	0	0
2	C	13	0	14	0	0
2	D	13	0	14	0	0
3	D	1	0	0	0	0
4	A	143	0	0	2	0
4	B	125	0	0	1	0
4	C	101	0	0	0	0
4	D	147	0	0	0	0
All	All	15034	0	14291	59	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 (59) 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:C:323:LYS:HD3	1:C:368:VAL:HG12	1.79	0.65
1:C:354:LEU:HD21	1:C:373:VAL:HG23	1.79	0.63
1:D:285:LEU:HD22	1:D:458:PRO:HB3	1.81	0.61
1:D:442:TRP:CZ2	1:D:448:SER:HB2	2.35	0.61
1:B:442:TRP:CZ2	1:B:448:SER:HB2	2.37	0.60
1:D:28:ASN:HB3	1:D:63:THR:HG23	1.85	0.57
1:D:354:LEU:HD21	1:D:373:VAL:HG23	1.86	0.57
1:A:442:TRP:CE2	1:A:448:SER:HB2	2.42	0.55
1:A:177:ASN:ND2	4:A:789:HOH:O	2.25	0.54
1:C:108:MET:HE3	1:C:167:ASN:HA	1.89	0.54
1:D:147:GLY:HA3	1:D:499:SER:HB3	1.89	0.53
1:A:147:GLY:O	1:A:498:TYR:N	2.41	0.53
1:A:66:LYS:NZ	4:A:701:HOH:O	2.42	0.53
1:C:20:ARG:NH2	1:C:24:GLU:OE1	2.33	0.53
1:A:442:TRP:CZ2	1:A:448:SER:HB2	2.43	0.52
1:C:442:TRP:CE2	1:C:448:SER:HB2	2.46	0.51
1:D:125:ILE:HG21	1:D:175:TRP:HA	1.94	0.50
1:D:442:TRP:CE2	1:D:448:SER:HB2	2.47	0.50
1:B:94:ARG:NH2	1:B:127:ASP:OD2	2.44	0.49
1:C:188:LEU:HD11	1:C:223:THR:HG23	1.93	0.49
1:A:126:CYS:O	1:A:130:VAL:HG23	2.13	0.48
1:D:254:VAL:HA	1:D:257:MET:HE3	1.95	0.48
1:A:336:ASP:HB2	1:A:339:VAL:HG21	1.96	0.47
1:B:291:LEU:HD11	1:B:329:ILE:HD11	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:ARG:NH1	1:A:487:GLN:O	2.47	0.47
1:C:383:PRO:HA	1:C:403:PRO:HB2	1.98	0.46
1:B:310:ILE:HG22	1:B:314:ILE:HG13	1.96	0.46
1:A:354:LEU:HD21	1:A:373:VAL:HG23	1.98	0.46
1:B:61:GLU:HG3	1:B:65:LYS:HE3	1.98	0.46
1:C:310:ILE:HG22	1:C:314:ILE:HG13	1.98	0.46
1:B:188:LEU:HD11	1:B:223:THR:HG23	1.99	0.44
1:C:264:ARG:NH2	1:C:487:GLN:O	2.50	0.44
1:C:126:CYS:O	1:C:130:VAL:HG23	2.18	0.44
1:A:494:CYS:HA	1:B:453:VAL:O	2.17	0.44
1:C:327:ALA:HA	1:C:369:TYR:OH	2.17	0.43
1:C:254:VAL:HA	1:C:257:MET:HE3	2.00	0.43
1:C:354:LEU:HD23	1:C:354:LEU:HA	1.89	0.42
1:B:108:MET:HE3	1:B:167:ASN:HA	2.00	0.42
1:D:352:MET:HE3	1:D:400:THR:HG22	2.01	0.42
1:D:126:CYS:O	1:D:130:VAL:HG23	2.20	0.42
1:B:360:ALA:HA	1:B:395:ILE:HG21	2.01	0.41
1:D:4:LEU:CD1	1:D:21:GLU:HG3	2.50	0.41
1:A:254:VAL:HA	1:A:257:MET:HE3	2.02	0.41
1:D:372:LYS:HE3	1:D:382:GLU:OE1	2.20	0.41
1:B:457:ILE:HB	4:B:817:HOH:O	2.20	0.41
1:A:279:GLU:H	1:A:279:GLU:CD	2.29	0.41
1:A:457:ILE:HD13	1:B:497:ASN:HB2	2.02	0.41
1:C:262:PHE:CZ	1:D:252:LYS:HB2	2.56	0.41
1:C:374:MET:O	1:C:376:ARG:N	2.54	0.41
1:D:431:ILE:HD11	1:D:455:VAL:HG22	2.03	0.41
1:C:453:VAL:O	1:D:494:CYS:HA	2.21	0.41
1:C:232:ALA:O	1:C:236:ASP:HB2	2.20	0.41
1:D:282:ASP:OD1	1:D:284:SER:HB3	2.21	0.40
1:C:436:LEU:HD12	1:C:436:LEU:HA	1.90	0.40
1:D:105:SER:OG	1:D:112:LEU:HA	2.22	0.40
1:B:189:TRP:CH2	1:B:199:SER:HA	2.57	0.40
1:C:445:PRO:HB3	1:D:156:PRO:HD2	2.03	0.40
1:C:442:TRP:CZ2	1:C:448:SER:HB2	2.56	0.40
1:A:232:ALA:O	1:A:236:ASP:HB2	2.22	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	469/513 (91%)	460 (98%)	9 (2%)	0	100	100
1	B	476/513 (93%)	465 (98%)	10 (2%)	1 (0%)	43	36
1	C	479/513 (93%)	471 (98%)	8 (2%)	0	100	100
1	D	481/513 (94%)	471 (98%)	9 (2%)	1 (0%)	43	36
All	All	1905/2052 (93%)	1867 (98%)	36 (2%)	2 (0%)	48	41

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	270	GLY
1	B	270	GLY

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	364/410 (89%)	360 (99%)	4 (1%)	65	64
1	B	369/410 (90%)	365 (99%)	4 (1%)	65	64
1	C	363/410 (88%)	359 (99%)	4 (1%)	65	64
1	D	369/410 (90%)	367 (100%)	2 (0%)	81	82
All	All	1465/1640 (89%)	1451 (99%)	14 (1%)	68	67

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	GLU
1	A	140	ILE
1	A	297	THR
1	A	443	LEU
1	B	47	GLU
1	B	285	LEU
1	B	295	VAL
1	B	297	THR
1	C	94	ARG
1	C	297	THR
1	C	368	VAL
1	C	381	VAL
1	D	24	GLU
1	D	297	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	120	GLN
1	A	213	ASN
1	A	328	GLN
1	A	338	ASN
1	A	454	ASN
1	B	88	GLN
1	B	213	ASN
1	B	454	ASN
1	C	46	ASN
1	C	148	HIS
1	C	155	ASN
1	C	177	ASN
1	C	213	ASN
1	C	338	ASN
1	D	46	ASN
1	D	177	ASN
1	D	412	ASN

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 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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	$\# Z > 2$
2	3W9	B	601	1	13,13,13	0.65	0	16,16,16	0.64	0
2	3W9	D	601	1	13,13,13	0.64	0	16,16,16	0.72	0
2	3W9	A	601	1	13,13,13	0.56	0	16,16,16	0.84	0
2	3W9	C	601	1	13,13,13	0.61	0	16,16,16	0.72	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	3W9	B	601	1	-	2/10/10/10	0/1/1/1
2	3W9	D	601	1	-	0/10/10/10	0/1/1/1
2	3W9	A	601	1	-	2/10/10/10	0/1/1/1
2	3W9	C	601	1	-	2/10/10/10	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	3W9	C11-C12-C13-O01
2	A	601	3W9	C10-C12-C13-O01
2	B	601	3W9	C11-C12-C13-O01
2	B	601	3W9	C10-C12-C13-O01
2	C	601	3W9	C11-C12-C13-O01
2	C	601	3W9	C10-C12-C13-O01

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	477/513 (92%)	0.01	18 (3%)	44	50	15, 28, 51, 67	0
1	B	484/513 (94%)	0.15	25 (5%)	33	38	17, 31, 54, 68	0
1	C	485/513 (94%)	0.43	51 (10%)	11	13	18, 34, 61, 72	0
1	D	487/513 (94%)	0.26	37 (7%)	20	22	15, 31, 58, 68	0
All	All	1933/2052 (94%)	0.21	131 (6%)	23	27	15, 31, 57, 72	0

All (131) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	335	TRP	5.2
1	C	272	ASN	5.1
1	B	392	ASP	4.5
1	D	476	GLY	4.5
1	C	400	THR	4.3
1	B	459	THR	4.2
1	C	504	LEU	4.2
1	B	468	PHE	4.1
1	B	499	SER	4.1
1	B	467	ALA	4.0
1	D	475	GLY	4.0
1	D	459	THR	3.9
1	D	477	GLY	3.9
1	A	272	ASN	3.8
1	D	481	GLY	3.8
1	D	498	TYR	3.7
1	B	423	VAL	3.7
1	C	499	SER	3.6
1	D	468	PHE	3.6
1	D	175	TRP	3.5
1	B	474	THR	3.5

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Mol	Chain	Res	Type	RSRZ
1	D	423	VAL	3.5
1	D	417	PHE	3.5
1	B	292	PHE	3.4
1	B	473	HIS	3.4
1	C	458	PRO	3.4
1	D	335	TRP	3.3
1	A	421	ASN	3.3
1	C	440	PHE	3.3
1	B	480	SER	3.3
1	C	341	TYR	3.3
1	A	510	PHE	3.3
1	D	458	PRO	3.2
1	C	417	PHE	3.2
1	C	425	GLN	3.2
1	C	394	SER	3.1
1	A	483	ASP	3.1
1	D	337	PRO	3.1
1	C	271	GLY	3.1
1	D	469	GLY	3.1
1	D	426	GLY	3.0
1	A	400	THR	2.9
1	A	3	THR	2.9
1	C	503	PRO	2.9
1	C	505	ALA	2.9
1	D	270	GLY	2.9
1	C	3	THR	2.8
1	C	506	GLN	2.8
1	C	510	PHE	2.8
1	D	478	ARG	2.8
1	C	262	PHE	2.8
1	D	351	SER	2.8
1	D	480	SER	2.8
1	B	424	LYS	2.8
1	C	381	VAL	2.8
1	C	401	PHE	2.8
1	B	470	GLY	2.8
1	C	502	LEU	2.8
1	A	498	TYR	2.7
1	B	469	GLY	2.7
1	A	417	PHE	2.7
1	C	501	ASP	2.7
1	A	511	GLN	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	423	VAL	2.7
1	A	507	GLY	2.7
1	B	498	TYR	2.7
1	D	271	GLY	2.6
1	C	395	ILE	2.6
1	B	270	GLY	2.6
1	B	399	GLU	2.6
1	A	398	THR	2.6
1	A	412	ASN	2.6
1	B	417	PHE	2.6
1	C	354	LEU	2.6
1	C	333	ASN	2.6
1	C	397	HIS	2.5
1	C	360	ALA	2.5
1	D	3	THR	2.5
1	A	271	GLY	2.4
1	D	272	ASN	2.4
1	D	473	HIS	2.4
1	B	472	LYS	2.4
1	C	393	ALA	2.4
1	D	400	THR	2.4
1	C	371	GLY	2.4
1	C	399	GLU	2.4
1	C	508	ILE	2.4
1	C	396	ALA	2.4
1	B	458	PRO	2.4
1	B	396	ALA	2.3
1	D	397	HIS	2.3
1	D	339	VAL	2.3
1	C	57	VAL	2.3
1	C	331	VAL	2.3
1	C	392	ASP	2.3
1	B	400	THR	2.3
1	D	440	PHE	2.3
1	C	337	PRO	2.3
1	A	508	ILE	2.3
1	B	448	SER	2.3
1	C	419	TRP	2.2
1	D	442	TRP	2.2
1	C	482	SER	2.2
1	C	457	ILE	2.2
1	C	113	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	368	VAL	2.2
1	D	336	ASP	2.2
1	D	396	ALA	2.2
1	B	401	PHE	2.2
1	D	338	ASN	2.2
1	D	421	ASN	2.2
1	B	3	THR	2.2
1	C	364	GLY	2.2
1	A	424	LYS	2.2
1	C	418	ALA	2.2
1	C	424	LYS	2.2
1	C	509	LYS	2.2
1	C	498	TYR	2.1
1	C	483	ASP	2.1
1	A	440	PHE	2.1
1	D	401	PHE	2.1
1	D	284	SER	2.1
1	C	398	THR	2.1
1	A	506	GLN	2.1
1	D	392	ASP	2.1
1	D	470	GLY	2.0
1	C	10	GLN	2.0
1	D	474	THR	2.0
1	C	9	PRO	2.0
1	B	57	VAL	2.0
1	C	373	VAL	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	3W9	B	601	13/13	0.81	0.16	31,44,52,56	0
2	3W9	C	601	13/13	0.81	0.15	41,43,49,54	0
2	3W9	A	601	13/13	0.84	0.14	36,42,46,47	0
2	3W9	D	601	13/13	0.86	0.13	30,38,43,47	0
3	MG	D	602	1/1	0.94	0.13	43,43,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.