



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:21 AM UTC

PDB ID : 1E0O / pdb_00001e0o
Title : CRYSTAL STRUCTURE OF A TERNARY FGF1-FGFR2-HEPARIN COMPLEX
Authors : Pellegrini, L.; Burke, D.F.; von Delft, F.; Mulloy, B.; Blundell, T.L.
Deposited on : 2000-04-03
Resolution : 2.80 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
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

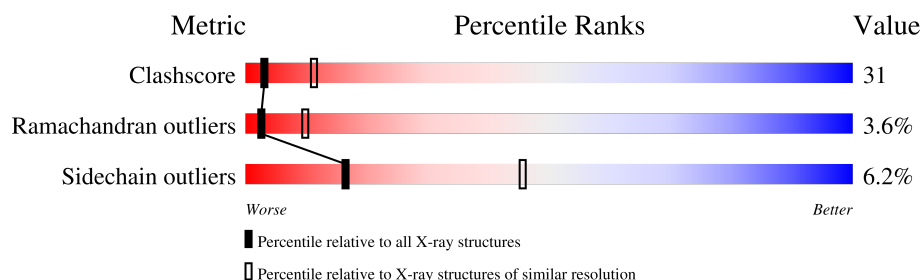
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)

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	140	34% 47% 10% 8%
1	C	140	41% 46% 6% 7%
2	B	219	46% 38% 6% 10%
2	D	219	44% 38% 7% 11%
3	E	10	60% 40%

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
5	SO4	D	401	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5247 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.

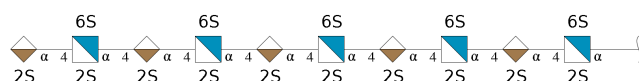
- Molecule 1 is a protein called FIBROBLAST GROWTH FACTOR 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	129	Total	C	N	O	S	0	0	0
			979	622	173	180	4			
1	C	130	Total	C	N	O	S	0	0	0
			1017	642	178	193	4			

- Molecule 2 is a protein called FIBROBLAST GROWTH FACTOR RECEPTOR 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	197	Total	C	N	O	S	0	0	0
			1505	955	266	276	8			
2	D	196	Total	C	N	O	S	0	0	0
			1500	950	266	276	8			

- Molecule 3 is an oligosaccharide called 2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose.

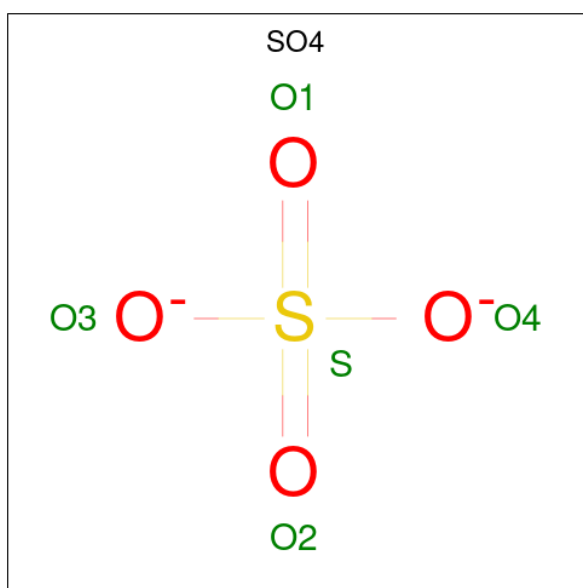


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	10	Total	C	N	O	S	0	0	0
			175	60	5	95	15			

- Molecule 4 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Ni 1 1	0	0
4	B	2	Total Ni 2 2	0	0
4	C	1	Total Ni 1 1	0	0
4	D	1	Total Ni 1 1	0	0

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	1	Total O S 5 4 1	0	0
5	C	1	Total O S 5 4 1	0	0
5	D	1	Total O S 5 4 1	0	0
5	D	1	Total O S 5 4 1	0	0
5	D	1	Total O S 5 4 1	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	7	Total O 7 7	0	0

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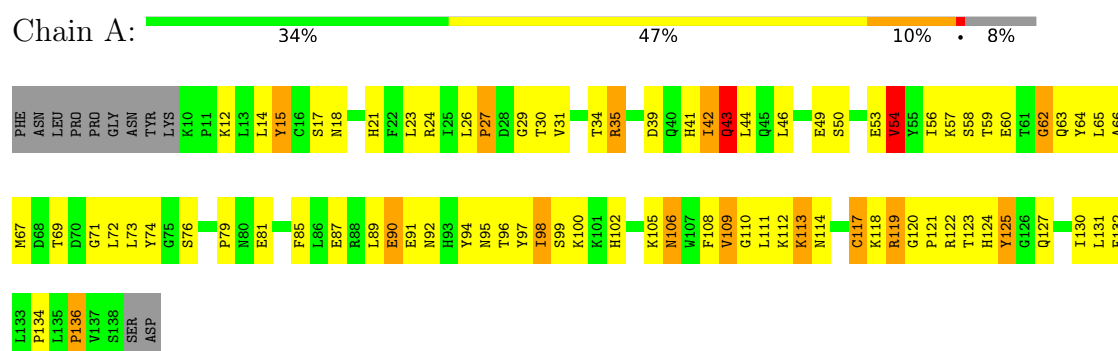
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	12	Total 12	O 12	0	0
6	C	7	Total 7	O 7	0	0
6	D	15	Total 15	O 15	0	0

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.

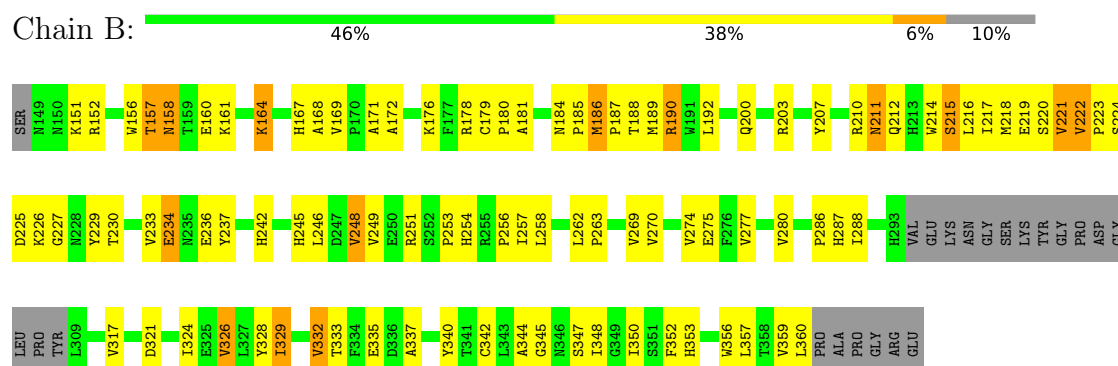
• Molecule 1: FIBROBLAST GROWTH FACTOR 1



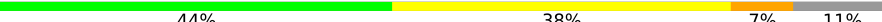
• Molecule 1: FIBROBLAST GROWTH FACTOR 1

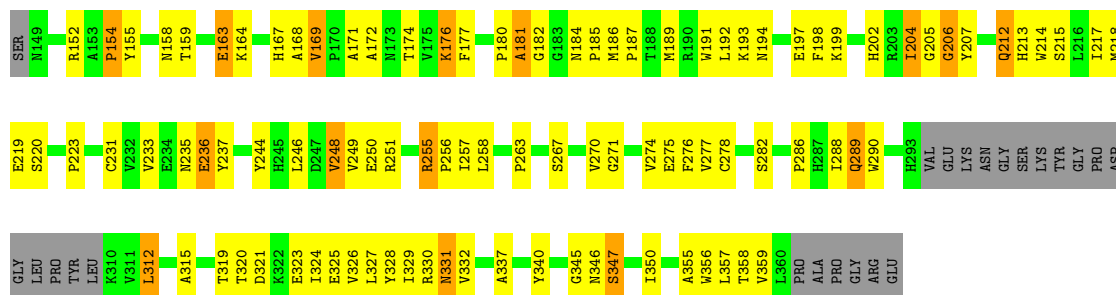


• Molecule 2: FIBROBLAST GROWTH FACTOR RECEPTOR 2

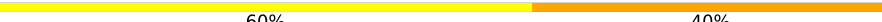


• Molecule 2: FIBROBLAST GROWTH FACTOR RECEPTOR 2

Chain D:  44% 38% 7% 11%



- Molecule 3: 2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose

Chain E:  60% 40%



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 62	Depositor
Cell constants a, b, c, α , β , γ	195.90 Å 195.90 Å 68.86 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	24.20 – 2.80	Depositor
% Data completeness (in resolution range)	98.3 (24.20-2.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.274 , 0.309	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5247	wwPDB-VP
Average B, all atoms (Å ²)	69.0	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: IDS, NI, SGN, SO4

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	0.49	0/1002	1.06	7/1360 (0.5%)
1	C	0.51	0/1040	1.05	5/1407 (0.4%)
2	B	0.52	0/1544	1.05	9/2109 (0.4%)
2	D	0.55	0/1539	1.03	7/2102 (0.3%)
All	All	0.52	0/5125	1.05	28/6978 (0.4%)

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	1

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	169	VAL	N-CA-C	7.18	117.12	109.01
1	C	54	VAL	N-CA-C	6.86	118.03	108.23
1	A	125	TYR	N-CA-C	6.58	119.00	111.11
1	A	54	VAL	N-CA-C	6.45	119.60	108.95
2	B	326	VAL	N-CA-C	6.21	117.15	107.28
1	A	109	VAL	N-CA-C	-6.20	101.00	109.55
2	B	234	GLU	N-CA-C	6.19	118.48	109.07
1	A	27	PRO	N-CA-C	-6.01	105.54	113.65
2	D	215	SER	N-CA-C	6.00	118.53	109.41
1	C	94	TYR	N-CA-C	-5.91	100.52	109.85
1	A	35	ARG	N-CA-C	-5.89	105.69	114.64
1	A	50	SER	N-CA-C	-5.87	100.73	109.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	90	GLU	N-CA-C	5.85	118.09	110.43
2	B	249	VAL	N-CA-C	5.77	116.17	107.75
1	C	100	LYS	N-CA-C	-5.64	104.83	110.97
1	A	42	ILE	N-CA-C	-5.60	108.39	113.71
2	D	249	VAL	CB-CA-C	-5.60	103.71	110.99
2	D	169	VAL	CA-C-N	5.59	125.59	119.89
2	D	169	VAL	C-N-CA	5.59	125.59	119.89
1	C	29	GLY	N-CA-C	-5.49	107.53	115.27
2	D	236	GLU	N-CA-C	5.48	119.46	112.34
2	D	164	LYS	N-CA-C	5.43	117.33	108.63
2	B	222	VAL	CB-CA-C	-5.29	103.46	110.97
2	B	186	MET	CA-C-N	5.24	125.52	119.92
2	B	186	MET	C-N-CA	5.24	125.52	119.92
2	D	181	ALA	N-CA-C	5.22	117.91	109.40
2	B	227	GLY	N-CA-C	5.13	116.95	110.38
2	B	215	SER	N-CA-C	5.10	117.42	109.52

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	15	TYR	Sidechain

5.2 Too-close contacts [i](#)

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	979	0	932	91	0
1	C	1017	0	982	60	0
2	B	1505	0	1419	88	0
2	D	1500	0	1412	86	0
3	E	175	0	56	8	0
4	A	1	0	0	0	0
4	B	2	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	C	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	15	0	0	2	0
6	A	7	0	0	2	0
6	B	12	0	0	5	0
6	C	7	0	0	0	0
6	D	15	0	0	3	0
All	All	5247	0	4801	315	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 31.

All (315) 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)
2:B:345:GLY:HA3	2:B:350:ILE:HG22	1.34	1.05
2:D:274:VAL:HG21	2:D:357:LEU:HD13	1.46	0.97
2:B:277:VAL:HG22	2:B:326:VAL:HG12	1.50	0.92
2:D:345:GLY:HA3	2:D:350:ILE:HG22	1.56	0.86
1:A:24:ARG:HG3	1:A:41:HIS:O	1.74	0.86
1:C:119:ARG:HH21	3:E:4:IDS:H2	1.41	0.84
2:B:211:ASN:N	2:B:211:ASN:HD22	1.78	0.82
2:D:176:LYS:HE3	2:D:217:ILE:HD11	1.64	0.79
2:B:274:VAL:HG21	2:B:357:LEU:HD13	1.65	0.77
2:D:186:MET:HE2	2:D:214:TRP:HZ2	1.50	0.76
1:A:74:TYR:HE1	1:A:76:SER:HB2	1.51	0.76
2:B:245:HIS:HB2	6:B:510:HOH:O	1.86	0.75
2:B:211:ASN:N	2:B:211:ASN:ND2	2.32	0.75
1:C:27:PRO:HG3	1:C:41:HIS:HE2	1.50	0.75
2:D:263:PRO:HB2	2:D:355:ALA:HB2	1.70	0.74
2:D:223:PRO:HA	2:D:248:VAL:HG11	1.69	0.72
2:D:282:SER:HB3	2:D:286:PRO:HG3	1.71	0.72
1:C:11:PRO:O	1:C:137:VAL:HG22	1.89	0.71
2:D:223:PRO:HA	2:D:248:VAL:CG1	2.21	0.70
2:D:337:ALA:HB2	2:D:359:VAL:HG23	1.73	0.70
2:B:211:ASN:ND2	2:B:211:ASN:H	1.86	0.70
2:D:186:MET:HE3	2:D:187:PRO:HD2	1.71	0.70
2:D:212:GLN:HG2	2:D:213:HIS:CD2	2.24	0.70
2:D:274:VAL:CG1	2:D:329:ILE:HB	2.22	0.70
2:D:189:MET:HE3	2:D:191:TRP:HE1	1.56	0.69
1:C:107:TRP:CZ2	1:C:121:PRO:HG3	2.28	0.69
2:B:321:ASP:O	2:B:324:ILE:HG22	1.92	0.69
2:B:256:PRO:HG2	2:B:345:GLY:HA2	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:GLU:HG3	1:A:97:TYR:CE1	2.29	0.68
1:A:35:ARG:NH1	2:B:164:LYS:HA	2.07	0.67
2:B:230:THR:HG23	2:B:242:HIS:O	1.95	0.67
1:C:123:THR:HA	1:C:127:GLN:OE1	1.95	0.67
2:D:321:ASP:O	2:D:324:ILE:HG22	1.96	0.66
1:C:92:ASN:OD1	1:C:94:TYR:HD2	1.79	0.66
1:C:118:LYS:HE3	1:C:122:ARG:O	1.96	0.65
1:A:64:TYR:CE1	1:A:79:PRO:HG3	2.32	0.65
1:A:73:LEU:HD11	1:A:109:VAL:HG13	1.77	0.65
1:C:35:ARG:NH1	2:D:163:GLU:O	2.30	0.64
2:B:350:ILE:HD12	2:B:350:ILE:O	1.97	0.64
1:C:12:LYS:NZ	1:C:46:LEU:HD12	2.12	0.64
1:C:105:LYS:HD3	1:C:107:TRP:CZ2	2.33	0.63
1:A:113:LYS:HE2	3:E:7:SGN:H3	1.80	0.63
1:C:12:LYS:CE	1:C:46:LEU:HD12	2.29	0.63
2:B:223:PRO:HA	2:B:248:VAL:HG13	1.79	0.63
2:B:236:GLU:HG3	2:B:237:TYR:CD1	2.34	0.62
1:C:27:PRO:HG3	1:C:41:HIS:NE2	2.13	0.62
2:D:235:ASN:OD1	2:D:235:ASN:C	2.42	0.62
2:B:345:GLY:HA3	2:B:350:ILE:CG2	2.22	0.61
1:C:18:ASN:HB2	1:C:128:LYS:O	2.00	0.61
1:A:67:MET:SD	1:A:71:GLY:HA2	2.41	0.61
1:C:92:ASN:O	1:C:93:HIS:HB2	2.00	0.61
2:D:171:ALA:O	2:D:172:ALA:HB3	2.01	0.61
2:D:329:ILE:HD11	2:D:340:TYR:CE2	2.35	0.61
2:B:176:LYS:HE3	2:B:178:ARG:HD3	1.83	0.60
1:A:72:LEU:HD21	1:A:119:ARG:HG2	1.82	0.60
2:B:350:ILE:HD12	2:B:350:ILE:C	2.26	0.60
2:D:257:ILE:HA	6:D:502:HOH:O	2.02	0.60
1:C:61:THR:O	1:C:63:GLN:HG3	2.02	0.59
2:D:167:HIS:O	2:D:246:LEU:HD12	2.02	0.59
2:B:236:GLU:HG3	2:B:237:TYR:CE1	2.38	0.59
2:D:320:THR:OG1	2:D:323:GLU:HG3	2.01	0.59
2:D:152:ARG:HG3	2:D:152:ARG:HH11	1.68	0.59
2:B:184:ASN:HA	2:B:185:PRO:C	2.28	0.58
2:D:350:ILE:HD12	2:D:350:ILE:C	2.28	0.58
1:A:64:TYR:CZ	1:A:79:PRO:HG3	2.38	0.58
1:A:130:ILE:HG13	1:A:131:LEU:HD23	1.86	0.58
2:D:236:GLU:HG3	2:D:237:TYR:CE1	2.39	0.58
1:A:134:PRO:HD2	2:B:251:ARG:NH1	2.19	0.58
2:B:256:PRO:HG2	2:B:345:GLY:CA	2.34	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:THR:HG22	1:A:59:THR:O	2.02	0.58
2:B:210:ARG:NH2	2:B:212:GLN:CB	2.66	0.58
2:B:337:ALA:HB2	2:B:359:VAL:HG23	1.86	0.58
2:B:157:THR:CG2	2:B:181:ALA:C	2.77	0.57
1:C:31:VAL:HG23	1:C:74:TYR:HA	1.85	0.57
1:C:50:SER:HB3	1:C:53:GLU:CD	2.30	0.57
1:A:34:THR:OG1	1:A:35:ARG:N	2.38	0.57
2:B:171:ALA:O	2:B:172:ALA:HB3	2.03	0.57
2:D:271:GLY:O	2:D:331:ASN:ND2	2.38	0.57
2:B:180:PRO:HA	2:B:214:TRP:CD1	2.40	0.57
2:B:157:THR:HG23	2:B:181:ALA:C	2.30	0.56
2:B:258:LEU:CD1	2:B:280:VAL:HG22	2.35	0.56
1:C:58:SER:C	1:C:60:GLU:H	2.14	0.56
1:A:119:ARG:HD2	3:E:3:SGN:O1S	2.06	0.56
1:A:110:GLY:O	1:A:111:LEU:HG	2.06	0.56
1:C:14:LEU:HG	1:C:44:LEU:HD12	1.86	0.56
1:C:64:TYR:CZ	1:C:79:PRO:HG3	2.40	0.56
2:B:222:VAL:O	2:B:248:VAL:HG11	2.05	0.56
1:A:27:PRO:HG3	1:A:41:HIS:NE2	2.21	0.55
1:A:35:ARG:HH11	2:B:164:LYS:HA	1.70	0.55
1:C:68:ASP:OD2	1:C:72:LEU:HB3	2.07	0.55
1:C:89:LEU:HD13	1:C:95:ASN:OD1	2.08	0.54
1:C:61:THR:HB	1:C:63:GLN:HG3	1.89	0.54
1:A:124:HIS:ND1	1:A:125:TYR:N	2.56	0.54
2:B:160:GLU:HG2	2:B:161:LYS:N	2.22	0.54
2:B:207:TYR:CD2	2:B:216:LEU:HD11	2.43	0.54
2:B:329:ILE:HD11	2:B:340:TYR:CE2	2.43	0.54
1:C:12:LYS:HE3	1:C:46:LEU:HD12	1.90	0.54
1:A:57:LYS:HA	1:A:63:GLN:O	2.08	0.54
1:C:89:LEU:HD13	2:D:251:ARG:NH2	2.23	0.54
1:A:24:ARG:NH1	1:A:39:ASP:OD2	2.40	0.53
1:A:98:ILE:CG2	1:A:99:SER:N	2.70	0.53
2:B:274:VAL:HG21	2:B:357:LEU:CD1	2.36	0.53
2:B:353:HIS:HE1	6:B:511:HOH:O	1.90	0.53
1:A:112:LYS:O	1:A:114:ASN:N	2.42	0.53
2:D:267:SER:HA	2:D:358:THR:O	2.08	0.53
1:A:58:SER:O	1:A:62:GLY:HA2	2.09	0.53
1:C:25:ILE:O	1:C:41:HIS:HB3	2.09	0.53
2:D:274:VAL:HG12	2:D:329:ILE:HB	1.91	0.53
1:A:113:LYS:HG3	3:E:6:IDS:O2S	2.09	0.53
2:D:256:PRO:HG2	2:D:345:GLY:HA2	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:ILE:O	1:A:44:LEU:HG	2.09	0.53
2:B:219:GLU:O	2:B:220:SER:C	2.52	0.53
1:A:72:LEU:CD2	1:A:119:ARG:HG2	2.38	0.53
1:C:57:LYS:HA	1:C:63:GLN:O	2.09	0.53
1:A:105:LYS:O	1:A:106:ASN:HB2	2.08	0.52
1:A:95:ASN:HB3	1:A:97:TYR:HE1	1.74	0.52
2:B:188:THR:HG22	2:B:234:GLU:O	2.10	0.52
1:A:71:GLY:HA3	1:A:121:PRO:HD3	1.91	0.52
2:D:155:TYR:CE1	2:D:182:GLY:C	2.88	0.52
2:D:236:GLU:HG3	2:D:237:TYR:CD1	2.45	0.51
1:C:61:THR:O	1:C:62:GLY:C	2.53	0.51
2:D:315:ALA:HA	2:D:319:THR:OG1	2.11	0.51
1:A:127:GLN:O	1:A:130:ILE:HG12	2.11	0.51
1:A:31:VAL:HG23	1:A:74:TYR:HA	1.92	0.51
2:B:176:LYS:HG3	2:B:217:ILE:HG12	1.92	0.51
2:D:176:LYS:HE3	2:D:217:ILE:CD1	2.37	0.51
1:A:12:LYS:CE	1:A:46:LEU:HD12	2.39	0.51
2:D:181:ALA:HB1	2:D:233:VAL:HG11	1.93	0.51
1:A:46:LEU:HD23	1:A:56:ILE:HG12	1.92	0.51
1:A:69:THR:O	1:A:69:THR:HG22	2.08	0.51
2:B:210:ARG:NH1	2:B:211:ASN:ND2	2.59	0.51
2:D:172:ALA:O	2:D:220:SER:HA	2.09	0.51
2:D:177:PHE:HB3	2:D:244:TYR:CD2	2.46	0.51
2:D:274:VAL:HG13	2:D:329:ILE:HB	1.91	0.51
1:A:112:LYS:HA	3:E:6:IDS:O3S	2.11	0.51
1:A:46:LEU:CD2	1:A:56:ILE:HG12	2.41	0.50
2:B:216:LEU:HD12	2:B:217:ILE:H	1.75	0.50
1:C:34:THR:OG1	1:C:35:ARG:N	2.44	0.50
1:A:53:GLU:OE2	1:A:100:LYS:HE3	2.11	0.50
1:C:94:TYR:OH	2:D:169:VAL:HG12	2.12	0.50
1:A:71:GLY:O	1:A:119:ARG:HA	2.11	0.50
2:D:174:THR:OG1	2:D:219:GLU:HA	2.11	0.50
2:B:262:LEU:HA	2:B:263:PRO:C	2.37	0.50
1:A:58:SER:C	1:A:60:GLU:H	2.19	0.50
2:B:171:ALA:O	2:B:172:ALA:CB	2.60	0.50
2:B:188:THR:O	2:B:233:VAL:HA	2.12	0.49
1:C:80:ASN:O	1:C:83:CYS:HB2	2.11	0.49
1:A:91:GLU:HG3	6:A:305:HOH:O	2.11	0.49
1:A:117:CYS:SG	1:A:118:LYS:N	2.85	0.49
2:D:276:PHE:HZ	2:D:357:LEU:HB2	1.76	0.49
1:C:66:ALA:HA	1:C:82:GLU:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:ILE:O	1:A:44:LEU:N	2.42	0.49
2:B:223:PRO:HA	2:B:248:VAL:CG1	2.42	0.49
1:C:105:LYS:O	1:C:106:ASN:HB2	2.12	0.49
2:D:154:PRO:HA	2:D:182:GLY:O	2.13	0.49
2:D:191:TRP:O	2:D:192:LEU:HD12	2.11	0.49
2:D:276:PHE:CZ	2:D:357:LEU:HB2	2.48	0.49
2:D:337:ALA:HB2	2:D:359:VAL:CG2	2.42	0.49
1:A:123:THR:HG22	1:A:124:HIS:N	2.26	0.49
2:B:345:GLY:CA	2:B:350:ILE:HG22	2.25	0.49
1:A:112:LYS:NZ	1:A:122:ARG:HH21	2.10	0.48
2:B:190:ARG:HD3	2:B:234:GLU:OE2	2.13	0.48
2:D:213:HIS:CE1	6:D:501:HOH:O	2.64	0.48
2:D:321:ASP:HA	2:D:324:ILE:HG22	1.94	0.48
2:B:254:HIS:HA	2:B:348:ILE:HG21	1.95	0.48
2:B:274:VAL:HG22	2:B:275:GLU:N	2.28	0.48
2:D:152:ARG:HG3	2:D:152:ARG:NH1	2.29	0.48
2:B:176:LYS:HD3	2:B:217:ILE:HD11	1.94	0.48
2:B:200:GLN:HA	2:B:207:TYR:CE1	2.49	0.48
1:A:74:TYR:CE1	1:A:76:SER:HB2	2.39	0.48
1:A:105:LYS:O	1:A:106:ASN:CB	2.61	0.48
1:A:124:HIS:CG	1:A:125:TYR:N	2.82	0.48
1:A:134:PRO:HD2	2:B:251:ARG:HH12	1.77	0.48
2:D:277:VAL:HG22	2:D:326:VAL:HG12	1.95	0.48
2:B:333:THR:O	2:B:359:VAL:HG21	2.14	0.47
1:A:98:ILE:HD13	1:A:106:ASN:HA	1.95	0.47
2:D:213:HIS:HE1	6:D:501:HOH:O	1.97	0.47
1:C:64:TYR:CE1	1:C:79:PRO:HG3	2.49	0.47
1:A:12:LYS:HE3	1:A:46:LEU:HD12	1.97	0.47
1:A:92:ASN:OD1	1:A:94:TYR:HD2	1.97	0.47
1:C:38:SER:O	1:C:39:ASP:C	2.58	0.47
2:D:174:THR:HG23	2:D:218:MET:O	2.15	0.47
2:D:177:PHE:HB3	2:D:244:TYR:CE2	2.50	0.47
2:D:329:ILE:HG22	2:D:329:ILE:O	2.14	0.47
1:A:26:LEU:HD12	1:A:30:THR:HB	1.96	0.47
1:C:50:SER:HB3	1:C:53:GLU:OE1	2.15	0.47
2:B:186:MET:HE3	2:B:187:PRO:HD2	1.97	0.46
2:B:353:HIS:CE1	6:B:511:HOH:O	2.66	0.46
2:D:193:LYS:O	2:D:194:ASN:C	2.58	0.46
1:A:113:LYS:CE	3:E:7:SGN:H3	2.42	0.46
2:D:326:VAL:HG23	2:D:326:VAL:O	2.16	0.46
1:A:73:LEU:CD1	1:A:109:VAL:HG13	2.43	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:176:LYS:HG3	2:B:217:ILE:CG1	2.45	0.46
1:C:64:TYR:CE2	1:C:79:PRO:HG3	2.51	0.46
2:D:255:ARG:NE	5:D:401:SO4:O4	2.48	0.46
1:A:43:GLN:HE21	1:A:43:GLN:HB3	1.55	0.46
2:D:278:CYS:O	2:D:325:GLU:HG3	2.16	0.46
2:B:207:TYR:CD2	2:B:216:LEU:CD1	2.99	0.46
1:C:80:ASN:O	1:C:83:CYS:N	2.49	0.46
1:A:65:LEU:O	1:A:85:PHE:HE2	1.98	0.45
1:A:113:LYS:HE3	3:E:6:IDS:O2S	2.16	0.45
1:A:98:ILE:HG23	1:A:99:SER:N	2.31	0.45
1:C:18:ASN:HB2	1:C:129:ALA:HA	1.98	0.45
2:D:184:ASN:HA	2:D:185:PRO:C	2.42	0.45
1:A:97:TYR:CD1	1:A:97:TYR:N	2.84	0.45
2:B:340:TYR:N	2:B:340:TYR:CD1	2.84	0.45
1:C:89:LEU:HD13	2:D:251:ARG:HH22	1.80	0.45
2:D:198:PHE:CE2	2:D:207:TYR:CD2	3.04	0.45
2:D:251:ARG:HD2	2:D:251:ARG:HA	1.80	0.45
1:A:112:LYS:HD2	1:A:118:LYS:HD2	1.98	0.45
1:C:105:LYS:HD3	1:C:107:TRP:CE2	2.51	0.45
2:D:267:SER:OG	2:D:358:THR:HB	2.16	0.45
1:A:54:VAL:HG21	1:A:87:GLU:HB2	1.97	0.45
2:D:290:TRP:NE1	2:D:327:LEU:HB2	2.32	0.45
1:A:119:ARG:HB3	1:A:121:PRO:HD2	1.99	0.45
1:A:99:SER:O	1:A:100:LYS:C	2.59	0.44
2:B:256:PRO:CG	2:B:345:GLY:HA2	2.46	0.44
1:C:64:TYR:HB2	1:C:83:CYS:SG	2.57	0.44
1:C:102:HIS:HB3	1:C:105:LYS:HD2	1.98	0.44
2:B:178:ARG:HD2	2:B:215:SER:HB3	1.98	0.44
2:B:269:VAL:HA	2:B:360:LEU:O	2.17	0.44
2:D:204:ILE:HD12	2:D:205:GLY:N	2.32	0.44
2:B:203:ARG:HD2	6:B:505:HOH:O	2.17	0.44
1:A:108:PHE:HB2	1:A:130:ILE:HG22	1.99	0.44
1:C:46:LEU:HD23	1:C:56:ILE:HG12	1.99	0.44
1:A:130:ILE:HG13	1:A:131:LEU:CD2	2.47	0.44
1:A:49:GLU:HA	1:A:49:GLU:OE1	2.18	0.44
1:A:124:HIS:CG	1:A:125:TYR:H	2.36	0.44
2:B:167:HIS:O	2:B:246:LEU:HD12	2.18	0.44
1:C:60:GLU:C	1:C:62:GLY:H	2.26	0.44
1:A:122:ARG:O	6:A:301:HOH:O	2.21	0.44
2:D:186:MET:HE3	2:D:187:PRO:CD	2.44	0.44
2:D:205:GLY:O	2:D:206:GLY:C	2.61	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:127:GLN:O	1:C:130:ILE:HG12	2.17	0.43
2:D:158:ASN:HB3	2:D:180:PRO:HG2	2.00	0.43
1:C:98:ILE:CG2	1:C:99:SER:N	2.81	0.43
2:D:270:VAL:HA	2:D:332:VAL:HG23	2.00	0.43
2:D:326:VAL:HB	2:D:328:TYR:HE1	1.84	0.43
1:C:17:SER:O	1:C:18:ASN:C	2.61	0.43
2:B:275:GLU:HG3	2:B:328:TYR:CD1	2.54	0.43
2:B:210:ARG:HH11	2:B:211:ASN:HD21	1.66	0.43
2:B:224:SER:HB2	6:B:502:HOH:O	2.19	0.43
1:C:43:GLN:OE1	1:C:137:VAL:HG13	2.19	0.43
1:A:125:TYR:O	1:A:130:ILE:HD11	2.18	0.43
2:B:225:ASP:O	2:B:226:LYS:C	2.62	0.43
1:C:61:THR:O	1:C:63:GLN:N	2.51	0.43
2:B:350:ILE:C	2:B:350:ILE:CD1	2.92	0.43
2:D:274:VAL:HG22	2:D:275:GLU:N	2.34	0.43
2:B:151:LYS:O	2:B:152:ARG:HB2	2.19	0.43
2:B:181:ALA:CB	2:B:233:VAL:HG11	2.49	0.43
1:C:49:GLU:HB3	1:C:50:SER:H	1.63	0.43
2:D:256:PRO:HG3	2:D:346:ASN:ND2	2.33	0.43
1:A:74:TYR:CD1	1:A:74:TYR:C	2.97	0.42
1:A:96:THR:C	1:A:97:TYR:HD1	2.28	0.42
1:C:15:TYR:CD1	1:C:15:TYR:C	2.96	0.42
1:C:64:TYR:CB	1:C:83:CYS:SG	3.08	0.42
2:D:258:LEU:N	2:D:258:LEU:HD12	2.35	0.42
1:A:95:ASN:HB3	1:A:97:TYR:CE1	2.53	0.42
1:A:73:LEU:HD21	1:A:109:VAL:HG22	2.01	0.42
2:B:286:PRO:HG2	2:B:288:ILE:HD11	2.01	0.42
2:D:199:LYS:H	2:D:202:HIS:CE1	2.38	0.42
1:A:14:LEU:HG	1:A:44:LEU:HD12	2.00	0.42
1:A:65:LEU:O	1:A:85:PHE:CE2	2.72	0.42
2:B:176:LYS:CD	2:B:217:ILE:HD11	2.49	0.42
2:B:189:MET:HE1	2:B:214:TRP:C	2.45	0.42
2:B:218:MET:HE1	2:B:229:TYR:CZ	2.54	0.42
2:B:269:VAL:O	2:B:270:VAL:C	2.61	0.42
2:D:191:TRP:CH2	2:D:231:CYS:HB3	2.54	0.42
1:A:14:LEU:O	1:A:132:PHE:HD2	2.02	0.42
1:A:123:THR:CG2	1:A:124:HIS:N	2.83	0.42
1:A:123:THR:HA	1:A:127:GLN:OE1	2.20	0.42
2:B:156:TRP:HA	2:B:156:TRP:CE3	2.55	0.42
2:B:157:THR:O	2:B:158:ASN:OD1	2.37	0.42
2:D:152:ARG:HB3	2:D:184:ASN:H	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:255:ARG:HG2	5:D:401:SO4:O1	2.20	0.42
1:A:27:PRO:C	1:A:29:GLY:H	2.26	0.42
1:C:15:TYR:CE1	2:D:168:ALA:HB3	2.55	0.42
1:A:89:LEU:O	2:B:257:ILE:HD13	2.20	0.41
2:B:287:HIS:O	2:B:344:ALA:HA	2.19	0.41
2:D:326:VAL:HB	2:D:328:TYR:CE1	2.55	0.41
1:A:15:TYR:CZ	2:B:168:ALA:HB3	2.55	0.41
2:B:164:LYS:HE2	2:B:167:HIS:NE2	2.35	0.41
2:B:179:CYS:O	2:B:214:TRP:HB3	2.20	0.41
2:D:197:GLU:O	2:D:202:HIS:CE1	2.73	0.41
1:A:66:ALA:HB3	1:A:74:TYR:CE1	2.56	0.41
1:C:67:MET:CG	1:C:71:GLY:HA2	2.50	0.41
2:B:221:VAL:HG12	2:B:248:VAL:HG21	2.02	0.41
2:D:155:TYR:CZ	2:D:182:GLY:HA3	2.55	0.41
3:E:3:SGN:O3	3:E:4:IDS:O5	2.35	0.41
1:A:17:SER:O	1:A:18:ASN:C	2.63	0.41
2:B:179:CYS:N	2:B:180:PRO:HD3	2.36	0.41
2:B:342:CYS:O	2:B:352:PHE:HA	2.21	0.41
1:A:64:TYR:CD1	1:A:79:PRO:HG3	2.55	0.41
1:C:23:LEU:HB3	1:C:44:LEU:HD11	2.03	0.41
1:C:134:PRO:O	1:C:135:LEU:HD23	2.20	0.41
1:A:69:THR:HA	1:A:102:HIS:CE1	2.56	0.41
1:A:108:PHE:HB2	1:A:130:ILE:CG2	2.51	0.41
1:C:23:LEU:HD21	1:C:65:LEU:HD22	2.02	0.41
1:C:32:ASP:OD1	1:C:32:ASP:C	2.62	0.41
2:D:223:PRO:HA	2:D:248:VAL:HG13	1.99	0.41
2:D:288:ILE:C	2:D:289:GLN:HG3	2.46	0.41
2:D:330:ARG:O	2:D:331:ASN:C	2.63	0.41
2:B:337:ALA:HB2	2:B:359:VAL:CG2	2.48	0.41
2:D:329:ILE:HD11	2:D:340:TYR:HE2	1.81	0.41
1:A:90:GLU:OE2	1:A:125:TYR:CE1	2.74	0.40
2:B:329:ILE:HG22	2:B:332:VAL:HG12	2.03	0.40
1:A:21:HIS:HB3	1:A:34:THR:O	2.20	0.40
2:B:237:TYR:CD1	2:B:237:TYR:N	2.87	0.40
1:A:23:LEU:HA	1:A:23:LEU:HD12	1.90	0.40
2:D:193:LYS:HB2	2:D:198:PHE:CD1	2.56	0.40
2:D:223:PRO:CA	2:D:248:VAL:HG11	2.47	0.40
1:A:119:ARG:O	1:A:120:GLY:C	2.65	0.40
1:C:61:THR:HG21	1:C:63:GLN:OE1	2.21	0.40
2:D:171:ALA:O	2:D:172:ALA:CB	2.67	0.40
2:B:335:GLU:C	2:B:337:ALA:H	2.30	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	127/140 (91%)	101 (80%)	19 (15%)	7 (6%)	1	4
1	C	128/140 (91%)	105 (82%)	16 (12%)	7 (6%)	1	4
2	B	193/219 (88%)	168 (87%)	21 (11%)	4 (2%)	5	20
2	D	192/219 (88%)	166 (86%)	21 (11%)	5 (3%)	4	15
All	All	640/718 (89%)	540 (84%)	77 (12%)	23 (4%)	2	10

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	106	ASN
1	A	113	LYS
2	B	317	VAL
2	B	332	VAL
2	D	312	LEU
1	A	43	GLN
1	A	81	GLU
1	A	136	PRO
2	B	158	ASN
1	C	18	ASN
1	C	41	HIS
1	C	49	GLU
1	C	113	LYS
2	D	206	GLY
2	D	347	SER
2	D	331	ASN
1	C	68	ASP
1	C	117	CYS
1	A	119	ARG
1	C	62	GLY

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Mol	Chain	Res	Type
2	B	329	ILE
2	D	154	PRO
1	A	62	GLY

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	100/123 (81%)	94 (94%)	6 (6%)	17	47
1	C	109/123 (89%)	105 (96%)	4 (4%)	30	65
2	B	155/186 (83%)	145 (94%)	10 (6%)	15	43
2	D	155/186 (83%)	143 (92%)	12 (8%)	12	36
All	All	519/618 (84%)	487 (94%)	32 (6%)	16	45

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	54	VAL
1	A	90	GLU
1	A	98	ILE
1	A	117	CYS
1	A	136	PRO
2	B	157	THR
2	B	164	LYS
2	B	190	ARG
2	B	192	LEU
2	B	211	ASN
2	B	221	VAL
2	B	248	VAL
2	B	253	PRO
2	B	347	SER
2	B	356	TRP
1	C	9	LYS
1	C	61	THR

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Mol	Chain	Res	Type
1	C	91	GLU
1	C	117	CYS
2	D	159	THR
2	D	163	GLU
2	D	176	LYS
2	D	204	ILE
2	D	212	GLN
2	D	248	VAL
2	D	250	GLU
2	D	255	ARG
2	D	289	GLN
2	D	312	LEU
2	D	347	SER
2	D	356	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
2	B	200	GLN
2	B	211	ASN
2	B	241	ASN
2	B	259	GLN
1	C	77	GLN
1	C	106	ASN
2	D	167	HIS
2	D	202	HIS
2	D	213	HIS
2	D	228	ASN
2	D	289	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 ⓘ

10 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
3	SGN	E	1	3	20,20,20	2.24	7 (35%)	25,31,31	1.15	2 (8%)
3	IDS	E	10	3	15,15,17	2.42	5 (33%)	14,22,26	1.70	4 (28%)
3	IDS	E	2	3	16,16,17	1.66	5 (31%)	16,24,26	1.24	2 (12%)
3	SGN	E	3	3	19,19,20	1.98	6 (31%)	23,29,31	1.15	1 (4%)
3	IDS	E	4	3	16,16,17	1.67	6 (37%)	16,24,26	1.52	4 (25%)
3	SGN	E	5	3	19,19,20	2.41	5 (26%)	23,29,31	1.07	1 (4%)
3	IDS	E	6	3	16,16,17	1.46	3 (18%)	16,24,26	1.45	5 (31%)
3	SGN	E	7	3	19,19,20	1.96	5 (26%)	23,29,31	1.06	2 (8%)
3	IDS	E	8	3	16,16,17	1.41	2 (12%)	16,24,26	1.31	2 (12%)
3	SGN	E	9	3	19,19,20	2.10	6 (31%)	23,29,31	1.14	2 (8%)

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	SGN	E	1	3	-	0/11/31/31	0/1/1/1
3	IDS	E	10	3	-	2/9/22/29	0/1/1/1
3	IDS	E	2	3	-	0/9/26/29	0/1/1/1
3	SGN	E	3	3	-	6/11/28/31	0/1/1/1
3	IDS	E	4	3	-	1/9/26/29	0/1/1/1
3	SGN	E	5	3	-	5/11/28/31	0/1/1/1
3	IDS	E	6	3	-	2/9/26/29	0/1/1/1
3	SGN	E	7	3	-	6/11/28/31	0/1/1/1
3	IDS	E	8	3	-	5/9/26/29	0/1/1/1
3	SGN	E	9	3	-	3/11/28/31	0/1/1/1

All (50) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	1	SGN	S1-N2	6.71	1.67	1.59
3	E	5	SGN	S1-N2	6.07	1.67	1.59
3	E	10	IDS	C4-C5	5.86	1.61	1.52
3	E	3	SGN	S1-N2	5.69	1.66	1.59
3	E	9	SGN	C1-C2	5.31	1.59	1.52
3	E	7	SGN	S1-N2	5.12	1.65	1.59
3	E	10	IDS	O5-C5	5.05	1.52	1.43
3	E	5	SGN	O2S-S1	4.92	1.47	1.42
3	E	5	SGN	C1-C2	4.36	1.58	1.52
3	E	9	SGN	S1-N2	4.25	1.64	1.59
3	E	1	SGN	C1-C2	3.80	1.57	1.52
3	E	7	SGN	O2S-S1	3.61	1.46	1.42
3	E	7	SGN	C1-C2	3.47	1.57	1.52
3	E	8	IDS	C4-C5	3.33	1.58	1.53
3	E	5	SGN	O1S-S1	3.28	1.45	1.42
3	E	2	IDS	O2-C2	-3.24	1.42	1.47
3	E	1	SGN	O2S-S1	3.22	1.45	1.42
3	E	3	SGN	O1S-S1	3.13	1.45	1.42
3	E	9	SGN	O2S-S1	2.92	1.45	1.42
3	E	9	SGN	C4-C5	2.85	1.59	1.53
3	E	10	IDS	O2-C2	-2.83	1.43	1.47
3	E	10	IDS	C5-C6	2.80	1.55	1.52
3	E	2	IDS	C4-C5	2.79	1.57	1.53
3	E	9	SGN	O1S-S1	2.79	1.45	1.42
3	E	1	SGN	O1S-S1	2.75	1.45	1.42
3	E	4	IDS	C1-C2	2.73	1.56	1.51
3	E	8	IDS	O6B-C6	-2.70	1.22	1.30
3	E	4	IDS	O2-C2	-2.70	1.43	1.47
3	E	6	IDS	O6B-C6	-2.68	1.22	1.30
3	E	3	SGN	O2S-S1	2.67	1.45	1.42
3	E	7	SGN	O1S-S1	2.66	1.45	1.42
3	E	2	IDS	C1-C2	2.66	1.55	1.51
3	E	4	IDS	O6B-C6	-2.63	1.22	1.30
3	E	2	IDS	O6B-C6	-2.58	1.22	1.30
3	E	6	IDS	C1-C2	2.45	1.55	1.51
3	E	10	IDS	O6B-C6	-2.44	1.22	1.30
3	E	6	IDS	C4-C5	2.44	1.57	1.53
3	E	5	SGN	O5-C5	2.39	1.48	1.43
3	E	9	SGN	O5-C5	2.28	1.47	1.43
3	E	2	IDS	C3-C2	2.25	1.57	1.53
3	E	4	IDS	O5-C5	2.24	1.47	1.43
3	E	1	SGN	C4-C5	2.23	1.57	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	3	SGN	O4-C4	2.17	1.48	1.43
3	E	4	IDS	O4-C4	-2.11	1.37	1.43
3	E	1	SGN	C3-C2	2.09	1.57	1.53
3	E	7	SGN	C4-C5	2.06	1.57	1.53
3	E	4	IDS	C5-C6	2.04	1.57	1.53
3	E	3	SGN	C4-C3	2.03	1.57	1.52
3	E	1	SGN	C2-N2	2.01	1.49	1.47
3	E	3	SGN	C1-C2	2.01	1.55	1.52

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	1	SGN	O1S-S1-O2S	-3.50	112.56	120.36
3	E	10	IDS	O6B-C6-C5	3.46	121.72	112.71
3	E	3	SGN	O1S-S1-O2S	-3.30	113.03	120.36
3	E	9	SGN	O1S-S1-O2S	-3.27	113.10	120.36
3	E	5	SGN	O1S-S1-O2S	-3.03	113.61	120.36
3	E	7	SGN	O1S-S1-O2S	-2.92	113.87	120.36
3	E	4	IDS	O4-C4-C5	-2.90	103.13	109.76
3	E	4	IDS	O6B-C6-O6A	-2.85	117.61	124.08
3	E	10	IDS	O6B-C6-O6A	-2.85	117.62	124.08
3	E	8	IDS	O6B-C6-O6A	-2.79	117.74	124.08
3	E	2	IDS	O6B-C6-O6A	-2.78	117.77	124.08
3	E	6	IDS	O6B-C6-O6A	-2.72	117.92	124.08
3	E	6	IDS	O4-C4-C3	-2.61	104.22	110.38
3	E	4	IDS	C4-C3-C2	2.58	114.92	110.23
3	E	10	IDS	C1-C2-C3	-2.39	106.46	109.84
3	E	10	IDS	C1-O5-C5	2.33	117.59	113.81
3	E	7	SGN	O2S-S1-N2	-2.25	105.12	108.88
3	E	6	IDS	O4-C4-C5	-2.24	104.65	109.76
3	E	4	IDS	O6B-C6-C5	2.23	121.70	113.64
3	E	6	IDS	O6B-C6-C5	2.20	121.58	113.64
3	E	9	SGN	C4-C3-C2	-2.14	107.89	111.02
3	E	8	IDS	O6B-C6-C5	2.13	121.32	113.64
3	E	6	IDS	C4-C3-C2	2.06	113.97	110.23
3	E	2	IDS	O2-C2-C3	2.03	109.78	106.95
3	E	1	SGN	C4-C3-C2	-2.01	107.47	110.40

There are no chirality outliers.

All (30) torsion outliers are listed below:

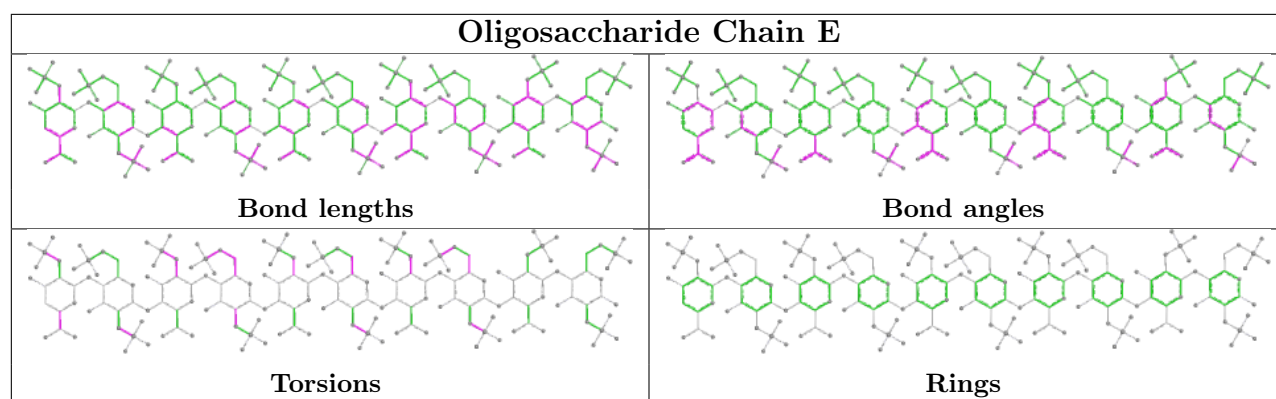
Mol	Chain	Res	Type	Atoms
3	E	3	SGN	C4-C5-C6-O6
3	E	3	SGN	O5-C5-C6-O6
3	E	4	IDS	C3-C2-O2-S
3	E	5	SGN	O5-C5-C6-O6
3	E	5	SGN	C2-N2-S1-O2S
3	E	5	SGN	C2-N2-S1-O3S
3	E	6	IDS	C1-C2-O2-S
3	E	6	IDS	C3-C2-O2-S
3	E	7	SGN	O5-C5-C6-O6
3	E	8	IDS	C1-C2-O2-S
3	E	8	IDS	C3-C2-O2-S
3	E	8	IDS	C2-O2-S-O2S
3	E	9	SGN	C2-N2-S1-O1S
3	E	9	SGN	C2-N2-S1-O3S
3	E	10	IDS	O5-C5-C6-O6A
3	E	3	SGN	C6-O6-S2-O4S
3	E	8	IDS	C2-O2-S-O1S
3	E	3	SGN	C6-O6-S2-O6S
3	E	8	IDS	C2-O2-S-O3S
3	E	7	SGN	C6-O6-S2-O6S
3	E	3	SGN	C6-O6-S2-O5S
3	E	7	SGN	C6-O6-S2-O4S
3	E	7	SGN	C6-O6-S2-O5S
3	E	3	SGN	C2-N2-S1-O2S
3	E	5	SGN	C2-N2-S1-O1S
3	E	5	SGN	C4-C5-C6-O6
3	E	7	SGN	C3-C2-N2-S1
3	E	9	SGN	C2-N2-S1-O2S
3	E	7	SGN	C5-C6-O6-S2
3	E	10	IDS	C2-O2-S-O1S

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	4	IDS	2	0
3	E	6	IDS	3	0
3	E	3	SGN	2	0
3	E	7	SGN	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.



5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 5 are 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	$\# Z > 2$
5	SO4	C	202	-	4,4,4	0.50	0	6,6,6	0.35	0
5	SO4	C	201	-	4,4,4	0.33	0	6,6,6	0.14	0
5	SO4	D	403	-	4,4,4	0.64	0	6,6,6	0.16	0
5	SO4	D	401	-	4,4,4	0.34	0	6,6,6	0.19	0
5	SO4	D	402	-	4,4,4	0.56	0	6,6,6	0.10	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	401	SO4	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

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.