



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 04:42 am GMT

PDB ID : 2C11
Title : CRYSTAL STRUCTURE OF THE 2-HYDRAZINOPYRIDINE OF SEMIC
ARBAZIDE-SENSITIVE AMINE OXIDASE
Authors : Jakobsson, E.; Kleywegt, G.J.
Deposited on : 2005-09-09
Resolution : 2.90 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

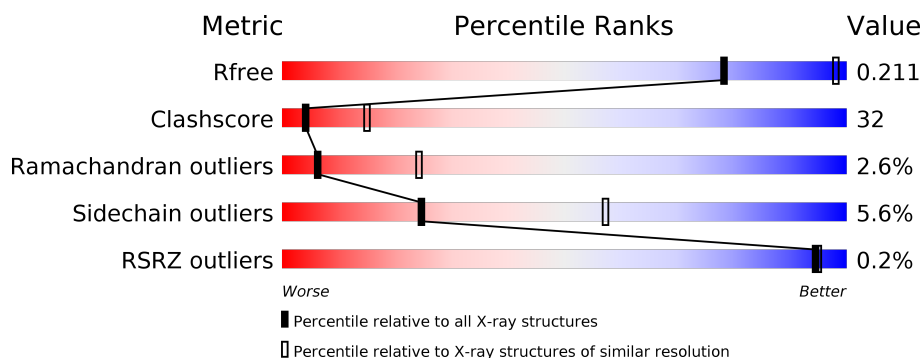
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	100719	1586 (2.90-2.90)
Clashscore	112137	1807 (2.90-2.90)
Ramachandran outliers	110173	1768 (2.90-2.90)
Sidechain outliers	110143	1770 (2.90-2.90)
RSRZ outliers	101464	1596 (2.90-2.90)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	735	
1	B	735	
1	C	735	
1	D	735	

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
1	PAQ	A	471	X	-	-	-
1	PAQ	B	471	X	-	-	-
1	PAQ	C	471	X	-	-	-
1	PAQ	D	471	X	-	-	-
6	CU	A	1744	-	-	-	X
6	CU	B	1748	-	-	-	X

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 21917 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 MEMBRANE COPPER AMINE OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	672	Total	C	N	O	S	0	0	1
			5340	3434	924	965	17			
1	B	672	Total	C	N	O	S	0	0	1
			5340	3434	924	965	17			
1	C	672	Total	C	N	O	S	0	0	1
			5340	3434	924	965	17			
1	D	672	Total	C	N	O	S	0	0	1
			5340	3434	924	965	17			

- Molecule 2 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	2	Total	C	N	O	0	0
			28	16	2	10		
2	C	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 3 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	2	Total	C	N	O	0	0
			24	14	1	9		

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	2	Total	Ca	0	0
			2	2		
5	A	2	Total	Ca	0	0
			2	2		
5	D	2	Total	Ca	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	2	Total	Ca	0	0
			2	2		

- Molecule 6 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	8	Total	Cu	0	0
			8	8		
6	A	7	Total	Cu	0	0
			7	7		
6	D	7	Total	Cu	0	0
			7	7		
6	C	8	Total	Cu	0	0
			8	8		

- Molecule 7 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 8 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	B	3	Total	C	N	O	0	0
			39	22	2	15		
8	D	3	Total	C	N	O	0	0
			39	22	2	15		

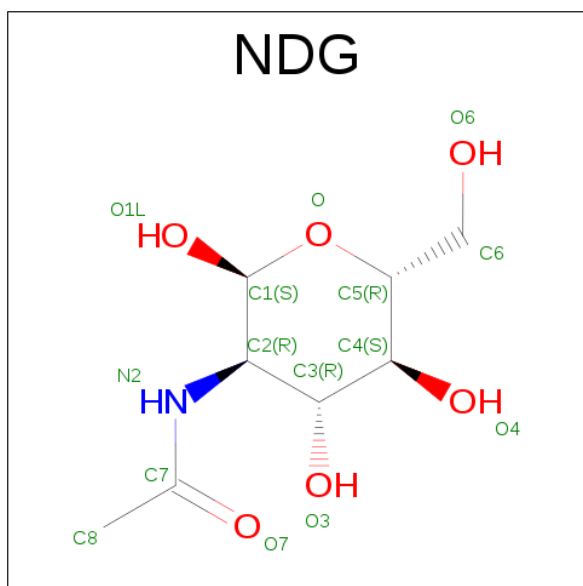
- Molecule 9 is a polymer of unknown type called SUGAR (5-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	B	5	Total	C	N	O	0	0
			60	34	2	24		
9	D	5	Total	C	N	O	0	0
			60	34	2	24		

- Molecule 10 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
10	C	3	Total	C	N	O	0	0
			38	22	2	14		

- Molecule 11 is SUGAR (2-(ACETYLAMINO)-2-DEOXY-A-D-GLUCOPYRANOSE) (three-letter code: NDG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
11	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	19	Total	O	0	0
			19	19		
12	B	20	Total	O	0	0
			20	20		
12	C	11	Total	O	0	0
			11	11		
12	D	9	Total	O	0	0
			9	9		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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: MEMBRANE COPPER AMINE OXIDASE



• Molecule 1: MEMBRANE COPPER AMINE OXIDASE





P377	G309	F227	L164	P91	GLY
	A378	G228	P165	S92	GLY
	A379	P310	P166	D93	ASP
	M380	A311	Y230	N94	GLY
	T381	P312	Y231	C95	GLY
		F313	R168	V96	GLU
	Y394	L314	P170	F97	PRO
	V395	Q315	V171	S98	SER
	D386	F316	L172	L101	LEU
	G387	Y317	H241	F173	PRO
P389	G388	H242	Q174	K106	HIS
	F389	P318	H243	A107	CYS
	G390	Q319	E175	A108	PRO
	M391	G320	Y176	A109	SER
	G392	P321	L177	L110	VAL
		R322	D180	A111	SER
	T395	Q326	L247	H112	PRO
	T396	G327	L248	H113	SER
	P397	S328	N250	D114	ALA
	R398	R329	H251	R115	GLM
P401	T399	V330	K252	R116	PRO
	R400	A331	A253	S117	THR
	G401	R400	R186	F118	HIS
	V402	L334	E187	P119	PRO
	D403	W335	L188	R122	GLY
	C404	T336	P189	E123	GLM
		F337	Q190	A124	SER
	L407	W259	A191	L125	Q58
	Y410	Q262	G193	V128	A61
	V411	K263	L194	F129	D62
P422	D412	V264	L195	L63	S64
		F265	H196	Q133	R65
	L416	Y266	H197	Q135	E66
	L417	Q267	C198	E67	L68
	L418	Y271	C199	V138	
	E418	D272	F200	S139	V71
	S419	S273	Y201	E140	R72
	R420	T274	K202	V143	R73
	Q421	F354	H203	G144	F74
	P422	F357	R204	L145	L75
P425	L425	Q276	R205	L146	T76
	R426	L277	R206	Y151	Q77
	D427	E278		M152	R78
	F432	F281	V209	G80	L79
	E433	T282	T210	P81	
	Q434	D283	R211	D154	G82
	M435	V288	T212	T155	L83
		L289	A214	V157	V84
	P439	Q366	L290	E158	D85
	L440	F367	T291	R159	A86
P442	R441	L369	R223	H160	Q87
	R442	A370	S301	G161	A88
		L365	L302	R224	S89
	S445	Y372	Q303	A224	R90
	Y448	G373	K304	T225	
	S449	G374	P305	V226	
	V450	C375	V226		



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	127.40Å 127.40Å 219.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.90 29.75 – 2.60	Depositor EDS
% Data completeness (in resolution range)	90.7 (20.00-2.90) 90.6 (29.75-2.60)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.61Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.215 , 0.276 0.216 , 0.211	Depositor DCC
R_{free} test set	3370 reflections (4.84%)	DCC
Wilson B-factor (Å ²)	55.3	Xtriage
Anisotropy	0.549	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 33.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.458 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	21917	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% 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: PAQ, BMA, NAG, CL, CA, NDG, FUL, CU, MAN

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	0.50	0/5488	0.72	2/7481 (0.0%)
1	B	0.48	0/5488	0.72	1/7481 (0.0%)
1	C	0.49	0/5488	0.71	2/7481 (0.0%)
1	D	0.48	0/5488	0.72	1/7481 (0.0%)
All	All	0.49	0/21952	0.72	6/29924 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	0
1	B	1	0
1	C	1	0
1	D	1	0
All	All	4	0

There are no bond length outliers.

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	59	LEU	CA-CB-CG	7.19	131.83	115.30
1	D	373	GLY	N-CA-C	-6.13	97.78	113.10
1	B	373	GLY	N-CA-C	-5.95	98.23	113.10
1	C	204	ARG	N-CA-C	5.47	125.78	111.00
1	A	204	ARG	N-CA-C	5.24	125.15	111.00

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	471	PAQ	CG
1	B	471	PAQ	CG
1	C	471	PAQ	CG
1	D	471	PAQ	CG

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	5340	0	5112	410	0
1	B	5340	0	5110	323	0
1	C	5340	0	5112	399	0
1	D	5340	0	5111	338	0
2	A	28	0	25	2	0
2	C	28	0	25	3	0
3	A	24	0	22	4	0
4	A	28	0	26	4	0
4	B	42	0	39	3	0
4	C	28	0	26	2	0
4	D	28	0	26	2	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
5	C	2	0	0	0	0
5	D	2	0	0	0	0
6	A	7	0	0	0	0
6	B	8	0	0	0	0
6	C	8	0	0	0	0
6	D	7	0	0	0	0
7	A	2	0	0	0	0
7	C	1	0	0	0	0
7	D	1	0	0	0	0
8	B	39	0	34	1	0
8	D	39	0	34	5	0
9	B	60	0	52	3	0
9	D	60	0	52	0	0
10	C	38	0	34	5	0
11	D	14	0	13	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
12	A	19	0	0	5	0
12	B	20	0	0	5	0
12	C	11	0	0	5	0
12	D	9	0	0	4	0
All	All	21917	0	20853	1388	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 32.

The worst 5 of 1388 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:360:VAL:HG11	1:B:363:ILE:HG13	1.37	1.04
1:C:90:ARG:HG2	1:C:91:PRO:HD2	1.42	1.02
1:B:106:LYS:HB2	1:B:637:ARG:HH21	1.27	0.99
1:A:90:ARG:HG2	1:A:91:PRO:HD2	1.43	0.96
1:A:344:PHE:HA	1:A:390:GLY:HA2	1.45	0.95

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	669/735 (91%)	571 (85%)	72 (11%)	26 (4%)	3	14
1	B	669/735 (91%)	585 (87%)	76 (11%)	8 (1%)	15	46
1	C	669/735 (91%)	575 (86%)	69 (10%)	25 (4%)	4	16
1	D	669/735 (91%)	587 (88%)	71 (11%)	11 (2%)	11	37
All	All	2676/2940 (91%)	2318 (87%)	288 (11%)	70 (3%)	6	24

5 of 70 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	380	MET
1	A	637	ARG
1	B	267	GLN
1	B	386	ASP
1	B	596	LYS

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	562/609 (92%)	523 (93%)	39 (7%)	18	46
1	B	562/609 (92%)	538 (96%)	24 (4%)	33	68
1	C	562/609 (92%)	527 (94%)	35 (6%)	21	52
1	D	562/609 (92%)	535 (95%)	27 (5%)	30	64
All	All	2248/2436 (92%)	2123 (94%)	125 (6%)	25	57

5 of 125 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	515	THR
1	C	203	HIS
1	D	513	GLU
1	B	528	VAL
1	C	88	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 55 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	510	GLN
1	C	219	GLN
1	D	420	GLN
1	B	560	GLN
1	C	160	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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$
1	PAQ	A	471	1	20,22,23	1.83	4 (20%)	13,29,31	2.17	3 (23%)
1	PAQ	B	471	1	20,22,23	1.83	4 (20%)	13,29,31	2.22	4 (30%)
1	PAQ	C	471	1	20,22,23	1.86	5 (25%)	13,29,31	2.24	3 (23%)
1	PAQ	D	471	1	20,22,23	1.83	4 (20%)	13,29,31	2.14	3 (23%)

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
1	PAQ	A	471	1	1/1/5/10	0/7/27/29	0/2/2/2
1	PAQ	B	471	1	1/1/5/10	0/7/27/29	0/2/2/2
1	PAQ	C	471	1	1/1/5/10	0/7/27/29	0/2/2/2
1	PAQ	D	471	1	1/1/5/10	0/7/27/29	0/2/2/2

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	471	PAQ	CG-CD2	-5.19	1.39	1.50
1	D	471	PAQ	CG-CD2	-5.16	1.39	1.50
1	C	471	PAQ	CG-CD2	-5.15	1.39	1.50
1	B	471	PAQ	CG-CD2	-5.14	1.39	1.50
1	C	471	PAQ	CE1-CD1	-3.22	1.39	1.45

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	471	PAQ	CD2-CE2-N1	-6.14	117.45	125.83
1	B	471	PAQ	CD2-CE2-N1	-5.90	117.78	125.83
1	A	471	PAQ	CD2-CE2-N1	-5.85	117.84	125.83
1	D	471	PAQ	CD2-CE2-N1	-5.50	118.33	125.83
1	B	471	PAQ	O-C-CA	-2.19	118.98	125.02

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	471	PAQ	CG
1	A	471	PAQ	CG
1	D	471	PAQ	CG
1	C	471	PAQ	CG

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	471	PAQ	5	0
1	B	471	PAQ	5	0
1	C	471	PAQ	4	0
1	D	471	PAQ	4	0

5.5 Carbohydrates [i](#)

25 carbohydrates 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	1730	1,2	14,14,15	0.52	0	15,19,21	1.09	1 (6%)
2	NAG	A	1731	2	14,14,15	0.54	0	15,19,21	0.86	1 (6%)
3	NAG	A	1732	1,3	14,14,15	0.80	1 (7%)	15,19,21	1.00	1 (6%)
3	FUL	A	1733	3	9,10,11	0.52	0	13,14,16	0.94	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	NAG	B	1730	1,8	14,14,15	0.75	0	15,19,21	0.91	0
8	NAG	B	1731	8	14,14,15	0.96	0	15,19,21	1.04	2 (13%)
9	NDG	B	1732	1,9	14,14,15	0.81	0	15,19,21	1.72	3 (20%)
9	NAG	B	1733	9	14,14,15	0.68	0	15,19,21	0.99	1 (6%)
9	FUL	B	1734	9	9,10,11	0.74	0	13,14,16	0.62	0
8	BMA	B	1737	8	11,11,12	0.94	0	13,15,17	0.61	0
9	BMA	B	1739	9	11,11,12	0.97	1 (9%)	13,15,17	1.26	1 (7%)
9	MAN	B	1740	9	11,11,12	0.73	0	13,15,17	0.90	1 (7%)
2	NAG	C	1730	1,2	14,14,15	0.60	0	15,19,21	0.79	0
2	NAG	C	1731	2	14,14,15	0.56	0	15,19,21	1.01	1 (6%)
10	NAG	C	1732	1,10	14,14,15	0.63	0	15,19,21	1.27	2 (13%)
10	NAG	C	1733	10	14,14,15	0.45	0	15,19,21	0.74	0
10	FUL	C	1734	10	9,10,11	0.50	0	13,14,16	0.67	0
8	NAG	D	1730	1,8	14,14,15	0.54	0	15,19,21	0.76	0
8	NAG	D	1731	8	14,14,15	0.75	0	15,19,21	0.84	0
9	NDG	D	1732	1,9	14,14,15	0.78	0	15,19,21	1.17	1 (6%)
9	NAG	D	1733	9	14,14,15	0.84	1 (7%)	15,19,21	0.97	1 (6%)
9	FUL	D	1734	9	9,10,11	0.84	0	13,14,16	0.63	0
8	BMA	D	1737	8	11,11,12	0.63	0	13,15,17	0.88	1 (7%)
9	BMA	D	1739	9	11,11,12	1.09	1 (9%)	13,15,17	0.65	0
9	MAN	D	1740	9	11,11,12	1.00	1 (9%)	13,15,17	0.70	1 (7%)

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	1730	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	1731	2	-	0/6/23/26	0/1/1/1
3	NAG	A	1732	1,3	-	0/6/23/26	0/1/1/1
3	FUL	A	1733	3	-	0/0/17/20	0/1/1/1
8	NAG	B	1730	1,8	-	0/6/23/26	0/1/1/1
8	NAG	B	1731	8	-	0/6/23/26	0/1/1/1
9	NDG	B	1732	1,9	-	0/6/23/26	0/1/1/1
9	NAG	B	1733	9	-	0/6/23/26	0/1/1/1
9	FUL	B	1734	9	-	0/0/17/20	0/1/1/1
8	BMA	B	1737	8	-	0/2/19/22	0/1/1/1
9	BMA	B	1739	9	-	0/2/19/22	0/1/1/1
9	MAN	B	1740	9	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	C	1730	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	1731	2	-	0/6/23/26	0/1/1/1
10	NAG	C	1732	1,10	-	0/6/23/26	0/1/1/1
10	NAG	C	1733	10	-	0/6/23/26	0/1/1/1
10	FUL	C	1734	10	-	0/0/17/20	0/1/1/1
8	NAG	D	1730	1,8	-	0/6/23/26	0/1/1/1
8	NAG	D	1731	8	-	0/6/23/26	0/1/1/1
9	NDG	D	1732	1,9	-	0/6/23/26	0/1/1/1
9	NAG	D	1733	9	-	1/6/23/26	0/1/1/1
9	FUL	D	1734	9	-	0/0/17/20	0/1/1/1
8	BMA	D	1737	8	-	0/2/19/22	0/1/1/1
9	BMA	D	1739	9	-	0/2/19/22	0/1/1/1
9	MAN	D	1740	9	-	0/2/19/22	1/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	D	1733	NAG	C1-C2	2.07	1.55	1.52
9	B	1739	BMA	C2-C3	2.15	1.55	1.52
3	A	1732	NAG	C1-C2	2.24	1.55	1.52
9	D	1740	MAN	C2-C3	2.49	1.55	1.52
9	D	1739	BMA	C2-C3	2.76	1.56	1.52

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	B	1732	NDG	C4-C3-C2	-4.13	104.97	111.02
9	B	1732	NDG	C3-C4-C5	-3.25	104.49	110.22
10	C	1732	NAG	C4-C3-C2	-3.14	106.42	111.02
8	D	1737	BMA	C1-C2-C3	-2.69	106.24	109.65
9	D	1732	NDG	C3-C4-C5	-2.62	105.61	110.22

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	D	1733	NAG	O7-C7-N2-C2

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	D	1740	MAN	C1-C2-C3-C4-C5-O5

16 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1730	NAG	2	0
2	A	1731	NAG	2	0
3	A	1732	NAG	2	0
3	A	1733	FUL	4	0
8	B	1730	NAG	1	0
9	B	1732	NDG	2	0
9	B	1733	NAG	2	0
9	B	1734	FUL	1	0
2	C	1730	NAG	3	0
2	C	1731	NAG	3	0
10	C	1732	NAG	5	0
10	C	1733	NAG	3	0
10	C	1734	FUL	2	0
8	D	1730	NAG	3	0
8	D	1731	NAG	3	0
8	D	1737	BMA	1	0

5.6 Ligand geometry

Of 52 ligands modelled in this entry, 42 are monoatomic - leaving 10 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	$\# Z > 2$
4	NAG	A	1734	1	14,14,15	0.68	0	15,19,21	0.65	0
4	NAG	A	1735	1	14,14,15	0.64	0	15,19,21	0.72	0
4	NAG	B	1735	1	14,14,15	0.70	0	15,19,21	0.82	0
4	NAG	B	1736	1	14,14,15	0.85	1 (7%)	15,19,21	0.67	0
4	NAG	B	1738	1	14,14,15	0.77	1 (7%)	15,19,21	0.58	0
4	NAG	C	1735	1	14,14,15	0.79	1 (7%)	15,19,21	0.70	0
4	NAG	C	1736	1	14,14,15	0.78	1 (7%)	15,19,21	0.75	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	D	1735	1	14,14,15	0.84	1 (7%)	15,19,21	0.71	0
11	NDG	D	1736	1	14,14,15	0.72	0	15,19,21	0.77	0
4	NAG	D	1738	1	14,14,15	1.03	1 (7%)	15,19,21	0.86	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
4	NAG	A	1734	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1735	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1735	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1736	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1738	1	-	0/6/23/26	0/1/1/1
4	NAG	C	1735	1	-	0/6/23/26	0/1/1/1
4	NAG	C	1736	1	-	0/6/23/26	0/1/1/1
4	NAG	D	1735	1	-	0/6/23/26	0/1/1/1
11	NDG	D	1736	1	-	0/6/23/26	0/1/1/1
4	NAG	D	1738	1	-	0/6/23/26	0/1/1/1

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1738	NAG	C1-C2	2.04	1.55	1.52
4	D	1735	NAG	C1-C2	2.16	1.55	1.52
4	C	1736	NAG	C1-C2	2.19	1.55	1.52
4	C	1735	NAG	C1-C2	2.25	1.55	1.52
4	B	1736	NAG	C1-C2	2.67	1.56	1.52

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	1736	NAG	C2-N2-C7	-2.01	120.01	122.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1734	NAG	4	0
4	B	1735	NAG	2	0
4	B	1738	NAG	1	0
4	C	1735	NAG	2	0
4	D	1735	NAG	2	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	671/735 (91%)	-0.24	0 100 100	43, 64, 79, 87	0
1	B	671/735 (91%)	-0.27	1 (0%) 95 95	41, 60, 75, 81	0
1	C	671/735 (91%)	-0.28	2 (0%) 93 93	43, 64, 79, 88	0
1	D	671/735 (91%)	-0.36	2 (0%) 93 93	41, 60, 75, 81	0
All	All	2684/2940 (91%)	-0.29	5 (0%) 94 95	41, 62, 77, 88	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	161	GLY	3.7
1	B	79	LEU	3.4
1	D	203	HIS	2.5
1	D	286	VAL	2.2
1	C	205	GLY	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
1	PAQ	B	471	21/22	0.81	0.35	-	55,59,61,61	0
1	PAQ	A	471	21/22	0.83	0.28	-	56,59,60,63	0
1	PAQ	D	471	21/22	0.89	0.27	-	54,59,60,61	0
1	PAQ	C	471	21/22	0.86	0.24	-	56,59,61,61	0

6.3 Carbohydrates

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
9	FUL	D	1734	10/11	0.91	0.26	1.88	84,85,86,86	0
2	NAG	A	1730	14/15	0.94	0.18	0.05	75,77,80,84	0
8	NAG	B	1730	14/15	0.92	0.18	-0.13	72,75,77,80	0
10	FUL	C	1734	10/11	0.96	0.15	-0.28	82,83,84,84	0
10	NAG	C	1732	14/15	0.89	0.15	-0.46	80,83,86,89	0
9	NDG	D	1732	14/15	0.94	0.13	-0.55	76,79,84,88	0
8	NAG	D	1730	14/15	0.95	0.16	-0.81	75,76,78,81	0
2	NAG	C	1730	14/15	0.94	0.12	-1.93	74,75,77,80	0
2	NAG	A	1731	14/15	0.94	0.15	-	88,91,92,92	0
3	NAG	A	1732	14/15	0.86	0.18	-	80,81,83,85	0
2	NAG	C	1731	14/15	0.94	0.15	-	82,84,85,85	0
10	NAG	C	1733	14/15	0.90	0.20	-	91,92,93,93	0
8	NAG	B	1731	14/15	0.90	0.13	-	85,86,89,92	0
9	NAG	D	1733	14/15	0.81	0.23	-	92,95,96,97	0
9	BMA	D	1739	11/12	0.94	0.12	-	97,98,98,98	0
9	MAN	B	1740	11/12	0.75	0.20	-	96,98,99,99	0
9	BMA	B	1739	11/12	0.93	0.10	-	97,97,98,98	0
8	BMA	D	1737	11/12	0.85	0.14	-	89,90,91,91	0
8	BMA	B	1737	11/12	0.88	0.10	-	94,94,95,95	0
8	NAG	D	1731	14/15	0.92	0.17	-	83,85,87,89	0
9	NAG	B	1733	14/15	0.93	0.11	-	95,97,98,98	0
3	FUL	A	1733	10/11	0.93	0.12	-	78,79,80,80	0
9	MAN	D	1740	11/12	0.79	0.17	-	96,98,98,98	0
9	FUL	B	1734	10/11	0.90	0.22	-	91,93,94,95	0
9	NDG	B	1732	14/15	0.91	0.15	-	82,86,91,93	0

6.4 Ligands

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
6	CU	A	1744	1/1	0.97	0.24	5.20	74,74,74,74	0
6	CU	B	1748	1/1	0.98	0.19	3.28	66,66,66,66	0
6	CU	B	1750	1/1	0.96	0.20	0.89	63,63,63,63	0
4	NAG	B	1735	14/15	0.91	0.20	0.44	69,70,73,73	0
5	CA	B	1743	1/1	0.98	0.18	0.25	47,47,47,47	0
6	CU	D	1749	1/1	0.97	0.17	0.18	67,67,67,67	0
6	CU	A	1745	1/1	0.98	0.14	-0.18	80,80,80,80	0
4	NAG	D	1735	14/15	0.91	0.15	-0.69	71,74,76,77	0
5	CA	A	1736	1/1	0.97	0.15	-0.78	52,52,52,52	0
4	NAG	C	1735	14/15	0.93	0.16	-0.79	70,72,73,74	0
5	CA	C	1739	1/1	0.99	0.13	-0.85	50,50,50,50	0
4	NAG	A	1734	14/15	0.91	0.14	-0.86	71,73,74,74	0
5	CA	B	1741	1/1	0.97	0.13	-1.08	46,46,46,46	0
6	CU	C	1747	1/1	0.97	0.10	-1.10	84,84,84,84	0
5	CA	D	1743	1/1	0.97	0.13	-1.12	46,46,46,46	0
6	CU	C	1745	1/1	0.99	0.14	-1.27	77,77,77,77	0
5	CA	A	1738	1/1	0.97	0.11	-1.28	49,49,49,49	0
5	CA	C	1737	1/1	0.99	0.11	-1.37	50,50,50,50	0
6	CU	D	1746	1/1	0.90	0.11	-1.38	78,78,78,78	0
7	CL	A	1746	1/1	0.96	0.10	-1.42	47,47,47,47	0
7	CL	D	1744	1/1	0.99	0.09	-1.61	45,45,45,45	0
5	CA	D	1741	1/1	0.93	0.11	-1.64	47,47,47,47	0
7	CL	A	1739	1/1	0.98	0.06	-2.19	47,47,47,47	0
7	CL	C	1740	1/1	0.98	0.10	-2.44	49,49,49,49	0
6	CU	D	1750	1/1	0.99	0.10	-2.93	67,67,67,67	0
6	CU	D	1748	1/1	0.99	0.13	-	74,74,74,74	0
4	NAG	D	1738	14/15	0.86	0.18	-	81,83,85,85	0
6	CU	B	1746	1/1	0.97	0.10	-	71,71,71,71	0
6	CU	D	1745	1/1	0.95	0.12	-	79,79,79,79	0
6	CU	C	1738	1/1	0.98	0.16	-	49,49,49,49	0
6	CU	C	1743	1/1	0.97	0.13	-	73,73,73,73	0
6	CU	A	1743	1/1	0.97	0.18	-	84,84,84,84	0
6	CU	B	1749	1/1	0.99	0.11	-	69,69,69,69	0
4	NAG	B	1738	14/15	0.86	0.14	-	81,83,84,85	0
6	CU	B	1744	1/1	0.97	0.20	-	76,76,76,76	0
11	NDG	D	1736	14/15	0.86	0.13	-	83,84,86,86	0
6	CU	C	1741	1/1	0.93	0.12	-	79,79,79,79	0
6	CU	C	1746	1/1	0.99	0.09	-	67,67,67,67	0
4	NAG	B	1736	14/15	0.90	0.10	-	84,86,88,88	0
4	NAG	A	1735	14/15	0.88	0.16	-	84,86,88,89	0
6	CU	A	1737	1/1	0.99	0.21	-	49,49,49,49	0
6	CU	B	1745	1/1	0.97	0.14	-	75,75,75,75	0
4	NAG	C	1736	14/15	0.86	0.14	-	82,85,88,88	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
6	CU	A	1740	1/1	0.89	0.18	-	81,81,81,81	0
6	CU	D	1747	1/1	0.95	0.05	-	82,82,82,82	0
6	CU	C	1744	1/1	0.98	0.14	-	90,90,90,90	0
6	CU	D	1742	1/1	0.98	0.15	-	52,52,52,52	0
6	CU	B	1747	1/1	0.96	0.15	-	81,81,81,81	0
6	CU	B	1742	1/1	0.96	0.20	-	55,55,55,55	0
6	CU	C	1742	1/1	0.96	0.09	-	84,84,84,84	0
6	CU	A	1741	1/1	0.98	0.10	-	83,83,83,83	0
6	CU	A	1742	1/1	0.96	0.13	-	77,77,77,77	0

6.5 Other polymers [i](#)

There are no such residues in this entry.