



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 08:27 AM BST

PDB ID : 1Y8J
Title : Crystal Structure of human NEP complexed with an imidazo[4,5-c]pyridine inhibitor
Authors : Sahli, S.; Frank, B.; Schweizer, W.B.; Diederich, F.; Blum-Kaelin, D.; Aebi, J.D.; Bohm, H.J.; Oefner, C.; Dale, G.E.
Deposited on : 2004-12-13
Resolution : 2.25 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

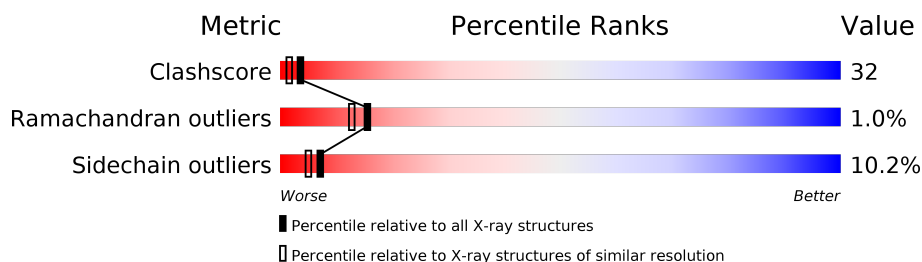
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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))
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)

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 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	696	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	ACT	A	801	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5873 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 Neprilysin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	696	5595	3538	957	1074	26	0	0	0

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



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

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

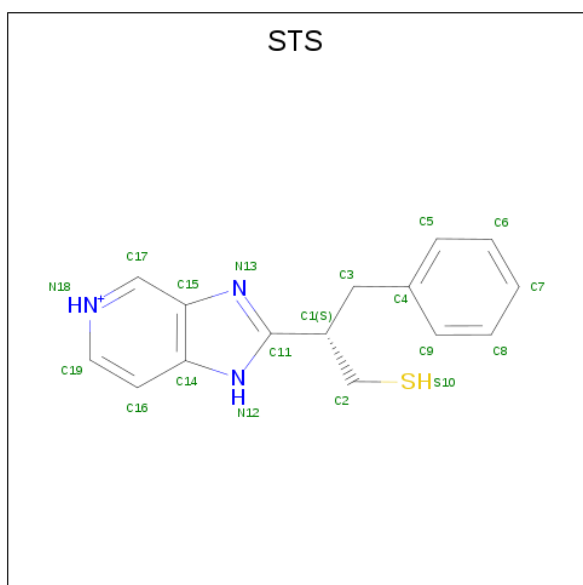
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



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

- Molecule 5 is 2-[(1S)-1-BENZYL-2-SULFANYLETHYL]-1H-IMIDAZO[4,5-C]PYRIDIN-5-ILUM (three-letter code: STS) (formula: C₁₅H₁₆N₃S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	S	0	0
			19	15	3	1		

- Molecule 6 is water.

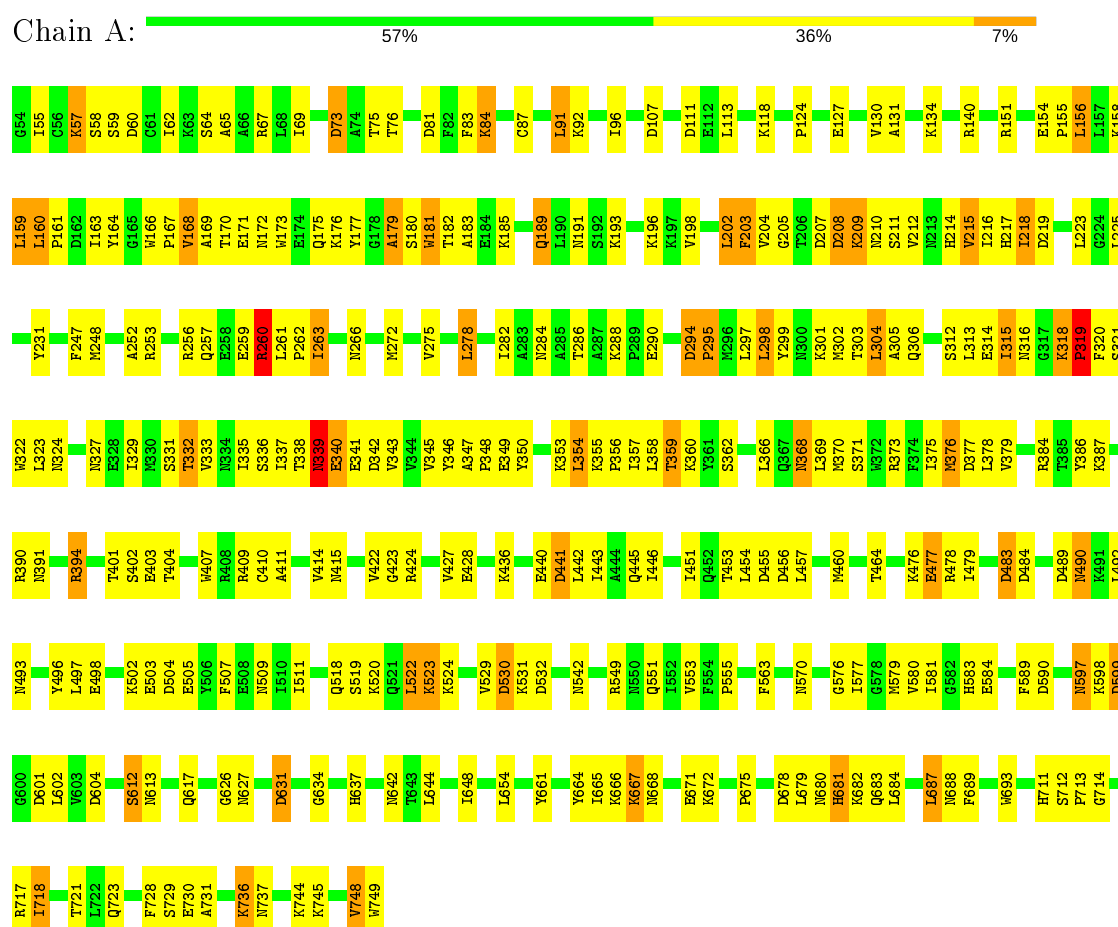
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	212	Total	O	0	0
			212	212		

3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

• Molecule 1: Neprilysin



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	107.36 Å 107.36 Å 112.47 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.25	Depositor
% Data completeness (in resolution range)	95.5 (20.00-2.25)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.226 , 0.296	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5873	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, STS, NAG, ACT

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	0.40	0/5713	0.63	18/7727 (0.2%)

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	0	2

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	294	ASP	CB-CG-OD2	5.93	123.64	118.30
1	A	530	ASP	CB-CG-OD2	5.89	123.60	118.30
1	A	107	ASP	CB-CG-OD2	5.83	123.54	118.30
1	A	678	ASP	CB-CG-OD2	5.58	123.32	118.30
1	A	441	ASP	CB-CG-OD2	5.50	123.25	118.30
1	A	111	ASP	CB-CG-OD2	5.48	123.23	118.30
1	A	377	ASP	CB-CG-OD2	5.46	123.21	118.30
1	A	219	ASP	CB-CG-OD2	5.38	123.14	118.30
1	A	208	ASP	CB-CG-OD2	5.36	123.13	118.30
1	A	631	ASP	CB-CG-OD2	5.35	123.11	118.30
1	A	599	ASP	CB-CG-OD2	5.31	123.08	118.30
1	A	504	ASP	CB-CG-OD2	5.30	123.07	118.30
1	A	455	ASP	CB-CG-OD2	5.28	123.05	118.30
1	A	73	ASP	CB-CG-OD2	5.27	123.04	118.30
1	A	483	ASP	CB-CG-OD2	5.24	123.02	118.30
1	A	60	ASP	CB-CG-OD2	5.20	122.98	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	456	ASP	CB-CG-OD2	5.17	122.95	118.30
1	A	489	ASP	CB-CG-OD2	5.10	122.89	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	181	TRP	Peptide
1	A	319	PRO	Peptide

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	5595	0	5446	353	4
2	A	42	0	39	1	0
3	A	1	0	0	0	0
4	A	4	0	3	5	0
5	A	19	0	15	4	0
6	A	212	0	0	21	0
All	All	5873	0	5503	353	4

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.

All (353) 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:522:LEU:CD2	1:A:522:LEU:H	1.31	1.32
1:A:523:LYS:NZ	1:A:523:LYS:CB	1.94	1.22
1:A:256:ARG:NH1	1:A:262:PRO:O	1.73	1.20
1:A:318:LYS:HB3	1:A:319:PRO:CA	1.70	1.19
1:A:353:LYS:O	1:A:357:ILE:CD1	1.94	1.15
1:A:313:LEU:O	1:A:319:PRO:HB2	1.45	1.14
1:A:522:LEU:N	1:A:522:LEU:HD23	1.50	1.13

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:LYS:CB	1:A:319:PRO:HA	1.70	1.12
1:A:338:THR:HG22	1:A:340:GLU:N	1.64	1.10
1:A:523:LYS:HZ3	1:A:523:LYS:CB	1.55	1.10
1:A:714:GLY:O	1:A:718:ILE:HD13	1.52	1.08
1:A:523:LYS:HZ2	1:A:523:LYS:HB2	0.98	1.08
1:A:338:THR:HG22	1:A:340:GLU:H	0.90	1.06
1:A:579:MET:HE2	6:A:909:HOH:O	1.53	1.05
1:A:523:LYS:HZ3	1:A:523:LYS:HB3	0.91	1.05
1:A:202:LEU:HD23	1:A:218:ILE:HD13	1.10	1.05
1:A:460:MET:CE	1:A:464:THR:HG22	1.86	1.05
1:A:523:LYS:NZ	1:A:523:LYS:HB2	1.55	1.03
1:A:318:LYS:HB3	1:A:319:PRO:C	1.78	1.03
1:A:460:MET:HE2	1:A:464:THR:HG22	1.37	1.03
1:A:202:LEU:HD23	1:A:218:ILE:CD1	1.89	1.02
1:A:522:LEU:CD2	1:A:522:LEU:N	2.05	0.99
1:A:626:GLY:O	1:A:637:HIS:HD2	1.44	0.99
1:A:202:LEU:CD2	1:A:218:ILE:HD13	1.93	0.98
1:A:159:LEU:HD12	1:A:159:LEU:O	1.63	0.98
1:A:661:TYR:CZ	1:A:665:ILE:HD11	1.99	0.97
1:A:209:LYS:HE2	1:A:299:TYR:CZ	1.98	0.96
1:A:333:VAL:HG12	1:A:333:VAL:O	1.66	0.95
1:A:671:GLU:HB2	1:A:681:HIS:CD2	2.01	0.95
1:A:579:MET:CE	6:A:909:HOH:O	2.10	0.95
1:A:318:LYS:HB2	1:A:319:PRO:HA	1.49	0.94
1:A:253:ARG:O	1:A:257:GLN:HG3	1.68	0.94
1:A:257:GLN:O	6:A:937:HOH:O	1.85	0.93
1:A:318:LYS:CB	1:A:319:PRO:CA	2.34	0.92
1:A:626:GLY:O	1:A:637:HIS:CD2	2.23	0.91
1:A:323:LEU:HD22	1:A:339:ASN:OD1	1.70	0.91
1:A:353:LYS:O	1:A:357:ILE:HD12	1.72	0.90
1:A:202:LEU:HD22	1:A:217:HIS:O	1.72	0.90
1:A:209:LYS:CE	1:A:299:TYR:CZ	2.34	0.89
1:A:202:LEU:CD2	1:A:218:ILE:CD1	2.50	0.89
1:A:338:THR:CG2	1:A:340:GLU:H	1.84	0.88
1:A:671:GLU:HB2	1:A:681:HIS:NE2	1.88	0.88
1:A:404:THR:CG2	1:A:409:ARG:HG3	2.04	0.87
1:A:661:TYR:CE2	1:A:665:ILE:HD11	2.10	0.87
1:A:323:LEU:HD22	1:A:339:ASN:CG	1.96	0.86
1:A:223:LEU:HD23	4:A:801:ACT:H3	1.58	0.85
1:A:523:LYS:NZ	1:A:523:LYS:HB3	1.68	0.85
1:A:391:ASN:HB2	6:A:1001:HOH:O	1.77	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:ASP:OD1	1:A:75:THR:HB	1.77	0.85
1:A:140:ARG:NH2	1:A:503:GLU:OE1	2.11	0.82
1:A:154:GLU:OE1	1:A:158:LYS:HD2	1.80	0.82
1:A:530:ASP:O	1:A:549:ARG:NH2	2.14	0.81
1:A:315:ILE:CD1	1:A:315:ILE:N	2.43	0.80
1:A:256:ARG:HB2	1:A:263:ILE:HD11	1.62	0.80
1:A:338:THR:CG2	1:A:340:GLU:CD	2.50	0.80
1:A:440:GLU:HG3	1:A:479:ILE:HD13	1.64	0.80
1:A:315:ILE:HD13	1:A:315:ILE:N	1.97	0.79
1:A:67:ARG:HH22	1:A:688:ASN:HD21	1.31	0.79
1:A:714:GLY:O	1:A:718:ILE:CD1	2.31	0.79
1:A:198:VAL:HG21	1:A:370:MET:HG2	1.64	0.79
1:A:131:ALA:CB	1:A:492:LEU:HD12	2.13	0.78
1:A:522:LEU:HD23	1:A:522:LEU:H	0.64	0.78
1:A:288:LYS:HD2	6:A:1058:HOH:O	1.81	0.78
1:A:202:LEU:HD11	1:A:216:ILE:HG23	1.66	0.78
1:A:177:TYR:O	1:A:180:SER:N	2.17	0.77
1:A:259:GLU:HA	1:A:259:GLU:OE2	1.84	0.77
1:A:693:TRP:HB2	1:A:718:ILE:HD11	1.67	0.77
1:A:202:LEU:HD11	1:A:216:ILE:CG2	2.14	0.77
1:A:318:LYS:HB3	1:A:319:PRO:O	1.84	0.77
1:A:131:ALA:HB1	1:A:492:LEU:CD1	2.15	0.76
1:A:168:VAL:H	1:A:368:ASN:HD21	1.33	0.76
1:A:131:ALA:CB	1:A:492:LEU:CD1	2.64	0.76
1:A:313:LEU:HB3	1:A:315:ILE:HD11	1.67	0.76
1:A:718:ILE:HD12	1:A:718:ILE:N	2.00	0.76
1:A:353:LYS:O	1:A:357:ILE:HD13	1.83	0.76
1:A:718:ILE:H	1:A:718:ILE:CD1	1.97	0.76
1:A:355:LYS:O	1:A:359:THR:HG23	1.86	0.75
1:A:189:GLN:HE22	1:A:193:LYS:HD2	1.52	0.75
1:A:248:MET:HE1	1:A:275:VAL:CG1	2.17	0.75
1:A:460:MET:CE	1:A:464:THR:CG2	2.64	0.75
1:A:748:VAL:HG22	1:A:749:TRP:N	2.01	0.75
1:A:597:ASN:HD22	1:A:597:ASN:C	1.86	0.74
1:A:260:ARG:HB2	1:A:260:ARG:HH11	1.52	0.74
1:A:522:LEU:N	1:A:522:LEU:HD22	2.03	0.74
1:A:202:LEU:CD1	1:A:216:ILE:HG23	2.18	0.73
1:A:140:ARG:HH21	1:A:503:GLU:CD	1.93	0.72
1:A:338:THR:HG21	1:A:340:GLU:CD	2.09	0.72
1:A:736:LYS:O	1:A:737:ASN:HB2	1.87	0.71
1:A:671:GLU:HB2	1:A:681:HIS:HE2	1.54	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:627:ASN:HA	1:A:637:HIS:CD2	2.25	0.71
1:A:748:VAL:HG22	1:A:749:TRP:H	1.54	0.71
1:A:563:PHE:HB3	1:A:577:ILE:HD13	1.73	0.71
1:A:477:GLU:HG3	1:A:479:ILE:HD11	1.71	0.70
1:A:248:MET:HE1	1:A:275:VAL:HG12	1.72	0.70
1:A:345:VAL:HG12	1:A:348:PRO:HD3	1.73	0.70
1:A:333:VAL:O	1:A:333:VAL:CG1	2.39	0.70
1:A:203:PHE:HE1	1:A:205:GLY:HA3	1.57	0.69
1:A:67:ARG:HH22	1:A:688:ASN:ND2	1.89	0.69
1:A:313:LEU:O	1:A:319:PRO:CB	2.33	0.69
1:A:579:MET:HE1	6:A:908:HOH:O	1.91	0.69
1:A:717:ARG:O	1:A:721:THR:HG23	1.91	0.69
1:A:661:TYR:OH	1:A:671:GLU:OE1	2.03	0.69
1:A:681:HIS:ND1	1:A:684:LEU:HD12	2.08	0.69
1:A:171:GLU:O	1:A:176:LYS:HE3	1.93	0.69
1:A:163:ILE:O	1:A:164:TYR:HB2	1.92	0.69
1:A:723:GLN:O	1:A:745:LYS:HD3	1.92	0.69
1:A:87:CYS:O	1:A:91:LEU:HD22	1.93	0.69
1:A:355:LYS:HB3	1:A:356:PRO:HD3	1.74	0.68
1:A:167:PRO:HD2	1:A:368:ASN:HD22	1.58	0.68
1:A:424:ARG:O	1:A:428:GLU:HG3	1.93	0.68
1:A:371:SER:O	1:A:375:ILE:HG13	1.93	0.68
1:A:260:ARG:O	1:A:261:LEU:HD23	1.94	0.68
1:A:597:ASN:ND2	1:A:599:ASP:H	1.91	0.68
1:A:333:VAL:HG11	1:A:523:LYS:NZ	2.09	0.67
1:A:718:ILE:H	1:A:718:ILE:HD12	1.57	0.67
1:A:204:VAL:HG21	1:A:522:LEU:HA	1.76	0.67
1:A:159:LEU:HD12	1:A:159:LEU:C	2.13	0.67
1:A:131:ALA:HB3	1:A:492:LEU:HD12	1.78	0.66
1:A:218:ILE:HD12	1:A:322:TRP:CZ3	2.31	0.66
1:A:181:TRP:CE3	1:A:182:THR:HA	2.31	0.66
1:A:355:LYS:O	1:A:359:THR:CG2	2.44	0.65
1:A:181:TRP:O	1:A:182:THR:CG2	2.43	0.65
1:A:718:ILE:N	1:A:718:ILE:CD1	2.58	0.65
1:A:748:VAL:CG2	1:A:749:TRP:N	2.58	0.65
1:A:327:ASN:O	1:A:331:SER:HB3	1.95	0.65
1:A:597:ASN:ND2	1:A:597:ASN:C	2.50	0.65
1:A:502:LYS:H	1:A:509:ASN:HD21	1.45	0.64
1:A:549:ARG:NH1	1:A:549:ARG:HG2	2.12	0.64
1:A:198:VAL:CG2	1:A:370:MET:HG2	2.27	0.64
1:A:196:LYS:HE3	1:A:373:ARG:O	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:209:LYS:HE2	1:A:299:TYR:CE2	2.33	0.63
1:A:282:ILE:O	1:A:286:THR:HG23	1.99	0.63
1:A:314:GLU:HA	1:A:319:PRO:HB3	1.78	0.63
1:A:218:ILE:HD12	1:A:322:TRP:HZ3	1.63	0.63
1:A:156:LEU:HD13	1:A:160:LEU:HD22	1.80	0.62
1:A:338:THR:HG21	1:A:340:GLU:OE1	2.00	0.62
1:A:404:THR:HG22	1:A:409:ARG:HG3	1.80	0.62
1:A:312:SER:OG	1:A:355:LYS:NZ	2.31	0.62
1:A:549:ARG:HG2	1:A:549:ARG:HH11	1.64	0.62
1:A:318:LYS:HD2	1:A:319:PRO:O	2.00	0.61
1:A:355:LYS:HB3	1:A:356:PRO:CD	2.30	0.61
1:A:156:LEU:HD21	1:A:378:LEU:HD12	1.83	0.61
1:A:394:ARG:HH21	1:A:401:THR:C	2.04	0.61
1:A:181:TRP:O	1:A:182:THR:HG22	2.01	0.60
1:A:387:LYS:HE2	6:A:939:HOH:O	2.00	0.60
1:A:290:GLU:HG3	1:A:642:ASN:OD1	2.01	0.60
1:A:248:MET:CE	1:A:275:VAL:HG11	2.31	0.60
1:A:484:ASP:HB2	6:A:1018:HOH:O	2.00	0.60
1:A:679:LEU:HA	1:A:683:GLN:OE1	2.01	0.60
1:A:338:THR:HG21	1:A:340:GLU:HB2	1.81	0.60
1:A:171:GLU:O	1:A:176:LYS:CE	2.49	0.60
1:A:347:ALA:N	1:A:348:PRO:CD	2.65	0.59
1:A:661:TYR:CE2	1:A:665:ILE:CD1	2.85	0.59
1:A:203:PHE:CE1	1:A:205:GLY:HA3	2.36	0.59
1:A:207:ASP:HB2	1:A:215:VAL:CG2	2.32	0.59
1:A:167:PRO:HD2	1:A:368:ASN:ND2	2.17	0.59
1:A:440:GLU:HG3	1:A:479:ILE:CD1	2.31	0.59
1:A:207:ASP:HB2	1:A:215:VAL:HG22	1.83	0.59
1:A:65:ALA:O	1:A:69:ILE:HG13	2.03	0.59
1:A:303:THR:OG1	1:A:305:ALA:HB3	2.03	0.59
1:A:379:VAL:HG22	1:A:386:TYR:HB3	1.85	0.58
1:A:248:MET:CE	1:A:275:VAL:CG1	2.81	0.58
1:A:248:MET:HE2	1:A:275:VAL:HG11	1.84	0.58
1:A:189:GLN:NE2	1:A:193:LYS:HD2	2.19	0.58
1:A:654:LEU:HD13	1:A:689:PHE:CD2	2.38	0.58
1:A:518:GLN:HG3	1:A:522:LEU:HD11	1.87	0.57
1:A:76:THR:HG21	1:A:84:LYS:HB3	1.86	0.57
1:A:282:ILE:HG21	4:A:801:ACT:CH3	2.35	0.56
1:A:338:THR:CG2	1:A:340:GLU:HB2	2.35	0.56
1:A:284:ASN:HB3	6:A:1051:HOH:O	2.06	0.56
1:A:589:PHE:HB3	1:A:749:TRP:CZ2	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:384:ARG:NH1	1:A:387:LYS:HD2	2.21	0.56
1:A:298:LEU:HD13	1:A:346:TYR:CD1	2.41	0.56
1:A:55:ILE:HG21	1:A:675:PRO:HB3	1.88	0.56
1:A:214:HIS:CE1	1:A:524:LYS:HB3	2.41	0.56
1:A:67:ARG:NH2	1:A:688:ASN:HD21	2.01	0.56
1:A:453:THR:O	1:A:457:LEU:HG	2.05	0.56
1:A:210:ASN:HA	1:A:601:ASP:OD1	2.05	0.55
1:A:484:ASP:CB	6:A:1018:HOH:O	2.55	0.55
1:A:680:ASN:OD1	1:A:683:GLN:HG3	2.06	0.55
1:A:211:SER:O	1:A:529:VAL:HG21	2.06	0.55
1:A:202:LEU:HD21	1:A:218:ILE:CD1	2.37	0.55
1:A:671:GLU:CB	1:A:681:HIS:CD2	2.84	0.54
1:A:67:ARG:HH12	1:A:688:ASN:ND2	2.06	0.54
1:A:332:THR:HG22	1:A:333:VAL:HG23	1.90	0.54
1:A:460:MET:HE3	1:A:464:THR:CG2	2.38	0.54
1:A:151:ARG:CZ	1:A:505:GLU:OE1	2.56	0.54
1:A:333:VAL:CG1	1:A:523:LYS:NZ	2.71	0.54
1:A:590:ASP:C	1:A:590:ASP:OD1	2.46	0.54
1:A:353:LYS:O	1:A:357:ILE:HD11	2.02	0.54
1:A:338:THR:C	1:A:340:GLU:H	2.11	0.54
1:A:127:GLU:O	1:A:127:GLU:OE2	2.26	0.53
1:A:314:GLU:C	1:A:315:ILE:CD1	2.77	0.53
1:A:423:GLY:O	1:A:427:VAL:HG22	2.08	0.53
1:A:154:GLU:N	1:A:155:PRO:CD	2.70	0.53
1:A:338:THR:HG22	1:A:340:GLU:CA	2.36	0.53
1:A:67:ARG:HH12	1:A:688:ASN:HD22	1.55	0.53
1:A:394:ARG:HH21	1:A:402:SER:N	2.07	0.53
1:A:127:GLU:CD	1:A:127:GLU:C	2.68	0.52
1:A:315:ILE:O	1:A:316:ASN:HB3	2.09	0.52
1:A:492:LEU:O	1:A:492:LEU:HD13	2.09	0.52
1:A:338:THR:HG23	1:A:340:GLU:OE2	2.10	0.52
1:A:333:VAL:CG1	1:A:523:LYS:HZ1	2.22	0.52
1:A:55:ILE:O	1:A:57:LYS:HD2	2.10	0.52
1:A:302:MET:HG2	1:A:306:GLN:HB2	1.92	0.52
1:A:329:ILE:O	1:A:332:THR:HB	2.10	0.52
1:A:208:ASP:O	1:A:602:LEU:N	2.33	0.52
1:A:202:LEU:HD11	1:A:216:ILE:HG21	1.88	0.52
1:A:183:ALA:HB2	1:A:315:ILE:HG21	1.90	0.51
1:A:375:ILE:O	1:A:376:MET:C	2.48	0.51
1:A:282:ILE:HG21	4:A:801:ACT:H1	1.93	0.51
1:A:217:HIS:CE1	6:A:986:HOH:O	2.63	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:263:ILE:O	1:A:263:ILE:HG22	2.10	0.51
1:A:338:THR:HG22	1:A:340:GLU:CB	2.40	0.51
1:A:260:ARG:C	1:A:261:LEU:HD23	2.31	0.51
1:A:314:GLU:HA	1:A:319:PRO:CB	2.41	0.51
1:A:362:SER:O	1:A:366:LEU:HG	2.11	0.51
1:A:597:ASN:ND2	1:A:601:ASP:H	2.09	0.51
1:A:215:VAL:HG12	6:A:931:HOH:O	2.11	0.50
1:A:492:LEU:HD13	1:A:492:LEU:C	2.32	0.50
1:A:604:ASP:OD2	6:A:1037:HOH:O	2.19	0.50
1:A:327:ASN:O	1:A:331:SER:CB	2.60	0.50
1:A:627:ASN:CA	1:A:637:HIS:CD2	2.94	0.50
1:A:693:TRP:CB	1:A:718:ILE:HD11	2.41	0.50
1:A:338:THR:CG2	1:A:340:GLU:CB	2.89	0.50
1:A:414:VAL:HG23	1:A:415:ASN:N	2.27	0.50
1:A:253:ARG:HG3	1:A:263:ILE:HG12	1.93	0.50
1:A:321:SER:HB3	1:A:324:ASN:HB3	1.94	0.50
1:A:387:LYS:HD3	1:A:403:GLU:OE2	2.12	0.50
1:A:156:LEU:HD13	1:A:160:LEU:CD2	2.42	0.50
1:A:338:THR:HG23	1:A:340:GLU:CD	2.32	0.49
1:A:345:VAL:O	1:A:345:VAL:HG12	2.13	0.49
1:A:442:LEU:O	1:A:446:ILE:HD12	2.13	0.49
1:A:55:ILE:N	1:A:55:ILE:HD13	2.26	0.49
1:A:257:GLN:NE2	6:A:946:HOH:O	2.46	0.48
1:A:338:THR:C	1:A:340:GLU:N	2.65	0.48
1:A:58:SER:O	1:A:62:ILE:HD13	2.13	0.48
1:A:181:TRP:O	1:A:182:THR:HG23	2.13	0.48
1:A:584:GLU:OE2	5:A:900:STS:H22	2.13	0.48
1:A:671:GLU:CG	1:A:681:HIS:HD2	2.26	0.48
1:A:414:VAL:HG23	1:A:415:ASN:H	1.79	0.48
1:A:577:ILE:O	1:A:581:ILE:HG13	2.14	0.47
1:A:443:ILE:HG12	1:A:581:ILE:HG21	1.96	0.47
1:A:83:PHE:O	1:A:87:CYS:HB2	2.14	0.47
1:A:358:LEU:C	1:A:360:LYS:H	2.16	0.47
1:A:130:VAL:O	1:A:134:LYS:HG3	2.14	0.47
1:A:460:MET:HE2	1:A:464:THR:CG2	2.26	0.47
1:A:530:ASP:OD1	1:A:532:ASP:N	2.31	0.47
1:A:189:GLN:HA	1:A:189:GLN:OE1	2.14	0.47
1:A:191:ASN:OD1	1:A:191:ASN:C	2.53	0.47
1:A:304:LEU:HB3	1:A:339:ASN:HD22	1.80	0.47
1:A:314:GLU:C	1:A:315:ILE:HD13	2.35	0.47
1:A:424:ARG:O	1:A:427:VAL:CG2	2.63	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:VAL:HG13	1:A:523:LYS:HZ1	1.80	0.46
1:A:304:LEU:CD2	1:A:343:VAL:HG21	2.45	0.46
1:A:301:LYS:HE2	1:A:342:ASP:HB3	1.98	0.46
1:A:712:SER:O	1:A:717:ARG:HD2	2.15	0.46
1:A:748:VAL:CG2	1:A:749:TRP:CD1	2.98	0.46
1:A:410:CYS:O	1:A:414:VAL:HG22	2.16	0.46
1:A:202:LEU:HD22	1:A:202:LEU:HA	1.77	0.46
1:A:305:ALA:O	1:A:306:GLN:C	2.54	0.46
1:A:327:ASN:HB3	2:A:753:NAG:H81	1.98	0.45
1:A:303:THR:HA	1:A:341:GLU:O	2.15	0.45
1:A:225:LEU:HD12	1:A:231:TYR:CZ	2.51	0.45
1:A:314:GLU:C	1:A:315:ILE:HD12	2.36	0.45
1:A:336:SER:O	1:A:337:ILE:HD13	2.16	0.45
1:A:350:TYR:CE1	1:A:354:LEU:HD22	2.52	0.45
1:A:313:LEU:HB3	1:A:315:ILE:CD1	2.43	0.45
1:A:350:TYR:CZ	1:A:354:LEU:HD22	2.51	0.45
1:A:214:HIS:ND1	1:A:524:LYS:O	2.49	0.45
1:A:354:LEU:HA	1:A:357:ILE:HD13	1.98	0.45
1:A:507:PHE:O	1:A:511:ILE:HG12	2.16	0.45
1:A:542:ASN:O	1:A:555:PRO:HD2	2.17	0.45
1:A:570:ASN:ND2	1:A:664:TYR:OH	2.49	0.45
1:A:627:ASN:HA	1:A:637:HIS:HD2	1.76	0.45
1:A:644:LEU:HG	1:A:648:ILE:HD12	1.99	0.45
1:A:667:LYS:HG2	1:A:668:ASN:ND2	2.32	0.45
1:A:172:ASN:O	1:A:176:LYS:HG3	2.16	0.45
1:A:667:LYS:O	1:A:667:LYS:HG3	2.15	0.45
1:A:282:ILE:HG21	4:A:801:ACT:H2	1.99	0.44
1:A:404:THR:HG21	1:A:409:ARG:HG3	1.97	0.44
1:A:530:ASP:OD1	1:A:532:ASP:HB2	2.17	0.44
1:A:576:GLY:O	1:A:580:VAL:HG23	2.17	0.44
1:A:256:ARG:CB	1:A:263:ILE:HD11	2.41	0.44
1:A:713:PRO:HG3	6:A:1022:HOH:O	2.17	0.44
1:A:167:PRO:O	1:A:169:ALA:N	2.51	0.44
1:A:612:SER:OG	6:A:1036:HOH:O	2.15	0.44
1:A:424:ARG:O	1:A:427:VAL:HG22	2.17	0.44
1:A:252:ALA:O	1:A:256:ARG:HG3	2.17	0.44
1:A:613:ASN:O	1:A:617:GLN:HG2	2.18	0.43
1:A:202:LEU:HD13	1:A:216:ILE:HG23	1.97	0.43
1:A:57:LYS:HB3	1:A:57:LYS:HE2	1.57	0.43
1:A:217:HIS:HE1	6:A:988:HOH:O	1.99	0.43
1:A:394:ARG:NH2	1:A:401:THR:C	2.71	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:711:HIS:CD2	5:A:900:STS:N12	2.86	0.43
1:A:272:MET:HA	1:A:272:MET:HE2	2.00	0.43
1:A:223:LEU:HD23	4:A:801:ACT:CH3	2.39	0.43
1:A:335:ILE:HG22	1:A:336:SER:N	2.33	0.43
1:A:577:ILE:HD12	1:A:577:ILE:HA	1.74	0.43
1:A:278:LEU:HD13	1:A:369:LEU:HD23	2.01	0.43
1:A:212:VAL:C	1:A:529:VAL:HG23	2.38	0.43
1:A:272:MET:HA	1:A:272:MET:CE	2.48	0.43
1:A:490:ASN:HA	1:A:490:ASN:HD22	1.67	0.43
1:A:687:LEU:HA	1:A:687:LEU:HD23	1.75	0.43
1:A:166:TRP:CH2	1:A:252:ALA:HB2	2.53	0.43
1:A:304:LEU:HB3	1:A:339:ASN:ND2	2.34	0.42
1:A:549:ARG:CG	1:A:549:ARG:HH11	2.30	0.42
1:A:530:ASP:OD1	1:A:531:LYS:N	2.51	0.42
1:A:583:HIS:CD2	5:A:900:STS:H9	2.54	0.42
1:A:202:LEU:HD21	1:A:218:ILE:HD11	2.00	0.42
1:A:530:ASP:OD1	1:A:530:ASP:C	2.54	0.42
1:A:579:MET:HE1	5:A:900:STS:H7	2.01	0.42
1:A:131:ALA:HB2	1:A:493:ASN:HA	2.01	0.42
1:A:247:PHE:CZ	1:A:390:ARG:HG3	2.54	0.42
1:A:302:MET:O	1:A:343:VAL:N	2.43	0.42
1:A:379:VAL:CG2	1:A:386:TYR:HB3	2.48	0.42
1:A:118:LYS:HB2	1:A:407:TRP:CE3	2.55	0.42
1:A:225:LEU:HD12	1:A:231:TYR:CE2	2.54	0.42
1:A:358:LEU:O	1:A:360:LYS:N	2.53	0.42
1:A:415:ASN:HA	1:A:422:VAL:HG21	2.02	0.42
1:A:627:ASN:CA	1:A:637:HIS:HD2	2.32	0.42
1:A:748:VAL:HG22	1:A:749:TRP:CD1	2.55	0.42
1:A:156:LEU:O	1:A:159:LEU:HB3	2.19	0.42
1:A:294:ASP:HA	1:A:295:PRO:HD3	1.89	0.41
1:A:627:ASN:O	1:A:637:HIS:NE2	2.53	0.41
1:A:81:ASP:C	1:A:81:ASP:OD1	2.58	0.41
1:A:728:PHE:O	1:A:731:ALA:HB3	2.20	0.41
1:A:124:PRO:HA	6:A:1014:HOH:O	2.20	0.41
1:A:394:ARG:NH2	1:A:402:SER:N	2.69	0.41
1:A:478:ARG:HB2	1:A:553:VAL:HG13	2.02	0.41
1:A:160:LEU:N	1:A:161:PRO:CD	2.83	0.41
1:A:318:LYS:HB3	1:A:319:PRO:HA	1.40	0.41
1:A:319:PRO:HB2	1:A:320:PHE:H	1.71	0.41
1:A:631:ASP:O	1:A:634:GLY:N	2.44	0.41
1:A:369:LEU:HD12	1:A:369:LEU:O	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:TRP:CZ2	1:A:181:TRP:CD1	3.09	0.41
1:A:177:TYR:O	1:A:179:ALA:C	2.59	0.41
1:A:275:VAL:HG13	1:A:369:LEU:HB2	2.03	0.41
1:A:441:ASP:OD2	1:A:445:GLN:NE2	2.54	0.41
1:A:478:ARG:HB2	1:A:553:VAL:HA	2.02	0.41
1:A:604:ASP:HA	6:A:1042:HOH:O	2.21	0.41
1:A:180:SER:O	1:A:181:TRP:C	2.60	0.40
1:A:202:LEU:CD2	1:A:218:ILE:HD11	2.44	0.40
1:A:203:PHE:CZ	1:A:217:HIS:CG	3.08	0.40
1:A:323:LEU:HD12	1:A:323:LEU:HA	1.90	0.40
1:A:411:ALA:O	1:A:414:VAL:CG2	2.68	0.40
1:A:151:ARG:NH2	1:A:505:GLU:OE1	2.54	0.40
1:A:451:ILE:O	1:A:454:LEU:HB2	2.20	0.40
1:A:496:TYR:OH	6:A:1003:HOH:O	2.16	0.40
1:A:151:ARG:HG2	1:A:151:ARG:NH1	2.36	0.40
1:A:644:LEU:HG	1:A:648:ILE:CD1	2.51	0.40
1:A:73:ASP:OD1	1:A:73:ASP:C	2.60	0.40
1:A:96:ILE:HG13	6:A:961:HOH:O	2.21	0.40

All (4) 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:A:175:GLN:OE1	1:A:260:ARG:O[5_555]	1.52	0.68
1:A:175:GLN:OE1	1:A:260:ARG:C[5_555]	1.58	0.62
1:A:175:GLN:OE1	1:A:261:LEU:N[5_555]	1.84	0.36
1:A:175:GLN:OE1	1:A:261:LEU:CA[5_555]	1.88	0.32

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	694/696 (100%)	656 (94%)	31 (4%)	7 (1%)	15 13

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	179	ALA
1	A	318	LYS
1	A	339	ASN
1	A	376	MET
1	A	260	ARG
1	A	168	VAL
1	A	319	PRO

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	605/605 (100%)	543 (90%)	62 (10%)	7 5

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	LYS
1	A	59	SER
1	A	64	SER
1	A	84	LYS
1	A	91	LEU
1	A	92	LYS
1	A	113	LEU
1	A	156	LEU
1	A	159	LEU
1	A	160	LEU
1	A	170	THR
1	A	185	LYS
1	A	189	GLN
1	A	202	LEU

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Mol	Chain	Res	Type
1	A	203	PHE
1	A	209	LYS
1	A	215	VAL
1	A	218	ILE
1	A	260	ARG
1	A	263	ILE
1	A	266	ASN
1	A	278	LEU
1	A	295	PRO
1	A	297	LEU
1	A	298	LEU
1	A	304	LEU
1	A	315	ILE
1	A	332	THR
1	A	339	ASN
1	A	340	GLU
1	A	349	GLU
1	A	354	LEU
1	A	359	THR
1	A	368	ASN
1	A	394	ARG
1	A	436	LYS
1	A	476	LYS
1	A	477	GLU
1	A	483	ASP
1	A	490	ASN
1	A	497	LEU
1	A	498	GLU
1	A	519	SER
1	A	520	LYS
1	A	522	LEU
1	A	523	LYS
1	A	551	GLN
1	A	597	ASN
1	A	598	LYS
1	A	612	SER
1	A	666	LYS
1	A	667	LYS
1	A	672	LYS
1	A	681	HIS
1	A	682	LYS
1	A	687	LEU

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Mol	Chain	Res	Type
1	A	718	ILE
1	A	729	SER
1	A	730	GLU
1	A	736	LYS
1	A	744	LYS
1	A	748	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	105	ASN
1	A	189	GLN
1	A	217	HIS
1	A	257	GLN
1	A	267	GLN
1	A	306	GLN
1	A	339	ASN
1	A	368	ASN
1	A	490	ASN
1	A	509	ASN
1	A	550	ASN
1	A	570	ASN
1	A	597	ASN
1	A	617	GLN
1	A	619	GLN
1	A	637	HIS
1	A	656	GLN
1	A	688	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 monosaccharides in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 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	NAG	A	753	1	14,14,15	0.62	0	17,19,21	1.12	2 (11%)
2	NAG	A	754	1	14,14,15	0.57	0	17,19,21	0.95	0
4	ACT	A	801	-	1,3,3	0.70	0	0,3,3	0.00	-
2	NAG	A	752	1	14,14,15	0.44	0	17,19,21	1.22	2 (11%)
5	STS	A	900	3	16,21,21	1.14	2 (12%)	19,28,28	1.04	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	753	1	-	1/6/23/26	0/1/1/1
2	NAG	A	754	1	-	0/6/23/26	0/1/1/1
5	STS	A	900	3	-	0/6/10/10	0/3/3/3
2	NAG	A	752	1	-	2/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	900	STS	C16-C19	2.16	1.39	1.36
5	A	900	STS	C16-C14	-2.16	1.38	1.41

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	753	NAG	C2-N2-C7	-2.89	118.79	122.90
2	A	752	NAG	C4-C3-C2	-2.88	106.80	111.02
2	A	752	NAG	O5-C1-C2	-2.52	107.30	111.29
2	A	753	NAG	O5-C5-C6	2.24	110.72	107.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	900	STS	C19-N18-C17	2.19	121.92	117.25

There are no chirality outliers.

All (3) torsion outliers are listed below:

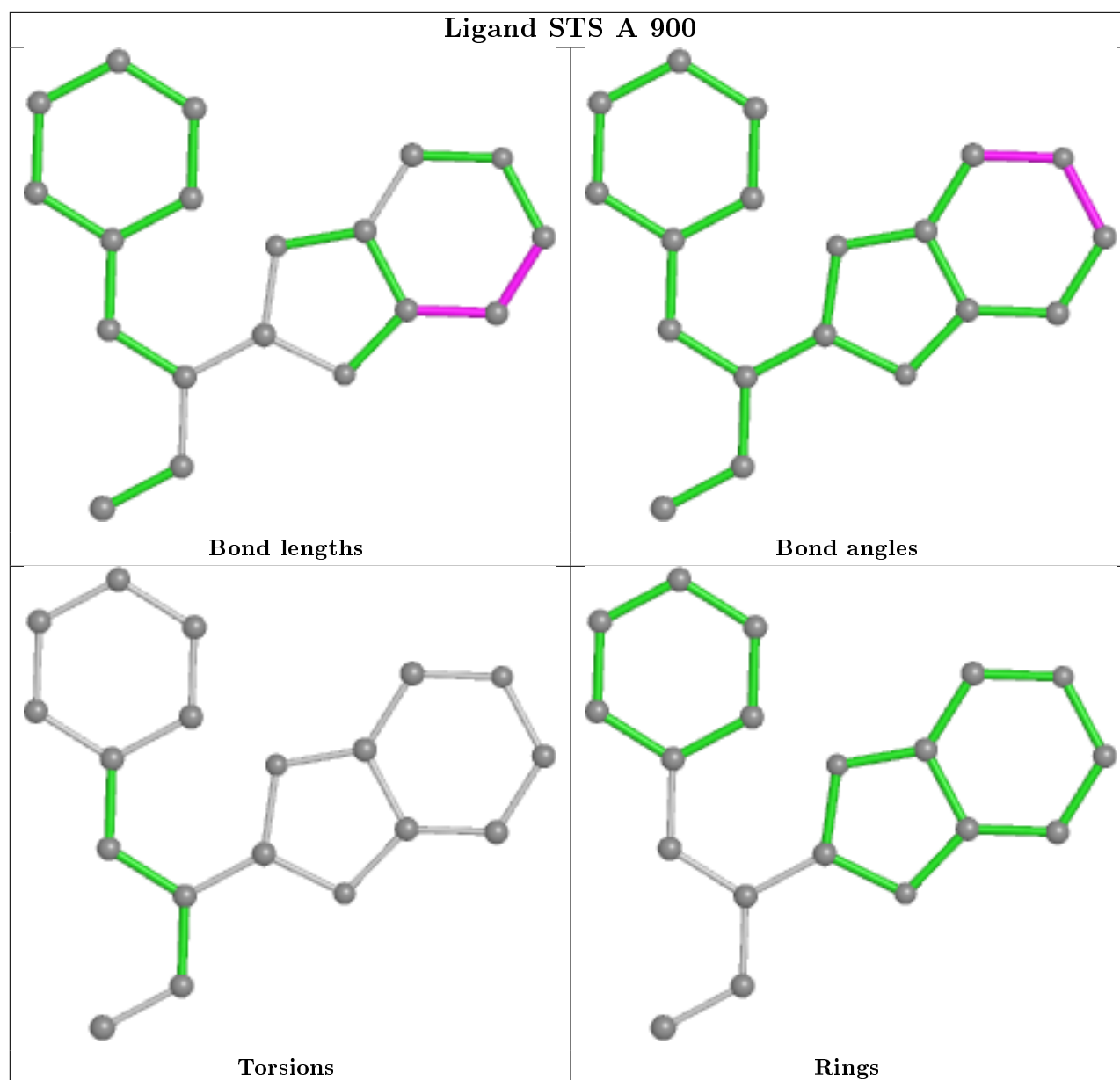
Mol	Chain	Res	Type	Atoms
2	A	752	NAG	O5-C5-C6-O6
2	A	752	NAG	C4-C5-C6-O6
2	A	753	NAG	O5-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	753	NAG	1	0
4	A	801	ACT	5	0
5	A	900	STS	4	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.