



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:36 AM UTC

PDB ID : 1USX / pdb_00001usx
Title : Crystal structure of the Newcastle disease virus hemagglutinin-neuraminidase complexed with thiosialoside
Authors : Zaitsev, V.; Itzstein, M.; Groves, D.; Kiefel, M.; Takimoto, T.; Portner, A.; Taylor, G.
Deposited on : 2003-12-01
Resolution : 2.70 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

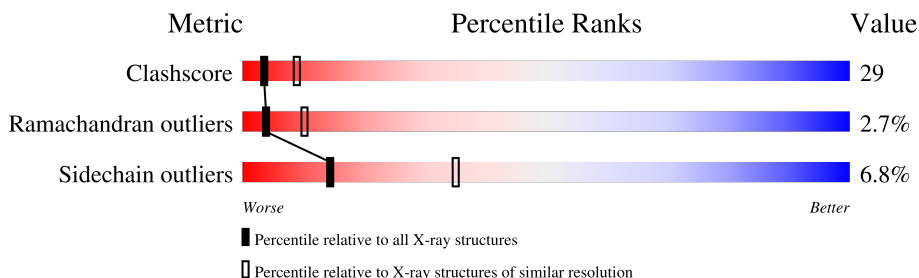
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)

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 $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	454	
1	B	454	
1	C	454	
2	D	2	
2	E	2	
2	F	2	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10497 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.

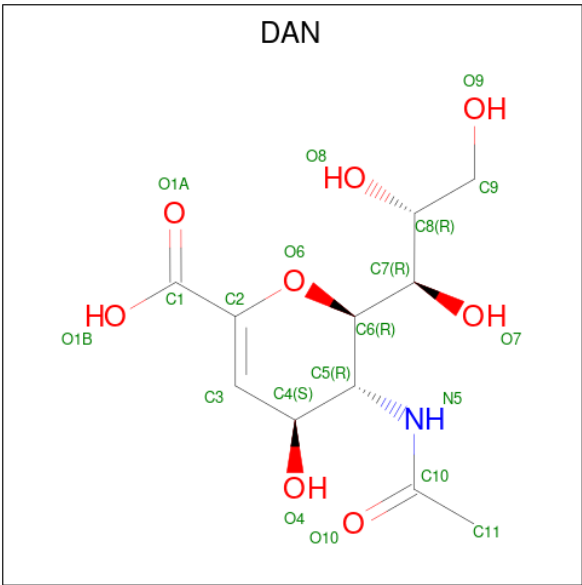
- Molecule 1 is a protein called HEMAGGLUTININ-NEURAMINIDASE GLYCOPROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	447	Total	C	N	O	S	0	0	1
			3446	2172	588	667	19			
1	B	447	Total	C	N	O	S	0	0	1
			3446	2172	588	667	19			
1	C	447	Total	C	N	O	S	0	0	1
			3446	2172	588	667	19			

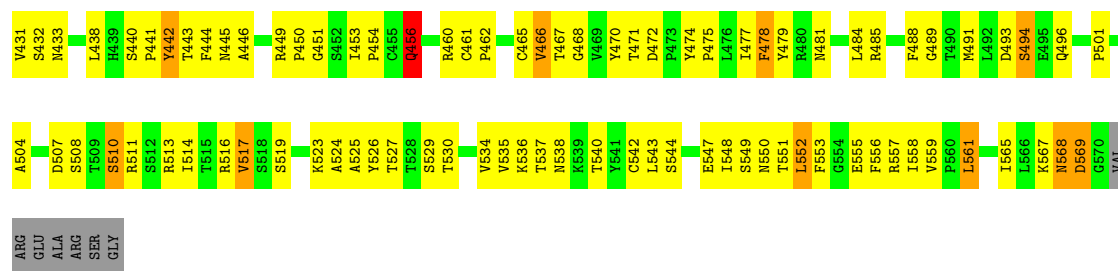
- Molecule 2 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-methyl 6-thio-beta-D-galactopyranoside.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	2	Total	C	N	O	S	0	0	0
			33	18	1	13	1			
2	E	2	Total	C	N	O	S	0	0	0
			33	18	1	13	1			
2	F	2	Total	C	N	O	S	0	0	0
			33	18	1	13	1			

- Molecule 3 is 2-DEOXY-2,3-DEHYDRO-N-ACETYL-NEURAMINIC ACID (CCD ID: DAN) (formula: C₁₁H₁₇NO₈).

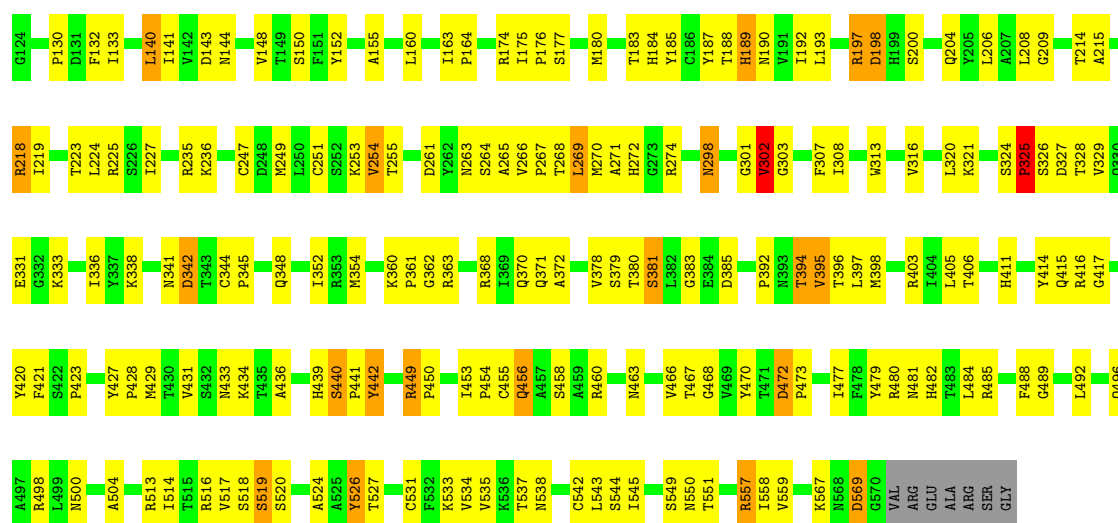


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			20	11	1	8		
3	B	1	Total	C	N	O	0	0
			20	11	1	8		
3	C	1	Total	C	N	O	0	0
			20	11	1	8		



• Molecule 1: HEMAGGLUTININ-NEURAMINIDASE GLYCOPROTEIN

Chain C: 57% 37% 5%



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	115.07Å 115.07Å 283.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	12.00 – 2.70	Depositor
% Data completeness (in resolution range)	79.6 (12.00-2.70)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.225 , 0.289	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	10497	wwPDB-VP
Average B, all atoms (Å ²)	30.0	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: DAN, WIA, SIA

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	0.48	1/3531 (0.0%)	0.99	15/4804 (0.3%)
1	B	0.50	1/3531 (0.0%)	1.02	17/4804 (0.4%)
1	C	0.51	1/3531 (0.0%)	1.05	19/4804 (0.4%)
All	All	0.50	3/10593 (0.0%)	1.02	51/14412 (0.4%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	569	ASP	C-N	-5.54	1.25	1.33
1	A	569	ASP	C-N	-5.43	1.25	1.33
1	C	569	ASP	C-N	-5.37	1.25	1.33

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	394	THR	N-CA-C	-8.24	103.22	113.28
1	A	442	TYR	N-CA-C	-7.65	96.76	109.07
1	C	254	VAL	N-CA-C	7.46	118.61	108.17
1	C	517	VAL	N-CA-C	-7.43	103.61	110.82
1	A	394	THR	N-CA-C	-7.38	103.73	112.89
1	C	519	SER	N-CA-C	-7.20	105.34	112.97
1	B	440	SER	CA-C-N	7.16	127.08	119.85
1	B	440	SER	C-N-CA	7.16	127.08	119.85
1	B	513	ARG	N-CA-C	-7.12	100.66	110.35
1	B	140	LEU	N-CA-C	7.08	118.89	111.03
1	C	513	ARG	N-CA-C	-7.03	101.22	110.43
1	C	140	LEU	N-CA-C	6.91	118.60	111.14
1	C	370	GLN	N-CA-C	6.57	119.64	109.07
1	B	472	ASP	N-CA-C	6.48	119.02	110.08
1	A	196	CYS	N-CA-C	6.43	119.98	111.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	263	ASN	N-CA-C	-6.42	103.97	113.61
1	A	514	ILE	N-CA-C	6.32	117.58	108.48
1	A	420	TYR	N-CA-C	6.30	118.14	111.28
1	B	255	THR	N-CA-C	-6.24	103.75	112.12
1	B	517	VAL	N-CA-C	-6.18	104.49	110.42
1	B	125	ALA	CA-C-N	6.13	127.50	119.84
1	B	125	ALA	C-N-CA	6.13	127.50	119.84
1	B	194	SER	N-CA-C	6.01	117.91	111.36
1	B	254	VAL	N-CA-C	6.01	118.51	109.20
1	A	559	VAL	N-CA-C	5.89	115.01	108.05
1	A	163	ILE	CA-C-N	5.88	125.88	119.89
1	A	163	ILE	C-N-CA	5.88	125.88	119.89
1	A	472	ASP	CA-C-N	5.81	127.10	119.84
1	A	472	ASP	C-N-CA	5.81	127.10	119.84
1	B	510	SER	N-CA-C	5.73	119.24	110.36
1	C	500	ASN	N-CA-C	5.68	122.37	109.81
1	C	155	ALA	N-CA-C	-5.56	103.04	110.55
1	A	472	ASP	N-CA-C	5.56	116.98	110.31
1	A	194	SER	N-CA-C	5.55	119.76	113.15
1	C	442	TYR	N-CA-C	-5.55	102.26	110.48
1	C	559	VAL	N-CA-C	5.54	114.59	108.05
1	A	193	LEU	N-CA-C	5.50	117.14	111.03
1	C	472	ASP	N-CA-C	5.50	118.63	110.39
1	B	371	GLN	N-CA-C	-5.45	103.70	110.41
1	C	193	LEU	N-CA-C	5.42	118.01	111.40
1	B	193	LEU	N-CA-C	5.40	120.23	111.37
1	B	442	TYR	N-CA-C	-5.37	101.72	110.20
1	A	152	TYR	N-CA-C	5.30	112.79	108.07
1	C	150	SER	N-CA-C	-5.26	106.86	113.28
1	A	348	GLN	N-CA-C	-5.25	105.67	111.71
1	C	255	THR	N-CA-C	-5.16	107.16	113.50
1	B	152	TYR	N-CA-C	5.10	114.16	108.25
1	C	440	SER	N-CA-C	-5.09	103.31	109.57
1	C	441	PRO	N-CA-C	5.08	119.30	111.57
1	B	441	PRO	N-CA-C	5.08	119.45	111.38
1	C	338	LYS	N-CA-C	5.05	117.97	110.14

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	3446	0	3342	192	0
1	B	3446	0	3342	270	0
1	C	3446	0	3342	163	0
2	D	33	0	27	2	0
2	E	33	0	24	0	0
2	F	33	0	24	1	0
3	A	20	0	16	0	0
3	B	20	0	16	1	0
3	C	20	0	16	1	0
All	All	10497	0	10149	608	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 29.

All (608) 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:219:ILE:HD11	1:A:562:LEU:HD21	1.43	1.00
1:B:456:GLN:H	1:B:456:GLN:NE2	1.58	1.00
1:A:550:ASN:HB2	1:A:557:ARG:HG2	1.44	0.99
1:A:500:ASN:HB3	1:A:516:ARG:HH22	1.29	0.96
1:C:268:THR:H	1:C:298:ASN:HD21	1.13	0.93
1:B:320:LEU:HD21	1:B:370:GLN:HB3	1.51	0.91
1:B:268:THR:H	1:B:298:ASN:HD21	1.20	0.88
1:B:320:LEU:HD11	1:B:370:GLN:HB2	1.55	0.87
1:B:456:GLN:H	1:B:456:GLN:HE21	1.20	0.87
1:B:301:GLY:O	1:B:302:VAL:HG22	1.75	0.86
1:B:253:LYS:HB2	1:B:269:LEU:HD12	1.58	0.86
1:B:203:HIS:CB	1:B:230:ASP:HB2	2.05	0.85
1:B:268:THR:N	1:B:298:ASN:HD21	1.75	0.85
1:B:392:PRO:HG2	1:B:395:VAL:HG12	1.59	0.84
1:B:550:ASN:HB2	1:B:557:ARG:HG2	1.61	0.83
1:A:447:PHE:CE1	1:A:491:MET:HB3	2.12	0.83
1:B:191:VAL:HB	1:B:203:HIS:CD2	2.13	0.83
1:A:453:ILE:HA	1:A:456:GLN:NE2	1.95	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:550:ASN:HA	1:A:557:ARG:NH1	1.95	0.81
1:C:371:GLN:HE21	1:C:415:GLN:HE21	1.27	0.81
1:C:354:MET:HE2	1:C:354:MET:HA	1.61	0.81
1:B:268:THR:H	1:B:298:ASN:ND2	1.79	0.80
1:B:252:SER:OG	1:B:300:PRO:HD3	1.82	0.79
1:B:256:GLU:HB2	1:B:260:GLU:CB	2.12	0.79
1:B:299:TYR:O	1:B:316:VAL:HG23	1.82	0.79
1:C:316:VAL:HG13	1:C:372:ALA:HB3	1.65	0.78
1:A:407:VAL:HG21	1:A:475:PRO:HB2	1.66	0.78
1:B:218:ARG:HH11	1:B:218:ARG:HG2	1.49	0.77
1:C:516:ARG:HG2	1:C:516:ARG:HH11	1.50	0.77
1:A:453:ILE:HG13	1:A:496:GLN:NE2	1.99	0.76
1:C:516:ARG:HE	1:C:519:SER:HA	1.50	0.76
1:A:371:GLN:HE21	1:A:399:GLY:HA2	1.51	0.75
1:A:342:ASP:OD2	1:A:460:ARG:HD3	1.85	0.75
1:C:268:THR:H	1:C:298:ASN:ND2	1.85	0.74
1:C:449:ARG:NH2	1:C:455:CYS:O	2.20	0.73
1:B:404:ILE:HG23	1:B:429:MET:HE1	1.68	0.73
1:B:218:ARG:HG2	1:B:218:ARG:NH1	2.02	0.73
1:A:205:TYR:C	1:A:206:LEU:HD23	2.13	0.72
1:C:132:PHE:CE1	1:C:534:VAL:HG22	2.24	0.72
1:B:456:GLN:HE21	1:B:456:GLN:N	1.86	0.72
1:A:456:GLN:CD	1:A:456:GLN:H	1.98	0.72
1:C:190:ASN:HD22	1:C:204:GLN:HE21	1.35	0.71
1:B:203:HIS:HB3	1:B:230:ASP:HB2	1.72	0.71
1:A:268:THR:H	1:A:298:ASN:HD21	1.38	0.71
1:C:190:ASN:HB2	1:C:204:GLN:HE22	1.54	0.71
1:B:371:GLN:NE2	1:B:415:GLN:HE21	1.89	0.71
1:A:371:GLN:HE22	1:A:400:ALA:H	1.38	0.70
1:B:320:LEU:CD1	1:B:368:ARG:HG2	2.21	0.70
1:A:550:ASN:CB	1:A:557:ARG:HG2	2.20	0.70
1:B:320:LEU:HG	1:B:368:ARG:O	1.92	0.70
1:A:190:ASN:HD22	1:A:204:GLN:HE21	1.39	0.70
1:B:453:ILE:HG23	1:B:454:PRO:HA	1.74	0.69
1:B:190:ASN:OD1	1:B:236:LYS:HE3	1.91	0.69
1:C:141:ILE:HB	1:C:477:ILE:HG12	1.75	0.69
1:C:453:ILE:HG23	1:C:454:PRO:HA	1.73	0.69
1:A:377:LYS:HD3	1:A:384:GLU:OE2	1.92	0.69
1:B:206:LEU:HD21	1:B:249:MET:SD	2.33	0.69
1:C:411:HIS:HB2	1:C:429:MET:O	1.93	0.69
1:B:256:GLU:HB2	1:B:260:GLU:HB3	1.73	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:467:THR:HG23	1:B:468:GLY:N	2.07	0.68
1:A:456:GLN:NE2	1:A:456:GLN:H	1.91	0.68
1:B:372:ALA:HB1	1:B:388:LEU:HD11	1.74	0.68
1:B:537:THR:HG22	1:C:537:THR:HG22	1.74	0.68
1:A:212:ARG:HB2	1:A:220:PHE:CZ	2.29	0.68
1:A:202:SER:OG	1:A:236:LYS:HE3	1.93	0.68
1:B:256:GLU:HB2	1:B:260:GLU:HB2	1.75	0.68
1:A:197:ARG:HD3	1:A:198:ASP:N	2.10	0.67
1:A:550:ASN:HB2	1:A:557:ARG:CG	2.21	0.67
1:B:320:LEU:HD12	1:B:368:ARG:HG2	1.76	0.67
1:A:411:HIS:HB2	1:A:429:MET:O	1.95	0.67
1:A:190:ASN:HD22	1:A:204:GLN:NE2	1.93	0.66
1:A:212:ARG:HB2	1:A:220:PHE:CE1	2.31	0.66
1:B:197:ARG:H	1:B:197:ARG:HD3	1.61	0.66
1:B:373:ILE:HD11	1:B:413:LEU:HD22	1.77	0.66
1:B:174:ARG:HH12	3:B:1572:DAN:H3	1.60	0.66
1:C:456:GLN:HG2	1:C:458:SER:H	1.61	0.65
1:A:479:TYR:CD1	1:A:485:ARG:HD2	2.31	0.65
1:A:231:ASP:HA	1:B:216:THR:HG23	1.79	0.64
1:B:467:THR:HG23	1:B:468:GLY:H	1.62	0.64
1:A:322:PRO:O	1:A:323:ASN:HB2	1.95	0.64
1:A:197:ARG:HD3	1:A:198:ASP:H	1.62	0.64
1:A:503:SER:O	1:A:504:ALA:HB2	1.97	0.64
1:A:136:ILE:HD12	1:A:180:MET:HE3	1.78	0.64
1:B:453:ILE:HG13	1:B:496:GLN:NE2	2.13	0.64
1:C:456:GLN:H	1:C:456:GLN:NE2	1.96	0.64
1:A:500:ASN:HB3	1:A:516:ARG:NH2	2.09	0.63
1:B:235:ARG:O	1:B:236:LYS:HG2	1.99	0.63
1:B:523:LYS:HB3	1:B:549:SER:HB3	1.81	0.63
1:B:413:LEU:HD23	1:B:438:LEU:HD11	1.80	0.63
1:A:453:ILE:HA	1:A:456:GLN:HE22	1.63	0.63
1:B:201:HIS:O	1:B:202:SER:HB3	1.99	0.63
1:A:301:GLY:C	1:A:302:VAL:HG13	2.25	0.62
1:A:297:ALA:C	1:A:298:ASN:HD22	2.08	0.62
1:B:132:PHE:CE1	1:B:534:VAL:HG22	2.35	0.62
1:B:203:HIS:HB2	1:B:230:ASP:HB2	1.82	0.61
1:B:189:HIS:ND1	1:B:558:ILE:HD12	2.14	0.61
1:B:291:LEU:HG	1:B:292:PHE:CE1	2.35	0.61
1:A:320:LEU:HD12	1:A:368:ARG:HG2	1.82	0.61
1:B:313:TRP:CZ2	1:B:431:VAL:HG21	2.35	0.61
1:B:504:ALA:HB2	1:B:514:ILE:HG22	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:206:LEU:HD21	1:C:249:MET:SD	2.41	0.61
1:C:219:ILE:C	1:C:219:ILE:HD12	2.26	0.61
1:B:535:VAL:HG23	1:C:215:ALA:O	2.00	0.60
1:B:370:GLN:HG3	1:B:371:GLN:O	2.01	0.60
1:C:185:TYR:OH	1:C:223:THR:HG22	2.02	0.60
1:B:451:GLY:O	1:B:496:GLN:HG2	2.01	0.60
1:C:224:LEU:O	1:C:225:ARG:HD3	2.02	0.60
1:B:525:ALA:HB3	1:B:547:GLU:HB2	1.84	0.60
1:C:264:SER:OG	1:C:264:SER:O	2.20	0.60
1:A:370:GLN:NE2	1:A:371:GLN:H	1.99	0.60
1:C:398:MET:HE1	1:C:466:VAL:HG22	1.82	0.60
1:A:136:ILE:HG23	1:A:180:MET:HE3	1.84	0.60
1:A:269:LEU:HD22	1:A:270:MET:H	1.67	0.60
1:A:268:THR:H	1:A:298:ASN:ND2	2.00	0.59
1:B:191:VAL:HB	1:B:203:HIS:HD2	1.61	0.59
1:B:371:GLN:HE22	1:B:400:ALA:H	1.48	0.59
1:B:416:ARG:HH21	1:B:468:GLY:C	2.11	0.59
1:C:481:ASN:O	1:C:482:HIS:HB2	2.02	0.59
1:A:249:MET:HE3	1:A:251:CYS:SG	2.43	0.59
1:B:346:ASP:HB3	1:B:350:TYR:HD2	1.68	0.59
1:B:449:ARG:HH12	1:B:451:GLY:HA3	1.67	0.58
1:B:396:THR:HA	1:B:427:TYR:HE2	1.69	0.58
1:B:516:ARG:HG2	1:B:516:ARG:HH11	1.69	0.58
1:C:144:ASN:OD1	1:C:480:ARG:HG3	2.04	0.58
1:A:555:GLU:HG2	1:A:556:PHE:CD2	2.39	0.58
1:B:316:VAL:HG22	1:B:317:TYR:N	2.18	0.58
1:B:396:THR:HA	1:B:427:TYR:CE2	2.39	0.58
1:C:253:LYS:HE3	1:C:271:ALA:HB2	1.84	0.58
1:C:342:ASP:OD2	1:C:460:ARG:HD3	2.04	0.58
1:B:491:MET:SD	1:B:511:ARG:NH1	2.76	0.58
1:B:516:ARG:HE	1:B:519:SER:HA	1.67	0.58
1:C:144:ASN:ND2	1:C:480:ARG:HG3	2.18	0.58
1:A:335:VAL:HG11	1:A:337:TYR:CE2	2.39	0.57
1:A:550:ASN:HA	1:A:557:ARG:HH12	1.68	0.57
1:B:386:PRO:O	1:B:387:VAL:HG23	2.05	0.57
1:A:447:PHE:CD1	1:A:491:MET:HB3	2.38	0.57
1:A:467:THR:HG23	1:A:468:GLY:N	2.20	0.57
1:B:404:ILE:CG2	1:B:429:MET:HE1	2.33	0.57
1:B:198:ASP:C	1:B:198:ASP:OD1	2.47	0.57
1:C:313:TRP:CZ2	1:C:431:VAL:HG21	2.40	0.57
1:B:456:GLN:NE2	1:B:456:GLN:N	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:518:SER:C	1:C:520:SER:H	2.12	0.57
1:C:160:LEU:C	1:C:160:LEU:HD23	2.30	0.57
1:A:479:TYR:CG	1:A:485:ARG:HD2	2.39	0.57
1:B:453:ILE:HG13	1:B:496:GLN:CD	2.29	0.57
1:A:193:LEU:CD1	1:B:218:ARG:HH21	2.17	0.57
1:A:132:PHE:CE2	1:A:211:LEU:HB3	2.40	0.57
1:B:392:PRO:O	1:B:395:VAL:HG12	2.05	0.57
1:A:148:VAL:HG13	1:A:477:ILE:HD12	1.86	0.56
1:A:371:GLN:NE2	1:A:400:ALA:H	2.03	0.56
1:B:225:ARG:HD2	1:B:279:GLY:CA	2.35	0.56
1:B:320:LEU:HD11	1:B:370:GLN:CB	2.31	0.56
1:B:334:TYR:O	1:B:394:THR:HA	2.03	0.56
1:C:197:ARG:NH1	1:C:198:ASP:HB2	2.20	0.56
1:B:345:PRO:O	1:B:346:ASP:C	2.48	0.56
1:C:144:ASN:CG	1:C:480:ARG:HG3	2.30	0.56
1:C:371:GLN:NE2	1:C:415:GLN:HE21	2.01	0.56
1:B:203:HIS:HB3	1:B:230:ASP:CB	2.36	0.56
1:B:360:LYS:HE3	1:B:368:ARG:CZ	2.34	0.56
1:A:268:THR:OG1	1:A:297:ALA:HA	2.05	0.56
1:B:192:ILE:HG12	1:B:202:SER:HB3	1.86	0.56
1:A:219:ILE:HD11	1:A:562:LEU:CD2	2.29	0.56
1:A:446:ALA:HB2	1:A:510:SER:HA	1.86	0.56
1:C:190:ASN:HD22	1:C:204:GLN:NE2	2.03	0.56
1:C:209:GLY:HA3	1:C:223:THR:HA	1.86	0.56
1:C:416:ARG:NH1	3:C:1572:DAN:H8	2.20	0.56
1:A:205:TYR:O	1:A:206:LEU:HD23	2.05	0.56
1:B:218:ARG:HH11	1:B:218:ARG:CG	2.18	0.56
1:B:388:LEU:HD12	1:B:389:THR:H	1.70	0.56
1:A:231:ASP:HA	1:B:216:THR:CG2	2.36	0.55
1:A:543:LEU:HD11	1:A:545:ILE:HG13	1.87	0.55
1:B:291:LEU:HG	1:B:292:PHE:CD1	2.41	0.55
1:B:418:SER:O	1:B:466:VAL:HA	2.06	0.55
1:B:191:VAL:O	1:B:202:SER:HB2	2.07	0.55
1:B:231:ASP:O	1:B:232:THR:C	2.49	0.55
1:C:249:MET:HE3	1:C:251:CYS:SG	2.46	0.55
1:C:479:TYR:CD1	1:C:485:ARG:HD2	2.42	0.55
1:C:567:LYS:HB3	1:C:569:ASP:O	2.06	0.55
1:B:270:MET:CE	1:B:300:PRO:HG3	2.37	0.55
1:C:268:THR:N	1:C:298:ASN:HD21	1.94	0.55
1:A:548:ILE:HG22	1:A:549:SER:N	2.22	0.55
1:B:201:HIS:CG	1:B:202:SER:H	2.24	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:326:SER:O	1:C:329:VAL:HG22	2.07	0.55
1:A:418:SER:HA	1:A:467:THR:O	2.07	0.55
1:A:549:SER:O	1:A:557:ARG:NH1	2.39	0.55
1:B:140:LEU:HD13	1:B:540:THR:HG23	1.88	0.55
1:B:360:LYS:HB3	1:B:366:GLY:O	2.05	0.54
1:B:442:TYR:CE1	1:B:484:LEU:HB3	2.42	0.54
1:C:403:ARG:HD3	1:C:405:LEU:HD11	1.89	0.54
1:A:447:PHE:HE1	1:A:491:MET:HB3	1.70	0.54
1:B:466:VAL:HG12	1:B:466:VAL:O	2.07	0.54
1:A:298:ASN:HD22	1:A:298:ASN:N	2.05	0.54
1:A:345:PRO:O	1:A:346:ASP:O	2.25	0.54
1:A:392:PRO:O	1:A:395:VAL:HG13	2.07	0.54
1:B:413:LEU:HD23	1:B:438:LEU:CD1	2.37	0.54
1:B:446:ALA:HB1	1:B:511:ARG:HB2	1.90	0.54
1:B:233:GLN:NE2	1:B:253:LYS:NZ	2.56	0.54
1:A:516:ARG:HH11	1:A:516:ARG:HG2	1.73	0.54
1:C:261:ASP:OD1	1:C:264:SER:OG	2.26	0.54
1:C:307:PHE:CE1	1:C:378:VAL:HG21	2.43	0.54
1:A:257:THR:OG1	1:A:260:GLU:HG3	2.07	0.53
1:B:297:ALA:C	1:B:298:ASN:HD22	2.16	0.53
1:C:192:ILE:HD11	1:C:236:LYS:NZ	2.23	0.53
1:C:328:THR:HA	1:C:331:GLU:HB2	1.89	0.53
1:A:193:LEU:HD11	1:B:218:ARG:HH21	1.73	0.53
1:B:190:ASN:HA	1:B:204:GLN:NE2	2.24	0.53
1:B:324:SER:O	1:B:328:THR:HG23	2.08	0.53
1:C:324:SER:O	1:C:325:PRO:C	2.50	0.53
1:A:234:ASN:HD22	1:A:254:VAL:HG23	1.73	0.53
1:C:385:ASP:OD2	1:C:434:LYS:NZ	2.35	0.53
1:A:298:ASN:ND2	1:A:298:ASN:N	2.55	0.53
1:B:199:HIS:CG	1:B:199:HIS:O	2.62	0.52
1:A:422:SER:OG	1:A:448:THR:HB	2.10	0.52
1:B:474:TYR:CD1	1:B:530:THR:HA	2.45	0.52
1:C:143:ASP:OD1	1:C:144:ASN:N	2.42	0.52
1:C:253:LYS:HD2	1:C:269:LEU:HD13	1.90	0.52
1:C:253:LYS:HE3	1:C:271:ALA:CB	2.40	0.52
1:C:417:GLY:O	1:C:449:ARG:NH1	2.42	0.52
1:C:479:TYR:CE1	1:C:485:ARG:HD2	2.44	0.52
1:C:504:ALA:HB2	1:C:514:ILE:HG22	1.90	0.52
1:C:516:ARG:NH2	1:C:518:SER:O	2.43	0.52
1:A:454:PRO:HA	1:A:456:GLN:HE22	1.75	0.52
1:C:453:ILE:CG2	1:C:454:PRO:HA	2.38	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:391:PRO:CG	1:B:396:THR:HG21	2.40	0.52
1:B:425:LEU:HD23	1:B:443:THR:HG22	1.92	0.52
1:B:551:THR:O	1:B:553:PHE:N	2.35	0.52
1:A:161:ASN:HD22	1:B:165:ALA:HB3	1.75	0.51
1:C:466:VAL:HG12	1:C:466:VAL:O	2.09	0.51
1:A:301:GLY:O	1:A:302:VAL:HG22	2.10	0.51
1:C:405:LEU:N	1:C:405:LEU:HD12	2.25	0.51
1:A:527:THR:HA	1:A:545:ILE:O	2.10	0.51
1:B:261:ASP:O	1:B:264:SER:OG	2.26	0.51
1:C:456:GLN:H	1:C:456:GLN:HE21	1.59	0.51
1:C:516:ARG:HG2	1:C:516:ARG:NH1	2.22	0.51
1:A:507:ASP:OD1	1:A:509:THR:N	2.33	0.51
1:A:143:ASP:CG	1:A:146:SER:HB2	2.36	0.51
1:A:152:TYR:HB2	1:A:153:PRO:HD2	1.93	0.51
1:A:453:ILE:HA	1:A:456:GLN:HE21	1.75	0.51
1:C:333:LYS:HD3	1:C:394:THR:CG2	2.41	0.51
1:B:479:TYR:HB2	1:B:481:ASN:OD1	2.11	0.50
1:B:547:GLU:O	1:B:548:ILE:HD13	2.11	0.50
1:B:371:GLN:HE21	1:B:399:GLY:HA2	1.76	0.50
1:C:325:PRO:O	1:C:329:VAL:HG13	2.11	0.50
1:B:147:ASP:O	1:B:148:VAL:C	2.54	0.50
1:B:371:GLN:NE2	1:B:400:ALA:H	2.09	0.50
1:A:371:GLN:NE2	1:A:399:GLY:HA2	2.23	0.50
1:C:174:ARG:C	1:C:175:ILE:HG12	2.36	0.50
1:A:150:SER:O	1:A:568:ASN:ND2	2.44	0.50
1:B:256:GLU:HG3	1:B:261:ASP:N	2.27	0.50
1:A:532:PHE:N	1:A:532:PHE:CD1	2.78	0.50
1:A:174:ARG:C	1:A:175:ILE:HG12	2.37	0.50
1:B:307:PHE:CZ	1:B:310:GLY:HA2	2.46	0.50
1:C:206:LEU:HB2	1:C:227:ILE:CG1	2.42	0.49
1:C:218:ARG:HB3	1:C:218:ARG:HH11	1.76	0.49
1:A:156:PHE:CE1	1:A:561:LEU:HB3	2.47	0.49
1:C:190:ASN:HB2	1:C:204:GLN:NE2	2.23	0.49
1:A:152:TYR:CD2	1:A:513:ARG:HD2	2.46	0.49
1:A:499:LEU:HD23	1:A:523:LYS:HA	1.93	0.49
1:B:233:GLN:NE2	1:B:253:LYS:HZ2	2.11	0.49
1:A:273:GLY:HA2	1:A:283:GLU:HA	1.94	0.49
1:A:538:ASN:O	1:A:538:ASN:CG	2.56	0.49
1:B:141:ILE:O	1:B:477:ILE:HA	2.12	0.49
1:B:555:GLU:HG2	1:B:556:PHE:CD2	2.47	0.49
1:B:268:THR:OG1	1:B:298:ASN:N	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:567:LYS:O	1:B:568:ASN:O	2.31	0.49
1:A:451:GLY:HA3	1:A:455:CYS:O	2.12	0.49
1:C:152:TYR:HE1	1:C:567:LYS:C	2.20	0.49
1:C:527:THR:HA	1:C:545:ILE:O	2.13	0.49
1:A:269:LEU:HD22	1:A:270:MET:N	2.28	0.49
1:A:476:LEU:HG	1:A:477:ILE:HG13	1.94	0.49
1:B:525:ALA:CB	1:B:547:GLU:OE1	2.61	0.49
1:B:209:GLY:HA3	1:B:223:THR:HA	1.94	0.49
1:B:357:SER:C	1:B:359:TYR:H	2.20	0.49
1:B:374:LEU:HD12	1:B:375:SER:N	2.28	0.49
1:A:295:TRP:N	1:A:295:TRP:CD1	2.80	0.48
1:B:235:ARG:HG3	1:B:235:ARG:HH11	1.77	0.48
1:C:144:ASN:HD21	1:C:480:ARG:HG3	1.78	0.48
1:A:523:LYS:HB3	1:A:549:SER:HB2	1.96	0.48
1:B:313:TRP:O	1:B:404:ILE:HG13	2.13	0.48
1:B:196:CYS:HB2	1:B:555:GLU:OE1	2.14	0.48
1:C:130:PRO:C	1:C:132:PHE:H	2.21	0.48
1:B:309:ASP:O	1:B:311:ARG:N	2.47	0.48
1:B:501:PRO:HB2	1:B:517:VAL:CG2	2.42	0.48
1:A:227:ILE:HA	1:B:220:PHE:CE2	2.48	0.48
1:A:548:ILE:HB	1:A:557:ARG:HD2	1.96	0.48
1:B:160:LEU:HD23	1:B:160:LEU:C	2.38	0.48
1:B:335:VAL:HG12	1:B:359:TYR:OH	2.13	0.48
1:B:507:ASP:OD1	1:B:510:SER:N	2.47	0.48
1:B:188:THR:HB	1:B:206:LEU:HD23	1.95	0.48
1:B:346:ASP:HB3	1:B:350:TYR:CD2	2.48	0.48
1:B:453:ILE:CG2	1:B:454:PRO:HA	2.42	0.48
1:C:261:ASP:HA	1:C:264:SER:HB3	1.95	0.48
1:C:264:SER:O	1:C:265:ALA:C	2.56	0.48
1:A:132:PHE:CE1	1:A:534:VAL:HG22	2.48	0.48
1:B:567:LYS:HG2	1:B:568:ASN:N	2.28	0.48
1:C:354:MET:HA	1:C:354:MET:CE	2.39	0.48
1:A:175:ILE:HD13	1:A:175:ILE:N	2.28	0.48
1:B:525:ALA:HB3	1:B:547:GLU:CB	2.44	0.48
1:A:432:SER:O	1:A:433:ASN:HB2	2.13	0.48
1:B:360:LYS:HE3	1:B:368:ARG:NH1	2.28	0.48
1:A:189:HIS:C	1:A:189:HIS:CD2	2.92	0.48
1:B:374:LEU:HD12	1:B:375:SER:H	1.78	0.48
1:C:190:ASN:CB	1:C:204:GLN:HE22	2.26	0.48
1:A:325:PRO:O	1:A:329:VAL:HG13	2.15	0.47
1:B:129:ASP:OD2	1:B:212:ARG:CB	2.62	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:298:ASN:ND2	1:B:298:ASN:N	2.62	0.47
1:B:350:TYR:CE1	1:B:354:MET:HE3	2.49	0.47
1:B:309:ASP:C	1:B:311:ARG:H	2.23	0.47
1:B:534:VAL:HG12	1:B:537:THR:H	1.79	0.47
1:C:183:THR:OG1	1:C:184:HIS:ND1	2.43	0.47
1:B:148:VAL:HG13	1:B:477:ILE:HD12	1.97	0.47
1:A:190:ASN:HB2	1:A:204:GLN:HE22	1.80	0.47
1:B:465:CYS:O	1:B:466:VAL:HG23	2.14	0.47
1:B:128:HIS:HA	1:B:210:VAL:HB	1.97	0.47
1:C:301:GLY:O	1:C:302:VAL:HG22	2.14	0.47
1:C:308:ILE:HG21	1:C:431:VAL:HG11	1.96	0.47
1:C:488:PHE:CG	1:C:489:GLY:N	2.83	0.47
1:A:250:LEU:HG	1:A:300:PRO:HG3	1.96	0.47
1:A:442:TYR:CE1	1:A:484:LEU:HB3	2.50	0.47
1:B:272:HIS:ND1	1:B:383:GLY:N	2.62	0.47
1:B:320:LEU:HD21	1:B:370:GLN:CB	2.34	0.47
1:B:449:ARG:NH1	1:B:451:GLY:HA3	2.28	0.47
1:B:467:THR:CG2	1:B:468:GLY:H	2.24	0.47
1:B:526:TYR:N	1:B:547:GLU:OE1	2.45	0.47
1:C:266:VAL:HB	1:C:267:PRO:HD2	1.97	0.47
1:C:313:TRP:CH2	1:C:431:VAL:CG2	2.98	0.47
1:B:190:ASN:HD22	1:B:204:GLN:HE21	1.61	0.47
1:B:268:THR:OG1	1:B:297:ALA:HA	2.15	0.47
1:B:132:PHE:CE2	1:B:211:LEU:HB3	2.50	0.47
1:C:141:ILE:O	1:C:477:ILE:HG23	2.15	0.47
1:C:320:LEU:HD12	1:C:368:ARG:HD2	1.97	0.47
1:A:158:GLU:HG2	1:B:193:LEU:HD23	1.96	0.47
1:A:161:ASN:ND2	1:B:165:ALA:HB3	2.30	0.47
1:B:173:THR:HA	1:B:190:ASN:O	2.15	0.47
1:C:130:PRO:C	1:C:132:PHE:N	2.72	0.47
1:C:398:MET:CE	1:C:466:VAL:HG22	2.44	0.47
1:C:431:VAL:HG13	1:C:431:VAL:O	2.14	0.47
1:C:498:ARG:HB3	1:C:524:ALA:O	2.14	0.47
1:B:398:MET:HE1	1:B:466:VAL:HG22	1.96	0.46
1:C:392:PRO:HG2	1:C:395:VAL:HG12	1.96	0.46
1:C:450:PRO:HD3	1:C:492:LEU:O	2.15	0.46
1:C:333:LYS:HD3	1:C:394:THR:HG23	1.96	0.46
1:A:181:SER:HB3	1:A:242:ALA:HB3	1.98	0.46
1:A:359:TYR:O	1:A:368:ARG:HA	2.15	0.46
1:A:234:ASN:HD22	1:A:254:VAL:CG2	2.28	0.46
1:A:352:ILE:O	1:A:356:LYS:HG2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:514:ILE:O	1:A:514:ILE:HG13	2.15	0.46
1:B:391:PRO:HG2	1:B:396:THR:HG21	1.97	0.46
1:C:549:SER:O	1:C:557:ARG:NH1	2.48	0.46
1:A:145:ALA:O	1:A:146:SER:C	2.58	0.46
1:A:264:SER:O	1:A:265:ALA:C	2.59	0.46
1:A:428:PRO:HB2	1:A:439:HIS:HB2	1.97	0.46
1:A:453:ILE:H	1:A:496:GLN:HE21	1.62	0.46
1:B:189:HIS:CD2	1:B:205:TYR:HB3	2.51	0.46
1:B:344:CYS:HB2	1:B:460:ARG:O	2.15	0.46
1:A:348:GLN:O	1:A:352:ILE:HG13	2.16	0.46
1:C:190:ASN:HA	1:C:204:GLN:NE2	2.31	0.46
1:B:478:PHE:CD1	1:B:478:PHE:N	2.83	0.46
1:B:537:THR:O	1:B:538:ASN:C	2.58	0.46
1:C:190:ASN:CB	1:C:204:GLN:NE2	2.79	0.46
1:C:152:TYR:CD1	1:C:152:TYR:N	2.84	0.46
1:A:471:THR:HG21	1:A:492:LEU:HD13	1.98	0.46
1:B:169:GLY:N	2:D:2:SIA:H111	2.31	0.46
1:B:219:ILE:O	1:B:219:ILE:HG13	2.14	0.46
1:B:317:TYR:CD2	1:B:400:ALA:HA	2.51	0.46
1:B:479:TYR:CE1	1:B:485:ARG:HD2	2.51	0.46
1:C:265:ALA:O	1:C:321:LYS:HE2	2.16	0.46
1:A:131:ASP:O	1:A:535:VAL:HG22	2.16	0.45
1:B:169:GLY:H	2:D:2:SIA:H111	1.81	0.45
1:C:362:GLY:O	1:C:363:ARG:C	2.60	0.45
1:A:261:ASP:OD1	1:A:264:SER:OG	2.28	0.45
1:A:452:SER:O	1:A:453:ILE:C	2.57	0.45
1:A:535:VAL:HG23	1:A:536:LYS:N	2.31	0.45
1:B:264:SER:O	1:B:265:ALA:C	2.57	0.45
1:B:516:ARG:HG2	1:B:516:ARG:NH1	2.31	0.45
1:C:308:ILE:HB	1:C:313:TRP:NE1	2.32	0.45
1:C:488:PHE:CD2	1:C:489:GLY:N	2.84	0.45
1:A:501:PRO:HB2	1:A:517:VAL:HG23	1.98	0.45
1:C:140:LEU:CD2	1:C:533:LYS:HB2	2.47	0.45
1:C:328:THR:HG22	1:C:331:GLU:OE1	2.17	0.45
1:C:354:MET:HE2	1:C:354:MET:CA	2.39	0.45
1:A:196:CYS:HB2	1:A:555:GLU:OE1	2.17	0.45
1:B:296:VAL:CG2	1:B:319:GLY:HA3	2.46	0.45
1:A:268:THR:N	1:A:298:ASN:HD21	2.08	0.45
1:B:212:ARG:HG3	1:B:220:PHE:CE2	2.52	0.45
1:C:537:THR:O	1:C:538:ASN:C	2.58	0.45
1:A:516:ARG:HG2	1:A:516:ARG:NH1	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:159:HIS:H	1:B:159:HIS:CD2	2.35	0.45
1:B:269:LEU:HD11	1:B:285:ASP:OD2	2.17	0.45
1:B:274:ARG:HB3	1:B:382:LEU:HD21	1.98	0.45
1:B:295:TRP:O	1:B:321:LYS:HE2	2.16	0.45
1:B:461:CYS:HA	1:B:462:PRO:HD3	1.87	0.45
1:C:414:TYR:CD2	1:C:472:ASP:HB3	2.52	0.45
1:C:453:ILE:HG13	1:C:496:GLN:NE2	2.32	0.45
1:A:185:TYR:CD1	1:A:185:TYR:C	2.95	0.45
1:B:371:GLN:HE21	1:B:415:GLN:HE21	1.63	0.45
1:B:493:ASP:O	1:B:494:SER:HB2	2.16	0.45
1:C:313:TRP:CH2	1:C:431:VAL:HG23	2.51	0.45
1:A:334:TYR:CD1	1:A:356:LYS:HD3	2.51	0.45
1:A:473:PRO:HB2	1:A:487:VAL:HG21	1.99	0.45
1:B:145:ALA:O	1:B:146:SER:O	2.34	0.45
1:B:363:ARG:HD3	1:B:466:VAL:HG11	1.99	0.45
1:C:133:ILE:HD13	1:C:133:ILE:N	2.32	0.45
1:A:531:CYS:HA	1:A:542:CYS:HA	1.98	0.44
1:B:268:THR:N	1:B:298:ASN:ND2	2.48	0.44
1:B:313:TRP:CD1	1:B:313:TRP:N	2.84	0.44
1:B:507:ASP:OD1	1:B:507:ASP:C	2.60	0.44
1:C:264:SER:O	1:C:266:VAL:O	2.35	0.44
1:C:313:TRP:CD1	1:C:313:TRP:N	2.84	0.44
1:A:203:HIS:HB2	1:A:229:LEU:O	2.17	0.44
1:B:139:GLU:HG2	1:B:478:PHE:CZ	2.52	0.44
1:B:316:VAL:CG2	1:B:317:TYR:N	2.80	0.44
1:B:334:TYR:CD1	1:B:356:LYS:HE3	2.52	0.44
1:B:488:PHE:CG	1:B:489:GLY:N	2.85	0.44
1:C:177:SER:HB3	1:C:188:THR:HG22	1.98	0.44
1:C:470:TYR:O	1:C:526:TYR:HA	2.17	0.44
1:A:426:LEU:HD23	1:A:426:LEU:HA	1.85	0.44
1:B:175:ILE:HD13	1:B:175:ILE:N	2.32	0.44
1:B:180:MET:HA	1:B:185:TYR:HA	1.98	0.44
1:C:208:LEU:HD23	1:C:208:LEU:C	2.42	0.44
1:B:561:LEU:N	1:B:561:LEU:HD23	2.32	0.44
1:C:442:TYR:CE1	1:C:484:LEU:HB3	2.53	0.44
1:A:153:PRO:HD3	1:A:504:ALA:HA	2.00	0.44
1:A:472:ASP:OD1	1:A:528:THR:HA	2.17	0.44
1:B:421:PHE:CE2	1:B:423:PRO:HB2	2.52	0.44
1:B:542:CYS:SG	1:B:565:ILE:HD11	2.58	0.44
1:A:491:MET:CE	1:A:493:ASP:HA	2.47	0.44
1:B:451:GLY:O	1:B:456:GLN:HA	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:301:GLY:C	1:C:302:VAL:HG13	2.42	0.44
1:A:301:GLY:O	1:A:302:VAL:HG13	2.17	0.44
1:B:262:TYR:O	1:B:367:LYS:NZ	2.50	0.44
1:C:307:PHE:HE1	1:C:378:VAL:HG21	1.80	0.44
2:F:1:WIA:H5	2:F:2:SIA:O6	2.18	0.44
1:A:140:LEU:HD13	1:A:540:THR:CG2	2.48	0.44
1:A:140:LEU:HD13	1:A:540:THR:HG23	2.00	0.44
1:A:304:GLY:O	1:A:403:ARG:HG3	2.18	0.44
1:B:203:HIS:CD2	1:B:203:HIS:H	2.35	0.44
1:B:320:LEU:CD2	1:B:370:GLN:HB3	2.36	0.44
1:B:174:ARG:C	1:B:175:ILE:HG12	2.43	0.43
1:B:320:LEU:N	1:B:320:LEU:HD23	2.33	0.43
1:B:529:SER:HA	1:B:543:LEU:O	2.18	0.43
1:C:189:HIS:C	1:C:189:HIS:CD2	2.96	0.43
1:A:148:VAL:HB	1:A:507:ASP:HB3	2.00	0.43
1:A:213:THR:HA	1:A:218:ARG:O	2.19	0.43
1:A:438:LEU:HD23	1:A:438:LEU:HA	1.86	0.43
1:B:413:LEU:CD2	1:B:438:LEU:HD11	2.48	0.43
1:C:209:GLY:CA	1:C:223:THR:HA	2.48	0.43
1:C:253:LYS:HB2	1:C:269:LEU:HB3	2.00	0.43
1:C:531:CYS:HA	1:C:542:CYS:HA	2.00	0.43
1:A:163:ILE:HA	1:A:164:PRO:HD3	1.86	0.43
1:B:250:LEU:HD21	1:B:374:LEU:HD23	1.99	0.43
1:B:516:ARG:NE	1:B:519:SER:HA	2.33	0.43
1:A:234:ASN:ND2	1:A:254:VAL:CG2	2.81	0.43
1:A:264:SER:O	1:A:264:SER:OG	2.36	0.43
1:A:536:LYS:C	1:A:538:ASN:H	2.25	0.43
1:A:543:LEU:HD12	1:A:544:SER:N	2.33	0.43
1:B:379:SER:C	1:B:381:SER:H	2.25	0.43
1:A:130:PRO:HA	1:A:133:ILE:HG12	2.00	0.43
1:B:301:GLY:C	1:B:303:GLY:H	2.26	0.43
1:A:209:GLY:HA3	1:A:223:THR:HA	2.00	0.43
1:B:474:TYR:HA	1:B:475:PRO:HD3	1.92	0.43
1:C:206:LEU:HD12	1:C:227:ILE:HD11	1.99	0.43
1:A:167:THR:OG1	1:B:158:GLU:HA	2.18	0.43
1:B:152:TYR:N	1:B:152:TYR:CD1	2.87	0.43
1:B:264:SER:OG	1:B:264:SER:O	2.37	0.43
1:B:282:HIS:O	1:B:382:LEU:HD13	2.18	0.43
1:B:392:PRO:C	1:B:394:THR:N	2.77	0.43
1:B:470:TYR:HB3	1:B:526:TYR:CD1	2.53	0.43
1:C:160:LEU:C	1:C:160:LEU:CD2	2.92	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:163:ILE:HA	1:C:164:PRO:HD3	1.81	0.43
1:A:173:THR:HG22	1:A:191:VAL:HG22	2.00	0.43
1:A:180:MET:HA	1:A:185:TYR:HA	2.01	0.43
1:A:423:PRO:HB3	1:A:445:ASN:HA	2.01	0.43
1:B:322:PRO:O	1:B:323:ASN:HB2	2.18	0.43
1:A:165:ALA:O	1:B:161:ASN:HB2	2.19	0.43
1:B:131:ASP:OD2	1:B:536:LYS:HG2	2.18	0.43
1:B:201:HIS:CG	1:B:202:SER:N	2.86	0.43
1:B:254:VAL:HG11	1:B:299:TYR:HE2	1.84	0.43
1:C:467:THR:OG1	1:C:468:GLY:N	2.48	0.43
1:B:413:LEU:HD12	1:B:414:TYR:N	2.34	0.42
1:C:235:ARG:NH1	1:C:235:ARG:HG3	2.33	0.42
1:C:550:ASN:OD1	1:C:557:ARG:CZ	2.67	0.42
1:B:139:GLU:HG2	1:B:478:PHE:CE2	2.54	0.42
1:B:197:ARG:H	1:B:197:ARG:CD	2.25	0.42
1:B:199:HIS:CE1	1:B:257:THR:HA	2.55	0.42
1:C:177:SER:HB3	1:C:188:THR:CG2	2.49	0.42
1:A:353:ARG:O	1:A:356:LYS:HB2	2.20	0.42
1:A:404:ILE:HD12	1:A:429:MET:SD	2.60	0.42
1:A:444:PHE:CZ	1:A:487:VAL:HG13	2.53	0.42
1:A:552:LEU:HB2	1:B:557:ARG:HH12	1.85	0.42
1:C:148:VAL:HG13	1:C:477:ILE:HD12	2.01	0.42
1:C:192:ILE:HD11	1:C:236:LYS:HZ1	1.84	0.42
1:A:135:GLY:HA3	1:A:533:LYS:O	2.20	0.42
1:A:165:ALA:H	1:B:161:ASN:HB3	1.85	0.42
1:B:290:THR:O	1:B:293:GLU:HB2	2.20	0.42
1:B:345:PRO:HG2	1:B:460:ARG:NH2	2.34	0.42
1:C:180:MET:HA	1:C:185:TYR:HA	2.01	0.42
1:C:327:ASP:OD1	1:C:368:ARG:NE	2.45	0.42
1:C:550:ASN:OD1	1:C:557:ARG:NE	2.52	0.42
1:C:550:ASN:CG	1:C:551:THR:H	2.26	0.42
1:A:337:TYR:HB3	1:A:421:PHE:CE1	2.55	0.42
1:B:444:PHE:CD1	1:B:444:PHE:N	2.88	0.42
1:C:176:PRO:HB3	1:C:558:ILE:HD13	2.02	0.42
1:C:371:GLN:HE21	1:C:415:GLN:NE2	2.05	0.42
1:A:151:PHE:HA	1:A:566:LEU:O	2.20	0.42
1:B:185:TYR:CD1	1:B:185:TYR:C	2.98	0.42
1:B:235:ARG:HG3	1:B:235:ARG:NH1	2.35	0.42
1:B:308:ILE:HD13	1:B:411:HIS:CE1	2.55	0.42
1:C:298:ASN:ND2	1:C:298:ASN:N	2.68	0.42
1:A:207:ALA:HA	1:A:225:ARG:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:PHE:CE2	1:B:227:ILE:HA	2.55	0.42
1:A:281:TYR:CD2	1:A:281:TYR:C	2.98	0.42
1:B:443:THR:O	1:B:508:SER:HA	2.20	0.42
1:B:525:ALA:HB1	1:B:547:GLU:OE1	2.20	0.42
1:B:548:ILE:HG22	1:B:549:SER:N	2.35	0.42
1:C:481:ASN:O	1:C:482:HIS:CB	2.67	0.42
1:A:161:ASN:HB2	1:B:165:ALA:O	2.20	0.42
1:B:189:HIS:CD2	1:B:189:HIS:C	2.98	0.42
1:B:237:SER:O	1:B:237:SER:OG	2.35	0.42
1:B:273:GLY:HA3	1:B:281:TYR:OH	2.19	0.42
1:B:427:TYR:O	1:B:428:PRO:C	2.63	0.42
1:B:169:GLY:HA2	1:B:552:LEU:O	2.20	0.41
1:C:360:LYS:CE	1:C:368:ARG:NH2	2.83	0.41
1:A:301:GLY:C	1:A:303:GLY:H	2.27	0.41
1:C:254:VAL:O	1:C:254:VAL:HG23	2.19	0.41
1:C:301:GLY:C	1:C:303:GLY:H	2.28	0.41
1:B:396:THR:HG22	1:B:438:LEU:HD12	2.02	0.41
1:C:348:GLN:O	1:C:352:ILE:HG13	2.19	0.41
1:C:361:PRO:HG3	1:C:466:VAL:CG2	2.50	0.41
1:C:379:SER:C	1:C:381:SER:H	2.29	0.41
1:C:439:HIS:O	1:C:440:SER:C	2.60	0.41
1:A:188:THR:CG2	1:A:240:VAL:HG13	2.50	0.41
1:B:357:SER:C	1:B:359:TYR:N	2.78	0.41
1:C:175:ILE:HD13	1:C:175:ILE:N	2.35	0.41
1:C:235:ARG:HG3	1:C:235:ARG:HH11	1.86	0.41
1:C:272:HIS:ND1	1:C:383:GLY:N	2.63	0.41
1:A:143:ASP:C	1:A:485:ARG:HH11	2.28	0.41
1:A:370:GLN:HE21	1:A:371:GLN:H	1.65	0.41
1:A:543:LEU:HD11	1:A:545:ILE:CG1	2.50	0.41
1:A:550:ASN:ND2	1:A:553:PHE:O	2.53	0.41
1:B:159:HIS:H	1:B:159:HIS:HD2	1.67	0.41
1:B:361:PRO:HG3	1:B:466:VAL:CG2	2.50	0.41
1:B:392:PRO:C	1:B:394:THR:H	2.29	0.41
1:C:406:THR:HG23	1:C:411:HIS:CE1	2.56	0.41
1:A:160:LEU:HD12	1:A:220:PHE:HD2	1.85	0.41
1:A:335:VAL:HG23	1:A:394:THR:O	2.20	0.41
1:B:307:PHE:CE1	1:B:378:VAL:HG21	2.55	0.41
1:B:398:MET:HE2	1:B:419:SER:CA	2.51	0.41
1:B:432:SER:O	1:B:433:ASN:HB2	2.20	0.41
1:B:523:LYS:HE2	1:B:556:PHE:CZ	2.56	0.41
1:A:169:GLY:HA2	1:A:552:LEU:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:216:THR:HG23	1:B:231:ASP:HA	2.02	0.41
1:A:413:LEU:HD22	1:A:438:LEU:HD11	2.02	0.41
1:A:484:LEU:HD12	1:A:485:ARG:N	2.34	0.41
1:B:201:HIS:O	1:B:202:SER:CB	2.67	0.41
1:B:389:THR:O	1:B:391:PRO:HD3	2.19	0.41
1:A:148:VAL:CG1	1:A:486:GLY:HA3	2.50	0.41
1:A:198:ASP:OD1	1:A:200:SER:N	2.48	0.41
1:A:470:TYR:CD1	1:A:470:TYR:C	2.99	0.41
1:C:187:TYR:CG	1:C:188:THR:N	2.89	0.41
1:C:247:CYS:O	1:C:274:ARG:HA	2.21	0.41
1:A:321:LYS:O	1:A:324:SER:OG	2.36	0.41
1:A:371:GLN:NE2	1:A:415:GLN:HE21	2.18	0.41
1:B:280:GLN:HE21	1:B:280:GLN:HA	1.86	0.41
1:B:475:PRO:HB3	1:B:484:LEU:HD13	2.02	0.41
1:B:501:PRO:HG3	1:B:524:ALA:CB	2.50	0.41
1:B:550:ASN:OD1	1:B:557:ARG:NH2	2.53	0.41
1:C:219:ILE:HG21	1:C:534:VAL:HG21	2.02	0.41
1:C:325:PRO:O	1:C:326:SER:C	2.63	0.41
1:C:421:PHE:CE1	1:C:423:PRO:HD2	2.56	0.41
1:C:429:MET:HE3	1:C:436:ALA:CB	2.50	0.41
1:C:543:LEU:HD12	1:C:544:SER:N	2.36	0.41
1:A:158:GLU:HB3	1:B:193:LEU:HD22	2.03	0.41
1:B:525:ALA:HB3	1:B:547:GLU:OE1	2.21	0.41
1:C:269:LEU:HD22	1:C:270:MET:N	2.36	0.41
1:C:344:CYS:HA	1:C:345:PRO:HD3	1.79	0.41
1:A:126:PRO:HB3	1:A:224:LEU:CD1	2.51	0.40
1:A:456:GLN:CD	1:A:456:GLN:N	2.73	0.40
1:A:470:TYR:HE1	1:A:528:THR:HB	1.86	0.40
1:A:513:ARG:HG2	1:A:514:ILE:N	2.36	0.40
1:B:149:THR:C	1:B:151:PHE:N	2.78	0.40
1:B:261:ASP:HA	1:B:264:SER:HB3	2.03	0.40
1:B:297:ALA:O	1:B:318:GLY:HA3	2.21	0.40
1:B:397:LEU:HD12	1:B:397:LEU:HA	1.91	0.40
1:C:336:ILE:HA	1:C:420:TYR:OH	2.21	0.40
1:C:427:TYR:O	1:C:428:PRO:C	2.64	0.40
1:C:526:TYR:N	1:C:526:TYR:CD1	2.89	0.40
1:B:249:MET:HE3	1:B:251:CYS:SG	2.61	0.40
1:B:364:PHE:HD2	1:B:369:ILE:HD11	1.87	0.40
1:B:416:ARG:NH2	1:B:468:GLY:O	2.54	0.40
1:B:449:ARG:HG2	1:B:450:PRO:O	2.21	0.40
1:B:568:ASN:HB3	1:B:569:ASP:H	1.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:256:GLU:HG3	1:A:261:ASP:HB2	2.03	0.40
1:A:299:TYR:CD1	1:A:299:TYR:N	2.89	0.40
1:A:444:PHE:CE2	1:A:487:VAL:HG13	2.56	0.40
1:A:449:ARG:O	1:A:449:ARG:HD3	2.21	0.40
1:A:534:VAL:O	1:A:535:VAL:C	2.62	0.40
1:B:178:PHE:CD1	1:B:178:PHE:C	2.98	0.40
1:B:309:ASP:C	1:B:311:ARG:N	2.76	0.40
1:C:518:SER:C	1:C:520:SER:N	2.78	0.40
1:C:449:ARG:O	1:C:449:ARG:HG2	2.21	0.40
1:C:516:ARG:NH1	1:C:516:ARG:CG	2.83	0.40
1:A:152:TYR:CD1	1:A:152:TYR:N	2.89	0.40
1:A:313:TRP:CD1	1:A:313:TRP:N	2.89	0.40
1:B:136:ILE:HG21	1:B:181:SER:O	2.22	0.40
1:B:151:PHE:CD1	1:B:567:LYS:HB2	2.56	0.40
1:B:342:ASP:OD2	1:B:460:ARG:HD3	2.22	0.40
1:C:214:THR:HG23	1:C:218:ARG:O	2.22	0.40

There are no symmetry-related clashes.

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	445/454 (98%)	391 (88%)	39 (9%)	15 (3%)	3	7
1	B	445/454 (98%)	376 (84%)	52 (12%)	17 (4%)	2	5
1	C	445/454 (98%)	394 (88%)	47 (11%)	4 (1%)	14	35
All	All	1335/1362 (98%)	1161 (87%)	138 (10%)	36 (3%)	4	10

All (36) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	200	SER

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Mol	Chain	Res	Type
1	A	302	VAL
1	A	346	ASP
1	B	146	SER
1	B	552	LEU
1	B	568	ASN
1	C	200	SER
1	A	304	GLY
1	A	342	ASP
1	A	393	ASN
1	A	473	PRO
1	B	148	VAL
1	B	202	SER
1	B	346	ASP
1	B	347	GLU
1	B	466	VAL
1	B	494	SER
1	A	345	PRO
1	A	504	ALA
1	A	536	LYS
1	B	145	ALA
1	B	232	THR
1	B	310	GLY
1	B	456	GLN
1	A	525	ALA
1	B	302	VAL
1	B	345	PRO
1	B	366	GLY
1	A	433	ASN
1	A	500	ASN
1	B	200	SER
1	C	473	PRO
1	A	459	ALA
1	A	475	PRO
1	C	302	VAL
1	C	325	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	387/392 (99%)	353 (91%)	34 (9%)	9	23
1	B	387/392 (99%)	364 (94%)	23 (6%)	18	42
1	C	387/392 (99%)	365 (94%)	22 (6%)	18	43
All	All	1161/1176 (99%)	1082 (93%)	79 (7%)	14	35

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	131	ASP
1	A	143	ASP
1	A	149	THR
1	A	160	LEU
1	A	170	SER
1	A	175	ILE
1	A	189	HIS
1	A	197	ARG
1	A	206	LEU
1	A	218	ARG
1	A	223	THR
1	A	266	VAL
1	A	269	LEU
1	A	298	ASN
1	A	302	VAL
1	A	315	SER
1	A	316	VAL
1	A	335	VAL
1	A	342	ASP
1	A	381	SER
1	A	395	VAL
1	A	396	THR
1	A	397	LEU
1	A	449	ARG
1	A	454	PRO
1	A	456	GLN
1	A	464	SER
1	A	473	PRO
1	A	500	ASN
1	A	518	SER
1	A	532	PHE
1	A	543	LEU

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Mol	Chain	Res	Type
1	A	557	ARG
1	A	559	VAL
1	B	143	ASP
1	B	149	THR
1	B	175	ILE
1	B	197	ARG
1	B	198	ASP
1	B	203	HIS
1	B	218	ARG
1	B	255	THR
1	B	256	GLU
1	B	266	VAL
1	B	298	ASN
1	B	320	LEU
1	B	342	ASP
1	B	380	THR
1	B	419	SER
1	B	445	ASN
1	B	456	GLN
1	B	471	THR
1	B	478	PHE
1	B	527	THR
1	B	544	SER
1	B	559	VAL
1	B	561	LEU
1	C	189	HIS
1	C	197	ARG
1	C	198	ASP
1	C	218	ARG
1	C	269	LEU
1	C	298	ASN
1	C	302	VAL
1	C	325	PRO
1	C	341	ASN
1	C	342	ASP
1	C	380	THR
1	C	381	SER
1	C	395	VAL
1	C	396	THR
1	C	397	LEU
1	C	433	ASN
1	C	449	ARG

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Mol	Chain	Res	Type
1	C	456	GLN
1	C	463	ASN
1	C	526	TYR
1	C	535	VAL
1	C	557	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	157	GLN
1	A	190	ASN
1	A	204	GLN
1	A	234	ASN
1	A	263	ASN
1	A	298	ASN
1	A	370	GLN
1	A	371	GLN
1	A	393	ASN
1	A	411	HIS
1	A	433	ASN
1	A	456	GLN
1	A	496	GLN
1	A	538	ASN
1	A	568	ASN
1	B	199	HIS
1	B	204	GLN
1	B	233	GLN
1	B	280	GLN
1	B	298	ASN
1	B	351	GLN
1	B	371	GLN
1	B	411	HIS
1	B	456	GLN
1	B	463	ASN
1	C	190	ASN
1	C	203	HIS
1	C	204	GLN
1	C	282	HIS
1	C	298	ASN
1	C	371	GLN
1	C	456	GLN
1	C	463	ASN

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Mol	Chain	Res	Type
1	C	496	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

6 monosaccharides 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	WIA	D	1	2	13,13,13	3.87	5 (38%)	17,18,18	1.66	2 (11%)
2	SIA	D	2	2	20,20,21	5.30	15 (75%)	21,28,31	2.51	9 (42%)
2	WIA	E	1	2	13,13,13	3.73	8 (61%)	17,18,18	1.26	2 (11%)
2	SIA	E	2	2	20,20,21	5.39	14 (70%)	21,28,31	3.01	10 (47%)
2	WIA	F	1	2	13,13,13	3.86	4 (30%)	17,18,18	1.53	1 (5%)
2	SIA	F	2	2	20,20,21	5.41	13 (65%)	21,28,31	2.63	11 (52%)

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	WIA	D	1	2	-	1/4/24/24	0/1/1/1
2	SIA	D	2	2	-	2/18/34/38	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	WIA	E	1	2	-	4/4/24/24	0/1/1/1
2	SIA	E	2	2	-	6/18/34/38	0/1/1/1
2	WIA	F	1	2	-	1/4/24/24	0/1/1/1
2	SIA	F	2	2	-	4/18/34/38	0/1/1/1

All (59) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	2	SIA	O8-C8	-16.63	1.08	1.43
2	D	2	SIA	O8-C8	-16.15	1.09	1.43
2	F	2	SIA	O8-C8	-15.88	1.10	1.43
2	F	1	WIA	C6-S6	-12.30	1.56	1.81
2	D	1	WIA	C6-S6	-11.90	1.56	1.81
2	E	2	SIA	C11-C10	-11.88	1.25	1.50
2	F	2	SIA	C11-C10	-11.84	1.26	1.50
2	D	2	SIA	C11-C10	-11.05	1.27	1.50
2	E	1	WIA	C6-S6	-10.21	1.60	1.81
2	F	2	SIA	O1B-C1	-6.72	1.09	1.30
2	D	2	SIA	O1B-C1	-6.57	1.09	1.30
2	D	2	SIA	C5-N5	-6.19	1.36	1.45
2	F	2	SIA	C5-N5	-5.96	1.36	1.45
2	E	2	SIA	C4-C5	5.39	1.58	1.53
2	E	2	SIA	C5-N5	-4.57	1.38	1.45
2	D	2	SIA	C4-C5	4.30	1.57	1.53
2	F	2	SIA	C2-C1	4.30	1.57	1.52
2	E	2	SIA	C2-C1	4.29	1.57	1.52
2	F	2	SIA	C8-C7	4.28	1.60	1.53
2	E	1	WIA	O1-C1	4.23	1.47	1.40
2	E	2	SIA	O1A-C1	-4.13	1.09	1.22
2	F	2	SIA	O6-C6	4.04	1.50	1.44
2	D	2	SIA	C2-C1	3.78	1.56	1.52
2	E	2	SIA	O6-C6	3.77	1.49	1.44
2	D	1	WIA	O1-C1	3.75	1.46	1.40
2	E	2	SIA	C7-C6	-3.68	1.48	1.52
2	E	2	SIA	C3-C2	3.52	1.58	1.52
2	F	2	SIA	C3-C2	3.46	1.58	1.52
2	E	1	WIA	O5-C1	3.43	1.50	1.41
2	F	1	WIA	O1-C1	3.33	1.45	1.40
2	D	2	SIA	O6-C6	3.29	1.49	1.44
2	F	2	SIA	C6-C5	3.24	1.58	1.53
2	E	1	WIA	C4-C5	3.22	1.59	1.53
2	E	2	SIA	C6-C5	3.05	1.57	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2	SIA	C8-C7	2.98	1.58	1.53
2	F	1	WIA	C4-C5	2.87	1.59	1.53
2	D	1	WIA	C4-C5	2.85	1.59	1.53
2	D	2	SIA	C3-C2	2.80	1.56	1.52
2	F	2	SIA	C4-C5	2.75	1.55	1.53
2	D	2	SIA	O6-C2	2.75	1.48	1.43
2	E	1	WIA	C6-C5	2.74	1.59	1.51
2	D	1	WIA	O5-C1	2.73	1.48	1.41
2	F	2	SIA	O6-C2	2.64	1.48	1.43
2	E	2	SIA	C9-C8	2.56	1.58	1.52
2	E	2	SIA	C8-C7	2.55	1.57	1.53
2	E	2	SIA	O6-C2	2.45	1.48	1.43
2	E	1	WIA	O5-C5	2.43	1.50	1.44
2	D	2	SIA	C6-C5	2.38	1.56	1.53
2	E	1	WIA	C4-C3	2.36	1.58	1.52
2	D	2	SIA	C9-C8	2.34	1.58	1.52
2	D	2	SIA	C7-C6	-2.30	1.50	1.52
2	F	1	WIA	O5-C1	2.25	1.47	1.41
2	D	2	SIA	O10-C10	2.20	1.28	1.23
2	F	2	SIA	C7-C6	-2.20	1.50	1.52
2	F	2	SIA	C9-C8	2.18	1.57	1.52
2	E	1	WIA	O3-C3	2.13	1.48	1.43
2	D	2	SIA	O1A-C1	2.09	1.28	1.22
2	D	1	WIA	C4-C3	2.03	1.57	1.52
2	E	2	SIA	C3-C4	2.00	1.56	1.52

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1	WIA	O1-C1-C2	5.73	114.75	108.14
2	F	2	SIA	C9-C8-C7	-5.58	100.79	112.17
2	D	2	SIA	C9-C8-C7	-5.55	100.85	112.17
2	E	2	SIA	C8-C7-C6	-5.35	102.99	113.05
2	F	1	WIA	O1-C1-C2	5.29	114.24	108.14
2	E	2	SIA	O1A-C1-C2	-5.11	111.81	122.85
2	E	2	SIA	O7-C7-C8	4.99	120.27	108.93
2	E	2	SIA	C9-C8-C7	-4.81	102.36	112.17
2	F	2	SIA	O10-C10-C11	-4.59	113.89	122.05
2	E	2	SIA	O10-C10-C11	-4.37	114.28	122.05
2	F	2	SIA	O7-C7-C8	4.29	118.67	108.93
2	D	2	SIA	O7-C7-C8	4.04	118.12	108.93
2	D	2	SIA	O10-C10-C11	-4.03	114.87	122.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	2	SIA	O10-C10-N5	3.64	128.42	121.98
2	F	2	SIA	O7-C7-C6	-3.57	101.72	109.44
2	E	2	SIA	O1B-C1-C2	3.56	121.98	112.71
2	E	1	WIA	O1-C1-C2	3.56	112.24	108.14
2	D	2	SIA	O7-C7-C6	-3.54	101.79	109.44
2	E	2	SIA	O7-C7-C6	-3.52	101.83	109.44
2	E	2	SIA	O8-C8-C9	3.42	116.81	109.03
2	D	2	SIA	O8-C8-C9	3.38	116.72	109.03
2	E	2	SIA	O10-C10-N5	3.30	127.81	121.98
2	D	2	SIA	O10-C10-N5	3.25	127.72	121.98
2	E	2	SIA	C5-N5-C10	3.19	130.57	123.11
2	D	2	SIA	C8-C7-C6	-3.14	107.15	113.05
2	F	2	SIA	O8-C8-C9	3.09	116.06	109.03
2	F	2	SIA	C8-C7-C6	-3.04	107.34	113.05
2	D	2	SIA	C5-N5-C10	2.91	129.92	123.11
2	F	2	SIA	C5-N5-C10	2.83	129.74	123.11
2	E	1	WIA	O5-C5-C6	2.67	110.96	106.49
2	F	2	SIA	C6-C5-N5	2.55	114.97	110.91
2	D	1	WIA	O5-C5-C6	2.50	110.67	106.49
2	F	2	SIA	O6-C2-C1	2.42	112.29	107.72
2	D	2	SIA	O6-C2-C1	2.40	112.25	107.72
2	F	2	SIA	O6-C2-C3	-2.06	107.79	110.56

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	2	SIA	C7-C8-C9-O9
2	D	2	SIA	O8-C8-C9-O9
2	E	1	WIA	C4-C5-C6-S6
2	E	1	WIA	O5-C5-C6-S6
2	E	2	SIA	O8-C8-C9-O9
2	F	1	WIA	O5-C5-C6-S6
2	E	1	WIA	O5-C1-O1-CM
2	E	2	SIA	C7-C8-C9-O9
2	E	1	WIA	C2-C1-O1-CM
2	E	2	SIA	C6-C7-C8-C9
2	F	2	SIA	O8-C8-C9-O9
2	E	2	SIA	C6-C7-C8-O8
2	F	2	SIA	C7-C8-C9-O9
2	F	2	SIA	O7-C7-C8-O8
2	F	2	SIA	O7-C7-C8-C9

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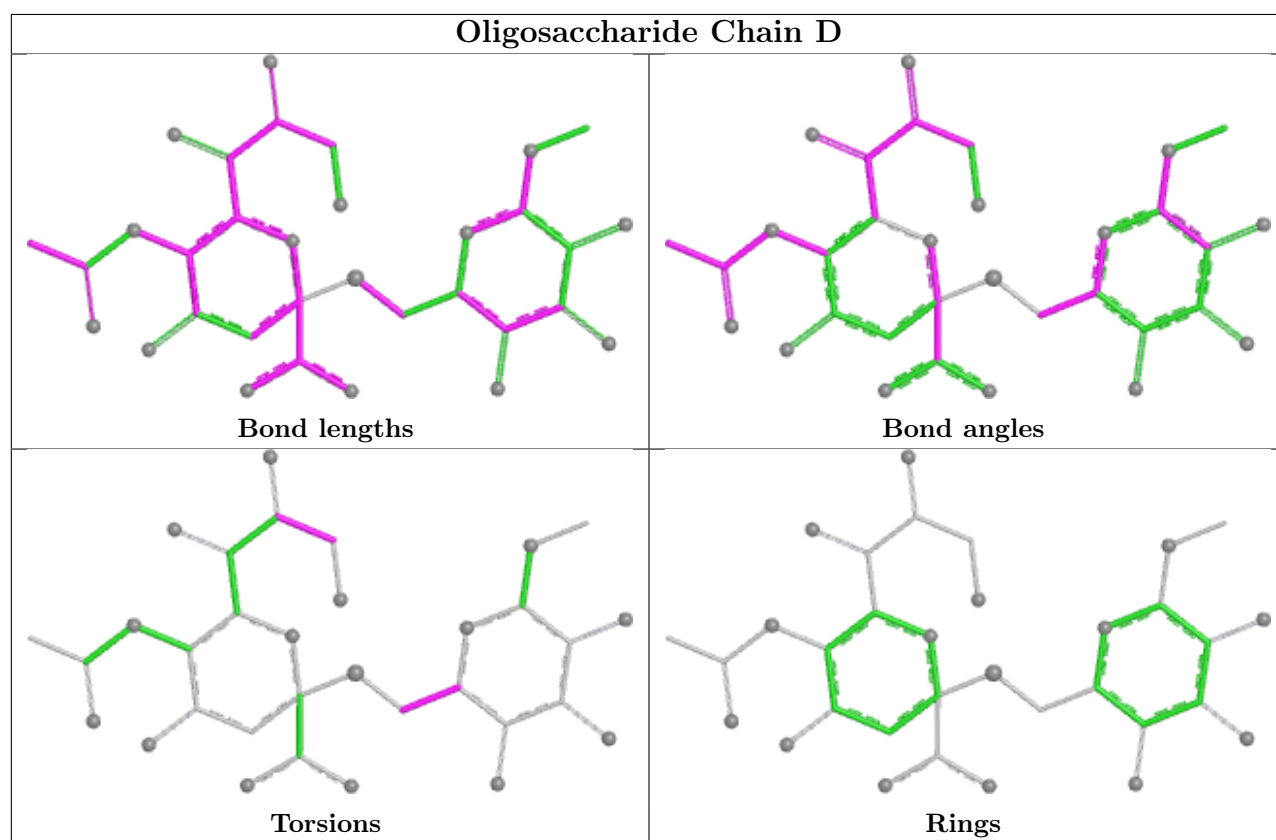
Mol	Chain	Res	Type	Atoms
2	D	1	WIA	O5-C5-C6-S6
2	E	2	SIA	C4-C5-N5-C10
2	E	2	SIA	O7-C7-C8-C9

There are no ring outliers.

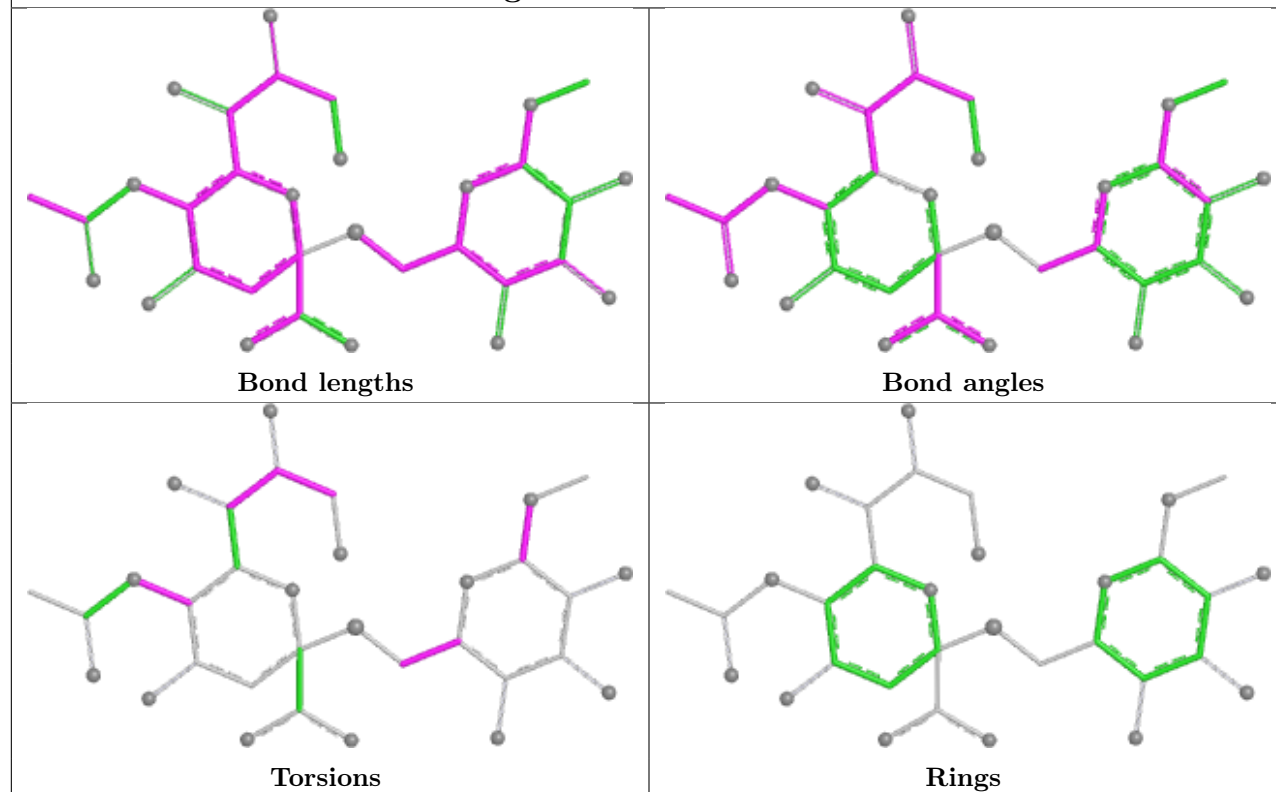
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	2	SIA	1	0
2	F	1	WIA	1	0
2	D	2	SIA	2	0

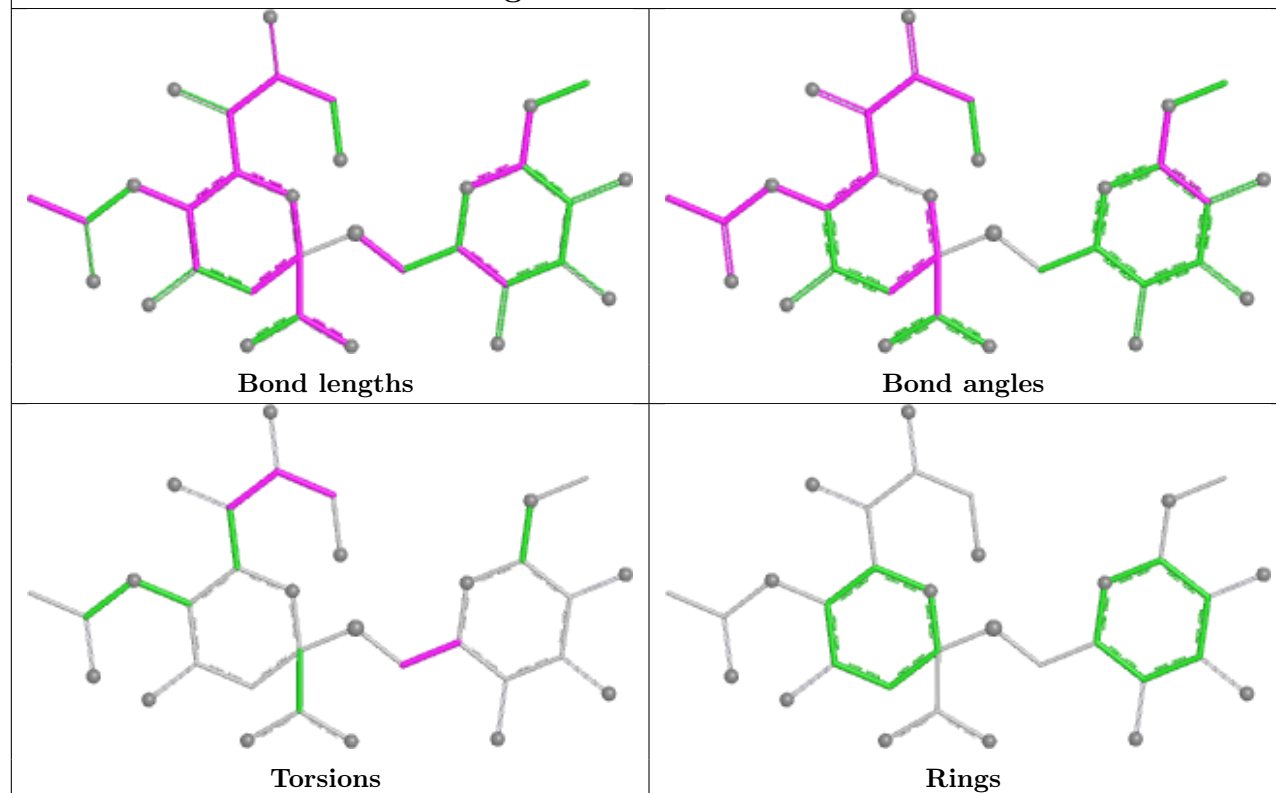
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



Oligosaccharide Chain E



Oligosaccharide Chain F



5.6 Ligand geometry

3 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
3	DAN	B	1572	-	20,20,20	2.00	6 (30%)	24,28,28	1.94	5 (20%)
3	DAN	C	1572	-	20,20,20	1.58	4 (20%)	24,28,28	2.54	5 (20%)
3	DAN	A	1570	-	20,20,20	1.58	4 (20%)	24,28,28	2.53	5 (20%)

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
3	DAN	B	1572	-	-	1/18/34/34	0/1/1/1
3	DAN	C	1572	-	-	0/18/34/34	0/1/1/1
3	DAN	A	1570	-	-	0/18/34/34	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1572	DAN	C2-C1	-6.41	1.34	1.48
3	A	1570	DAN	C3-C2	3.89	1.40	1.33
3	C	1572	DAN	C3-C2	3.89	1.40	1.33
3	C	1572	DAN	C2-C1	-3.44	1.41	1.48
3	A	1570	DAN	C2-C1	-3.44	1.41	1.48
3	B	1572	DAN	O8-C8	3.15	1.49	1.43
3	B	1572	DAN	O6-C6	-2.94	1.41	1.46
3	B	1572	DAN	O1B-C1	-2.24	1.24	1.30
3	A	1570	DAN	O8-C8	2.12	1.47	1.43
3	C	1572	DAN	O8-C8	2.11	1.47	1.43
3	A	1570	DAN	O1B-C1	-2.03	1.25	1.30
3	B	1572	DAN	O4-C4	-2.03	1.39	1.43
3	C	1572	DAN	O1B-C1	-2.00	1.25	1.30
3	B	1572	DAN	C3-C2	2.00	1.36	1.33

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1572	DAN	O1B-C1-C2	9.13	134.75	114.06
3	A	1570	DAN	O1B-C1-C2	9.12	134.72	114.06
3	A	1570	DAN	O1A-C1-C2	-5.27	107.67	120.57
3	C	1572	DAN	O1A-C1-C2	-5.25	107.70	120.57
3	B	1572	DAN	C4-C3-C2	-4.89	113.18	121.45
3	B	1572	DAN	C3-C2-C1	-4.25	114.36	123.56
3	C	1572	DAN	O4-C4-C3	-4.03	100.16	109.27
3	A	1570	DAN	O4-C4-C3	-4.03	100.17	109.27
3	B	1572	DAN	C8-C7-C6	-3.12	107.19	113.05
3	B	1572	DAN	O6-C2-C3	2.99	127.96	124.62
3	C	1572	DAN	O1B-C1-O1A	-2.75	117.35	123.90
3	A	1570	DAN	O1B-C1-O1A	-2.73	117.41	123.90
3	B	1572	DAN	O6-C2-C1	2.36	116.72	112.02
3	C	1572	DAN	C4-C3-C2	-2.03	118.03	121.45
3	A	1570	DAN	C4-C3-C2	-2.00	118.06	121.45

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1572	DAN	O8-C8-C9-O9

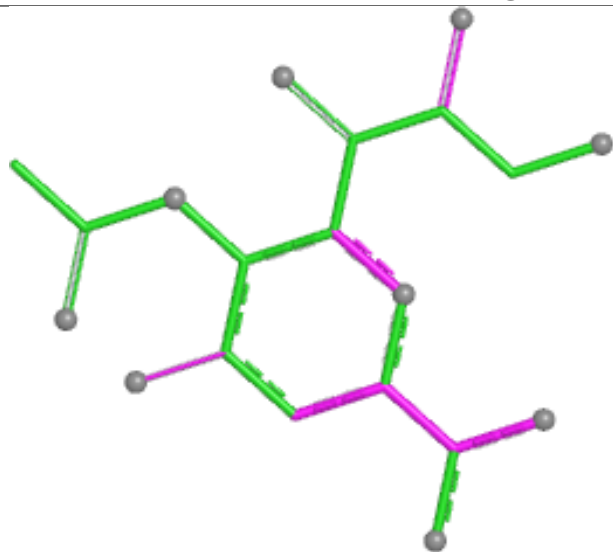
There are no ring outliers.

2 monomers are involved in 2 short contacts:

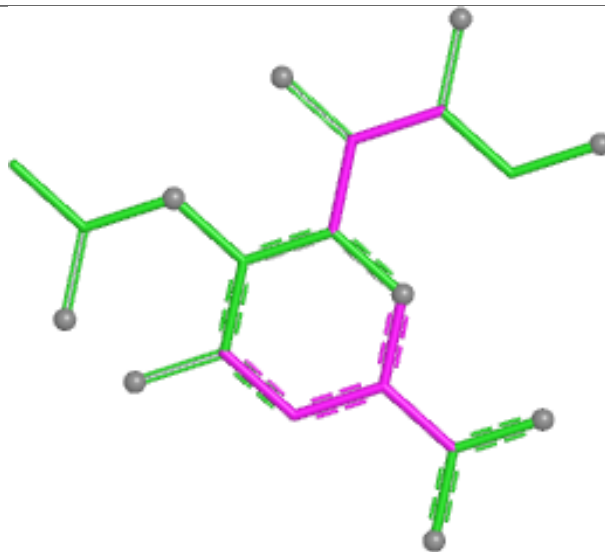
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1572	DAN	1	0
3	C	1572	DAN	1	0

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.

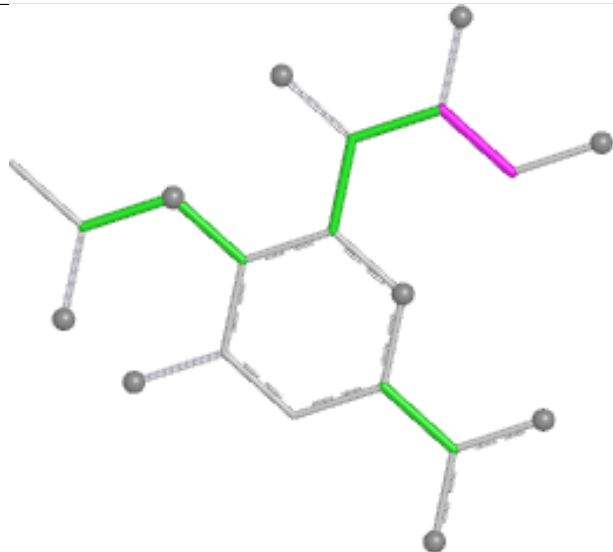
Ligand DAN B 1572



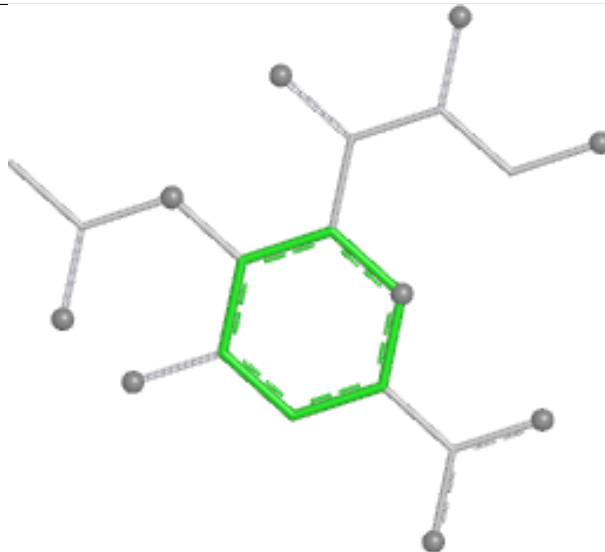
Bond lengths



Bond angles

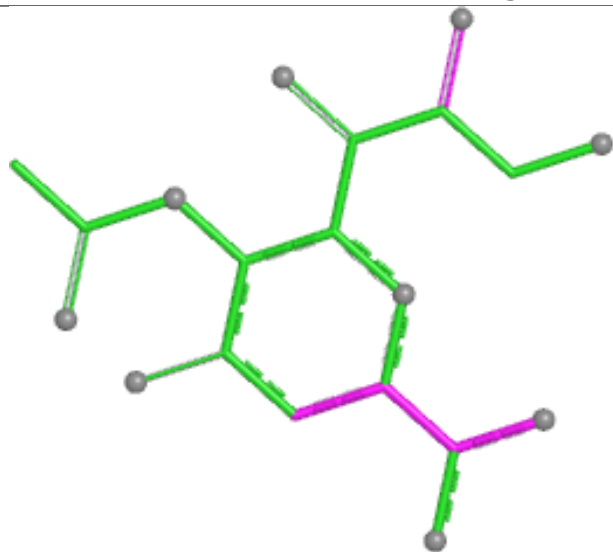


Torsions

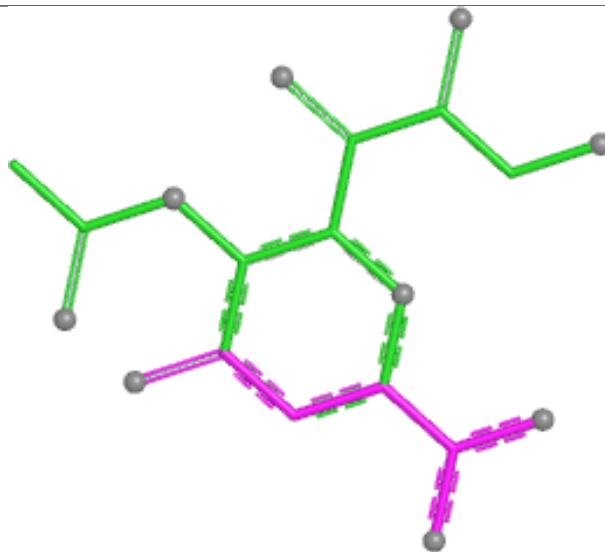


Rings

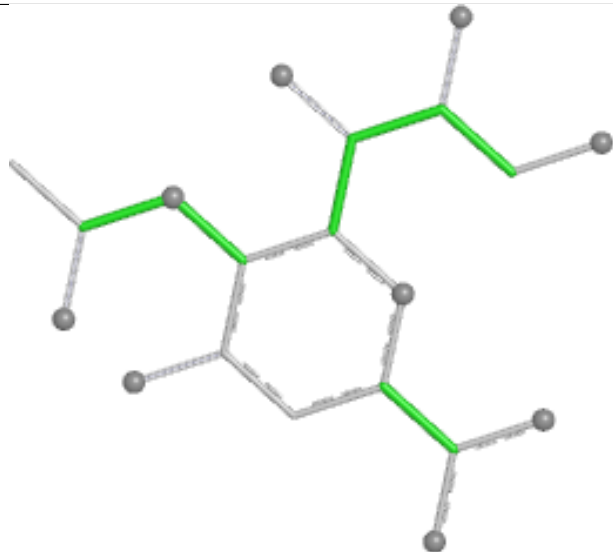
Ligand DAN C 1572



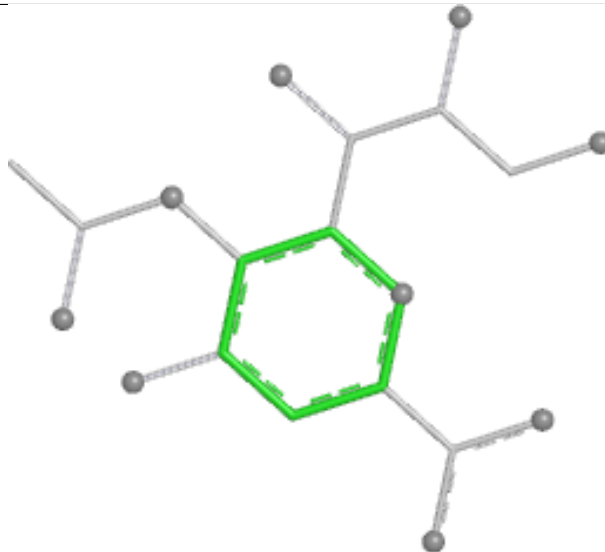
Bond lengths



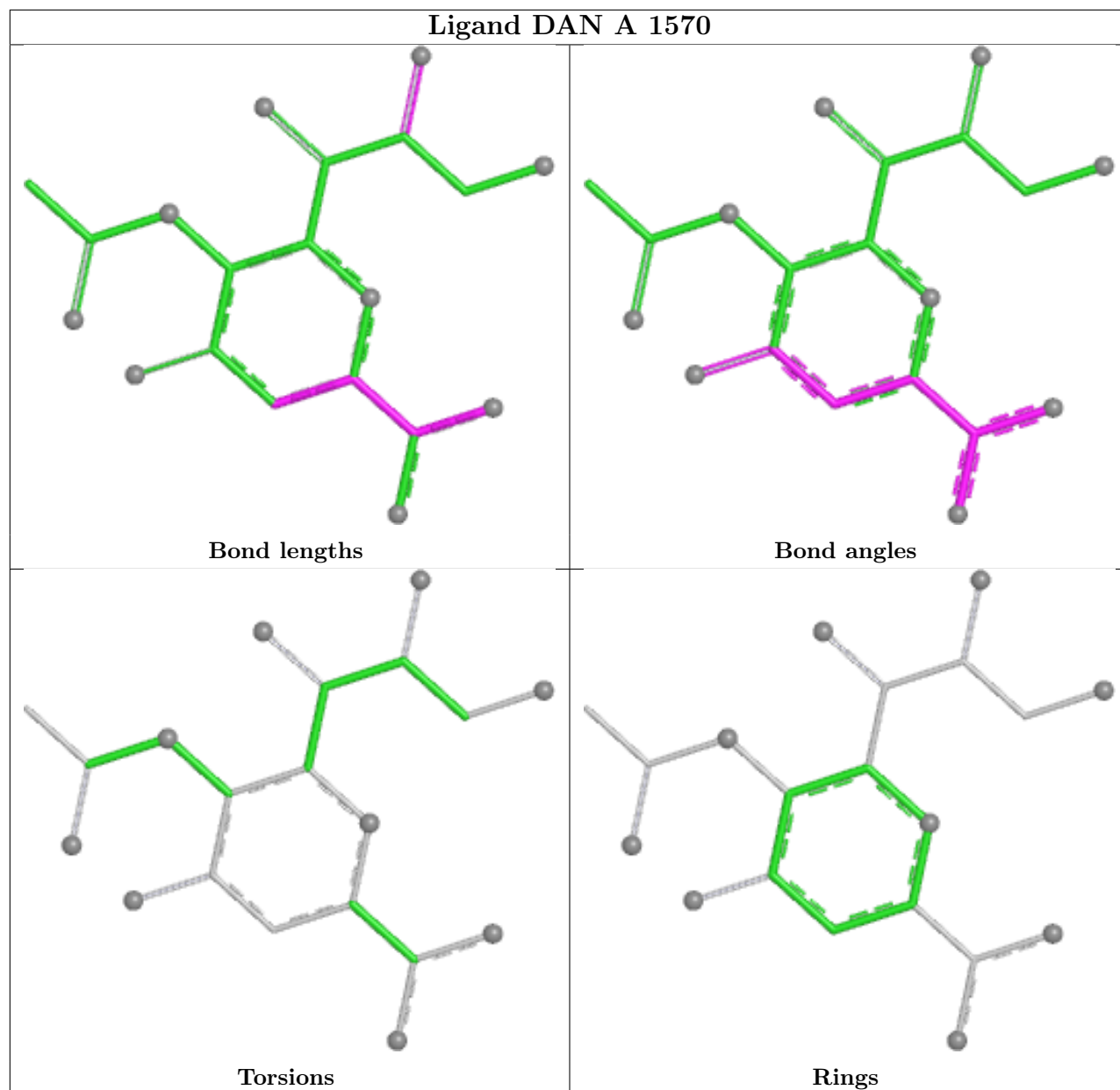
Bond angles



Torsions



Rings



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.