



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:45 PM UTC

PDB ID : 1QCW / pdb_00001qcw
Title : Flavocytochrome B2, ARG289LYS mutant
Authors : Mowat, C.G.; Durley, R.C.E.; Pike, A.D.; Barton, J.D.; Chen, Z.-W.; Mathews, F.S.; Lederer, F.; Reid, G.A.; Chapman, S.K.
Deposited on : 1999-05-07
Resolution : 2.75 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

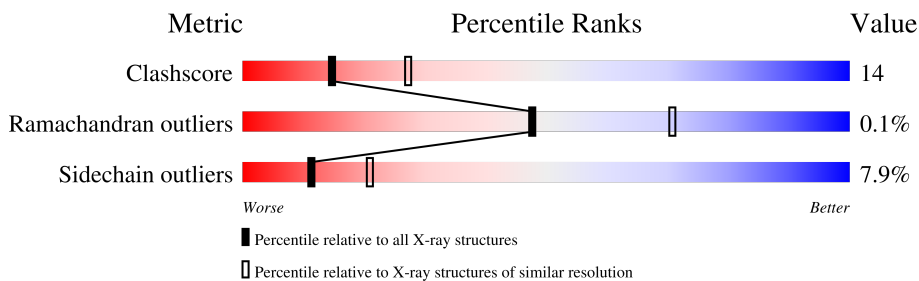
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)

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	410	 60% 30% • 6%
1	B	410	 60% 30% • 6%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6171 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.

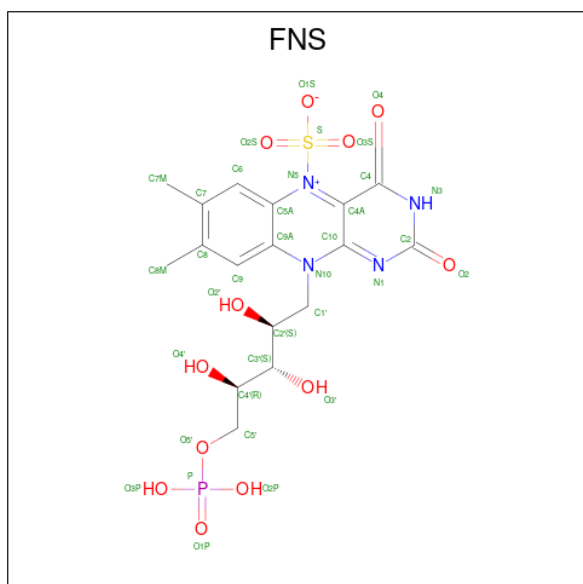
- Molecule 1 is a protein called FLAVOCYTOCHROME B2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	385	Total	C	N	O	S	0	0	0
			3012	1914	509	578	11			
1	B	385	Total	C	N	O	S	0	0	0
			3012	1914	509	578	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	289	LYS	ARG	engineered mutation	UNP P00175
B	289	LYS	ARG	engineered mutation	UNP P00175
A	304	PHE	PRO	engineered mutation	UNP P00175
B	304	PHE	PRO	engineered mutation	UNP P00175

- Molecule 2 is N-SULFO-FLAVIN MONONUCLEOTIDE (CCD ID: FNS) (formula: $C_{17}H_{21}N_4O_{12}PS$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			35	17	4	12	1	1		
2	B	1	Total	C	N	O	P	S	0	0
			35	17	4	12	1	1		

- Molecule 3 is water.

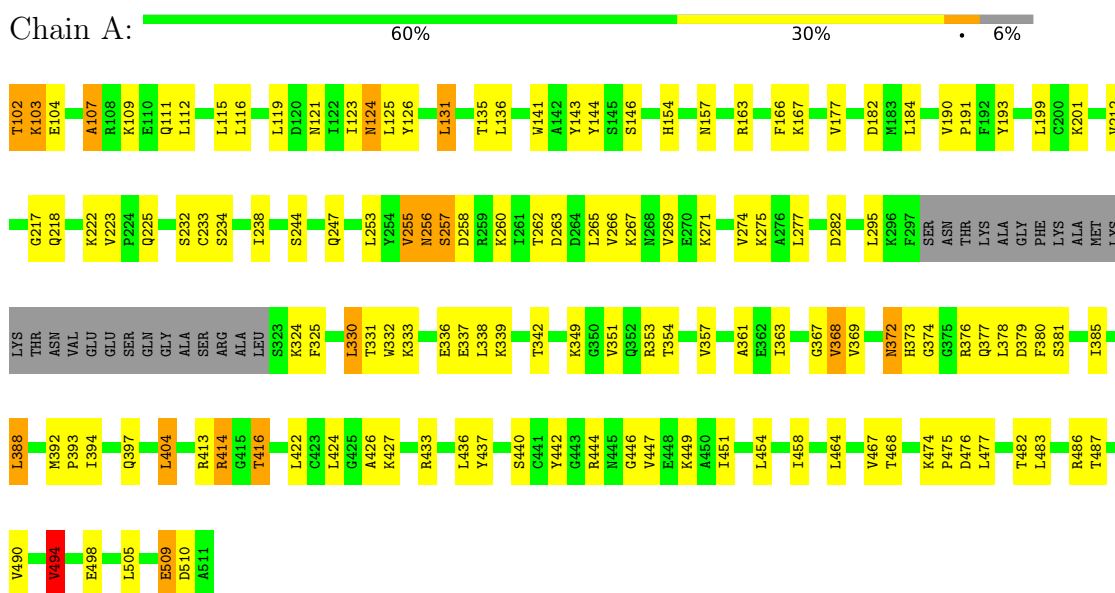
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	50	Total	O	0	0
			50	50		
3	B	27	Total	O	0	0
			27	27		

3 Residue-property plots

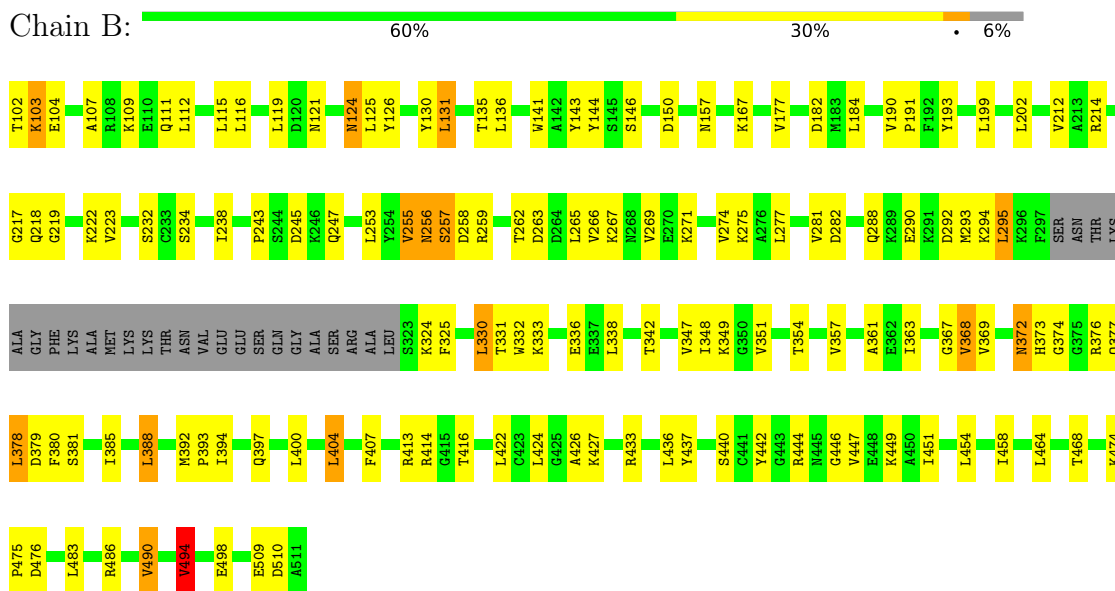
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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: FLAVOCYTOCHROME B2



• Molecule 1: FLAVOCYTOCHROME B2



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	162.57Å 162.57Å 112.01Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 – 2.75	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.75)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.204 , 0.257	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6171	wwPDB-VP
Average B, all atoms (Å ²)	27.0	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: FNS

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	0.51	0/3061	1.00	12/4136 (0.3%)
1	B	0.47	0/3061	1.00	13/4136 (0.3%)
All	All	0.49	0/6122	1.00	25/8272 (0.3%)

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	324	LYS	N-CA-C	-7.58	104.16	113.41
1	A	324	LYS	N-CA-C	-7.29	104.52	113.41
1	A	381	SER	N-CA-C	-6.82	98.51	109.96
1	B	381	SER	N-CA-C	-6.71	98.75	109.76
1	A	275	LYS	N-CA-C	6.64	120.64	112.54
1	B	257	SER	N-CA-C	-6.32	104.39	111.28
1	A	494	VAL	N-CA-C	6.30	117.05	110.62
1	A	167	LYS	CA-C-N	6.21	126.16	119.76
1	A	167	LYS	C-N-CA	6.21	126.16	119.76
1	B	167	LYS	CA-C-N	6.19	126.14	119.76
1	B	167	LYS	C-N-CA	6.19	126.14	119.76
1	B	275	LYS	N-CA-C	6.11	120.00	112.54
1	B	494	VAL	N-CA-C	5.91	116.65	110.62
1	B	107	ALA	N-CA-C	-5.71	105.20	111.82
1	B	357	VAL	N-CA-C	-5.59	105.08	110.72
1	B	146	SER	N-CA-C	5.57	118.72	110.48
1	A	212	VAL	N-CA-C	-5.56	104.94	111.00
1	A	257	SER	N-CA-C	-5.55	105.31	111.36
1	A	107	ALA	N-CA-C	-5.40	105.56	111.82
1	B	212	VAL	N-CA-C	-5.30	105.55	110.53
1	A	357	VAL	N-CA-C	-5.23	105.44	110.72
1	B	202	LEU	N-CA-C	-5.18	105.71	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	146	SER	N-CA-C	5.14	118.09	110.48
1	A	487	THR	N-CA-C	5.11	118.02	110.46
1	B	245	ASP	N-CA-C	-5.08	106.59	112.89

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3012	0	3078	87	0
1	B	3012	0	3078	89	0
2	A	35	0	19	1	0
2	B	35	0	19	2	0
3	A	50	0	0	5	0
3	B	27	0	0	1	0
All	All	6171	0	6194	171	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 14.

All (171) 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:102:THR:HB	1:A:104:GLU:HG3	1.43	1.01
1:B:102:THR:HB	1:B:104:GLU:HG3	1.46	0.97
1:A:256:ASN:HD22	1:A:258:ASP:H	1.18	0.92
1:B:256:ASN:HD22	1:B:258:ASP:H	1.18	0.86
1:B:256:ASN:ND2	1:B:258:ASP:H	1.84	0.75
1:A:256:ASN:ND2	1:A:258:ASP:H	1.84	0.74
1:B:256:ASN:HD22	1:B:257:SER:N	1.86	0.73
1:B:218:GLN:HE22	1:B:444:ARG:HE	1.36	0.72
1:A:218:GLN:HE22	1:A:444:ARG:HE	1.41	0.69
1:A:256:ASN:HD22	1:A:257:SER:N	1.90	0.67
1:A:182:ASP:O	1:A:427:LYS:HE3	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:VAL:O	1:A:274:VAL:HG22	1.96	0.65
1:B:124:ASN:HD22	1:B:126:TYR:H	1.44	0.65
1:B:157:ASN:HD21	1:B:372:ASN:HD21	1.43	0.65
1:B:269:VAL:O	1:B:274:VAL:HG22	1.99	0.63
1:A:143:TYR:O	1:A:376:ARG:NH2	2.33	0.62
1:A:474:LYS:HB3	1:A:475:PRO:HD2	1.82	0.62
1:B:474:LYS:HB3	1:B:475:PRO:HD2	1.82	0.62
1:A:154:HIS:HE1	3:A:10:HOH:O	1.82	0.61
1:B:361:ALA:HB2	1:B:404:LEU:HD12	1.83	0.61
1:A:373:HIS:O	1:A:376:ARG:HG3	2.02	0.60
1:A:361:ALA:HB2	1:A:404:LEU:HD12	1.85	0.59
1:A:486:ARG:HG2	1:B:490:VAL:HG12	1.84	0.58
1:B:331:THR:HG22	1:B:332:TRP:N	2.19	0.58
1:B:182:ASP:O	1:B:427:LYS:HE3	2.03	0.58
1:A:367:GLY:HA2	1:A:404:LEU:HD23	1.86	0.57
1:B:354:THR:HG21	1:B:394:ILE:HD12	1.86	0.57
1:B:136:LEU:HD21	1:B:440:SER:HB3	1.86	0.57
1:A:263:ASP:O	1:A:267:LYS:HG2	2.05	0.57
1:A:509:GLU:O	1:B:115:LEU:HD22	2.05	0.56
1:A:190:VAL:HG21	1:A:223:VAL:HG22	1.88	0.56
1:B:190:VAL:HG21	1:B:223:VAL:HG22	1.88	0.56
1:B:256:ASN:HD22	1:B:258:ASP:N	1.97	0.56
1:A:253:LEU:HD11	1:A:262:THR:HG23	1.88	0.56
2:A:570:FNS:O4'	2:A:570:FNS:H9	2.06	0.55
1:B:333:LYS:HD2	1:B:336:GLU:OE2	2.07	0.55
1:B:255:VAL:HG21	1:B:330:LEU:HD21	1.87	0.55
1:A:256:ASN:HA	1:A:325:PHE:O	2.06	0.55
1:A:136:LEU:HD21	1:A:440:SER:HB3	1.88	0.54
1:A:255:VAL:HG21	1:A:330:LEU:HD21	1.88	0.54
1:A:102:THR:HG22	1:A:103:LYS:H	1.73	0.54
1:A:505:LEU:HD13	1:B:130:TYR:HB2	1.90	0.54
1:B:143:TYR:O	1:B:376:ARG:NH2	2.41	0.54
1:B:373:HIS:O	1:B:376:ARG:HG3	2.06	0.54
1:A:119:LEU:O	1:A:449:LYS:HE2	2.07	0.53
1:B:193:TYR:HA	1:B:223:VAL:CG1	2.39	0.53
1:A:447:VAL:O	1:A:451:ILE:HG13	2.08	0.53
1:B:367:GLY:HA2	1:B:404:LEU:HD23	1.91	0.53
1:B:121:ASN:HD22	1:B:121:ASN:N	2.04	0.53
1:B:256:ASN:ND2	1:B:257:SER:N	2.54	0.53
1:B:394:ILE:HA	1:B:397:GLN:HG2	1.89	0.53
1:A:331:THR:HG22	1:A:332:TRP:N	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:THR:HG22	1:B:103:LYS:H	1.73	0.53
1:A:121:ASN:N	1:A:121:ASN:HD22	2.06	0.52
1:A:414:ARG:HD2	1:B:150:ASP:OD2	2.09	0.52
1:A:354:THR:HG21	1:A:394:ILE:HD12	1.90	0.52
1:A:394:ILE:HA	1:A:397:GLN:HG2	1.91	0.52
1:B:119:LEU:O	1:B:449:LYS:HE2	2.10	0.52
1:A:388:LEU:HD11	1:A:426:ALA:HB2	1.92	0.52
1:B:256:ASN:HA	1:B:325:PHE:O	2.10	0.51
1:B:494:VAL:HA	3:B:25:HOH:O	2.10	0.51
1:B:454:LEU:O	1:B:458:ILE:HG13	2.09	0.51
1:B:217:GLY:HA2	1:B:222:LYS:HG2	1.92	0.51
1:B:385:ILE:HD11	1:B:424:LEU:HD12	1.93	0.51
1:A:256:ASN:ND2	1:A:257:SER:N	2.58	0.51
1:A:333:LYS:HD2	1:A:336:GLU:OE2	2.09	0.51
1:B:447:VAL:O	1:B:451:ILE:HG13	2.11	0.51
1:A:332:TRP:HE3	1:A:363:ILE:HD13	1.76	0.51
1:B:124:ASN:HD22	1:B:126:TYR:N	2.06	0.51
1:A:177:VAL:HG22	1:A:468:THR:HA	1.94	0.50
1:B:442:TYR:HB2	1:B:446:GLY:HA3	1.93	0.50
1:A:217:GLY:HA2	1:A:222:LYS:HG2	1.93	0.50
1:A:392:MET:HB2	1:A:393:PRO:HD3	1.94	0.50
1:B:116:LEU:CD2	1:B:131:LEU:HG	2.42	0.49
1:B:263:ASP:O	1:B:267:LYS:HG2	2.11	0.49
1:A:123:ILE:HD12	1:B:294:LYS:NZ	2.27	0.49
1:B:388:LEU:HD11	1:B:426:ALA:HB2	1.93	0.49
1:B:392:MET:HB2	1:B:393:PRO:HD3	1.94	0.49
1:A:199:LEU:HA	1:A:232:SER:OG	2.13	0.49
1:B:144:TYR:O	1:B:433:ARG:HD3	2.13	0.49
1:A:116:LEU:CD2	1:A:131:LEU:HG	2.42	0.49
1:A:163:ARG:HA	3:A:22:HOH:O	2.12	0.49
1:A:256:ASN:HD22	1:A:258:ASP:N	1.98	0.49
1:B:351:VAL:HG21	1:B:368:VAL:CG2	2.43	0.49
1:A:166:PHE:CE2	1:A:416:THR:HG22	2.48	0.48
1:A:385:ILE:HD11	1:A:424:LEU:HD12	1.94	0.48
1:B:332:TRP:HE3	1:B:363:ILE:HD13	1.78	0.48
1:B:253:LEU:HD11	1:B:262:THR:HG23	1.94	0.48
1:A:385:ILE:HG22	3:A:24:HOH:O	2.14	0.48
1:A:374:GLY:O	1:A:433:ARG:NH2	2.45	0.48
1:A:442:TYR:HB2	1:A:446:GLY:HA3	1.96	0.48
1:A:144:TYR:O	1:A:433:ARG:HD3	2.14	0.48
1:A:494:VAL:O	1:A:498:GLU:HB2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:218:GLN:HG3	1:B:219:GLY:H	1.79	0.47
1:B:338:LEU:O	1:B:342:THR:HG22	2.15	0.47
1:B:369:VAL:HG22	1:B:407:PHE:HB2	1.96	0.47
1:B:374:GLY:O	1:B:433:ARG:NH2	2.45	0.47
1:A:201:LYS:HE3	1:A:233:CYS:SG	2.55	0.47
1:B:379:ASP:O	1:B:380:PHE:HB2	2.15	0.47
1:B:112:LEU:HD13	1:B:135:THR:HB	1.97	0.47
1:A:217:GLY:HA3	1:A:247:GLN:OE1	2.15	0.47
1:B:119:LEU:HD13	1:B:442:TYR:CD2	2.51	0.46
1:B:141:TRP:HZ3	1:B:437:TYR:CE1	2.34	0.46
1:A:351:VAL:HG21	1:A:368:VAL:CG2	2.45	0.46
1:A:124:ASN:C	1:A:124:ASN:HD22	2.24	0.46
3:A:9:HOH:O	1:B:494:VAL:HG13	2.15	0.46
1:B:349:LYS:HA	1:B:369:VAL:HB	1.97	0.46
1:A:123:ILE:HG21	1:B:290:GLU:HG2	1.98	0.46
1:A:379:ASP:O	1:A:380:PHE:HB2	2.15	0.46
1:A:193:TYR:HA	1:A:223:VAL:CG1	2.46	0.45
1:B:177:VAL:HG22	1:B:468:THR:HA	1.97	0.45
1:B:256:ASN:ND2	1:B:257:SER:H	2.13	0.45
1:B:331:THR:HG22	1:B:332:TRP:H	1.81	0.45
1:A:157:ASN:HD21	1:A:372:ASN:HD21	1.65	0.45
1:B:217:GLY:HA3	1:B:247:GLN:OE1	2.15	0.45
1:A:103:LYS:NZ	1:A:103:LYS:HB3	2.32	0.45
1:A:112:LEU:HD13	1:A:135:THR:HB	1.98	0.45
1:A:166:PHE:HE2	1:A:416:THR:HG22	1.81	0.45
1:B:271:LYS:HB3	1:B:271:LYS:HE2	1.82	0.45
1:A:123:ILE:CG2	1:B:290:GLU:HG2	2.47	0.45
1:A:234:SER:O	1:A:238:ILE:HG13	2.17	0.45
2:B:571:FNS:H9	2:B:571:FNS:O4'	2.17	0.45
1:A:454:LEU:O	1:A:458:ILE:HG13	2.17	0.44
1:B:109:LYS:NZ	1:B:112:LEU:HD23	2.32	0.44
1:A:107:ALA:O	1:A:111:GLN:HG2	2.17	0.44
1:B:191:PRO:HB2	1:B:422:LEU:HD11	1.99	0.43
1:B:494:VAL:O	1:B:498:GLU:HB2	2.18	0.43
1:B:288:GLN:CG	1:B:293:MET:SD	3.07	0.43
1:A:109:LYS:NZ	1:A:112:LEU:HD23	2.32	0.43
1:A:266:VAL:HG13	1:A:277:LEU:HD11	2.00	0.43
1:B:157:ASN:ND2	1:B:372:ASN:HD21	2.15	0.43
1:B:256:ASN:O	1:B:259:ARG:HD3	2.19	0.43
1:A:256:ASN:ND2	1:A:257:SER:H	2.16	0.43
1:B:111:GLN:HB2	1:B:135:THR:HG22	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:GLN:O	1:A:115:LEU:HB2	2.19	0.43
1:B:292:ASP:O	1:B:295:LEU:HB2	2.18	0.43
1:A:338:LEU:O	1:A:342:THR:HG22	2.18	0.43
1:B:266:VAL:HG13	1:B:277:LEU:HD11	2.00	0.43
1:A:244:SER:O	1:A:247:GLN:HG2	2.19	0.42
1:B:214:ARG:HA	1:B:243:PRO:HD3	2.01	0.42
1:B:281:VAL:HG11	1:B:348:ILE:HG23	2.00	0.42
1:A:482:THR:OG1	1:B:486:ARG:NH1	2.52	0.42
1:B:111:GLN:O	1:B:115:LEU:HB2	2.19	0.42
1:A:349:LYS:HA	1:A:369:VAL:HB	2.01	0.42
1:A:222:LYS:HE3	1:A:244:SER:CB	2.49	0.42
1:B:234:SER:O	1:B:238:ILE:HG13	2.19	0.42
1:B:199:LEU:HA	1:B:232:SER:OG	2.19	0.42
1:B:141:TRP:CZ3	1:B:437:TYR:CE1	3.08	0.42
1:A:124:ASN:HD22	1:A:126:TYR:N	2.18	0.42
1:A:141:TRP:HZ3	1:A:437:TYR:CE1	2.38	0.42
1:B:330:LEU:HD23	1:B:330:LEU:HA	1.80	0.41
1:A:282:ASP:O	1:A:377:GLN:HG3	2.20	0.41
1:B:121:ASN:N	1:B:121:ASN:ND2	2.68	0.41
1:B:282:ASP:O	1:B:377:GLN:HG3	2.20	0.41
2:B:571:FNS:HO2'	2:B:571:FNS:C10	2.31	0.41
1:A:339:LYS:HE2	1:A:339:LYS:HB3	1.84	0.41
1:A:332:TRP:O	1:A:336:GLU:HG3	2.20	0.41
1:B:347:VAL:CG1	1:B:369:VAL:HG23	2.50	0.41
1:A:225:GLN:HE21	1:A:225:GLN:HB3	1.72	0.41
1:A:111:GLN:HB2	1:A:135:THR:HG22	2.03	0.41
1:A:271:LYS:HB3	1:A:271:LYS:HE2	1.82	0.41
1:A:353:ARG:NE	3:A:53:HOH:O	2.52	0.41
1:A:124:ASN:HD22	1:A:126:TYR:H	1.69	0.41
1:A:191:PRO:HB2	1:A:422:LEU:HD11	2.03	0.41
1:A:467:VAL:HG11	1:A:477:LEU:HD21	2.03	0.41
1:B:119:LEU:HD13	1:B:442:TYR:CG	2.55	0.40
1:B:378:LEU:HD22	1:B:379:ASP:O	2.21	0.40
1:B:124:ASN:HD22	1:B:124:ASN:C	2.28	0.40
1:A:258:ASP:OD1	1:A:260:LYS:HB3	2.21	0.40
1:A:333:LYS:O	1:A:337:GLU:HG2	2.22	0.40
1:B:424:LEU:HD23	1:B:424:LEU:HA	1.95	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	381/410 (93%)	353 (93%)	28 (7%)	0	100	100
1	B	381/410 (93%)	356 (93%)	24 (6%)	1 (0%)	36	56
All	All	762/820 (93%)	709 (93%)	52 (7%)	1 (0%)	48	71

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	400	LEU

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	334/353 (95%)	307 (92%)	27 (8%)	11	20
1	B	334/353 (95%)	308 (92%)	26 (8%)	11	21
All	All	668/706 (95%)	615 (92%)	53 (8%)	11	21

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	102	THR
1	A	103	LYS
1	A	124	ASN
1	A	125	LEU
1	A	131	LEU

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Mol	Chain	Res	Type
1	A	184	LEU
1	A	255	VAL
1	A	256	ASN
1	A	265	LEU
1	A	295	LEU
1	A	330	LEU
1	A	368	VAL
1	A	372	ASN
1	A	378	LEU
1	A	388	LEU
1	A	404	LEU
1	A	413	ARG
1	A	414	ARG
1	A	416	THR
1	A	436	LEU
1	A	464	LEU
1	A	476	ASP
1	A	483	LEU
1	A	490	VAL
1	A	494	VAL
1	A	509	GLU
1	A	510	ASP
1	B	103	LYS
1	B	124	ASN
1	B	125	LEU
1	B	131	LEU
1	B	184	LEU
1	B	255	VAL
1	B	256	ASN
1	B	265	LEU
1	B	295	LEU
1	B	330	LEU
1	B	368	VAL
1	B	372	ASN
1	B	378	LEU
1	B	388	LEU
1	B	404	LEU
1	B	413	ARG
1	B	414	ARG
1	B	416	THR
1	B	436	LEU
1	B	464	LEU

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Mol	Chain	Res	Type
1	B	476	ASP
1	B	483	LEU
1	B	490	VAL
1	B	494	VAL
1	B	509	GLU
1	B	510	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	121	ASN
1	A	124	ASN
1	A	154	HIS
1	A	218	GLN
1	A	225	GLN
1	A	249	GLN
1	A	256	ASN
1	A	268	ASN
1	A	372	ASN
1	A	397	GLN
1	B	121	ASN
1	B	124	ASN
1	B	218	GLN
1	B	225	GLN
1	B	249	GLN
1	B	256	ASN
1	B	268	ASN
1	B	372	ASN
1	B	397	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	$\# Z > 2$
2	FNS	A	570	-	32,37,37	1.77	5 (15%)	40,58,58	1.99	10 (25%)
2	FNS	B	571	-	32,37,37	1.70	4 (12%)	40,58,58	1.85	11 (27%)

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	FNS	A	570	-	-	2/18/24/24	0/3/3/3
2	FNS	B	571	-	-	2/18/24/24	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	570	FNS	C1'-N10	-6.54	1.32	1.47
2	B	571	FNS	C1'-N10	-6.07	1.33	1.47
2	A	570	FNS	C5'-C4'	-3.93	1.46	1.51
2	B	571	FNS	C5'-C4'	-3.88	1.46	1.51
2	B	571	FNS	C10-N1	2.93	1.39	1.33
2	A	570	FNS	C2-N3	2.41	1.44	1.39
2	A	570	FNS	C9-C8	-2.40	1.36	1.39
2	B	571	FNS	C5A-N5	-2.22	1.35	1.41
2	A	570	FNS	C4'-C3'	-2.07	1.49	1.53

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	570	FNS	C5'-C4'-C3'	-5.19	102.43	112.22
2	B	571	FNS	C5'-C4'-C3'	-4.72	103.32	112.22
2	A	570	FNS	C10-N1-C2	4.70	127.02	116.85
2	A	570	FNS	O2S-S-O3S	-4.63	108.46	119.69
2	B	571	FNS	O2S-S-O3S	-4.63	108.47	119.69
2	A	570	FNS	O2-C2-N3	4.03	126.31	118.58
2	B	571	FNS	O4'-C4'-C3'	3.53	117.52	109.25
2	A	570	FNS	N10-C10-N1	3.45	128.65	118.51
2	B	571	FNS	C5A-C9A-N10	-3.38	115.05	119.04
2	B	571	FNS	N10-C10-N1	3.33	128.31	118.51
2	A	570	FNS	O4'-C4'-C3'	3.32	117.03	109.25
2	B	571	FNS	C10-N1-C2	3.23	123.85	116.85
2	B	571	FNS	O2-C2-N3	3.14	124.61	118.58
2	B	571	FNS	O3P-P-O5'	-3.11	98.55	106.67
2	A	570	FNS	C5A-C9A-N10	-2.92	115.60	119.04
2	A	570	FNS	N3-C2-N1	-2.41	114.37	119.50
2	B	571	FNS	C9-C9A-N10	2.27	124.91	121.85
2	A	570	FNS	O3P-P-O5'	-2.25	100.81	106.67
2	B	571	FNS	N3-C2-N1	-2.11	115.01	119.50
2	A	570	FNS	C4'-C3'-C2'	-2.07	110.12	113.57
2	B	571	FNS	O3P-P-O2P	2.01	115.33	107.80

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	570	FNS	C5'-O5'-P-O1P
2	A	570	FNS	C4'-C5'-O5'-P
2	B	571	FNS	C4'-C5'-O5'-P
2	B	571	FNS	C5'-O5'-P-O1P

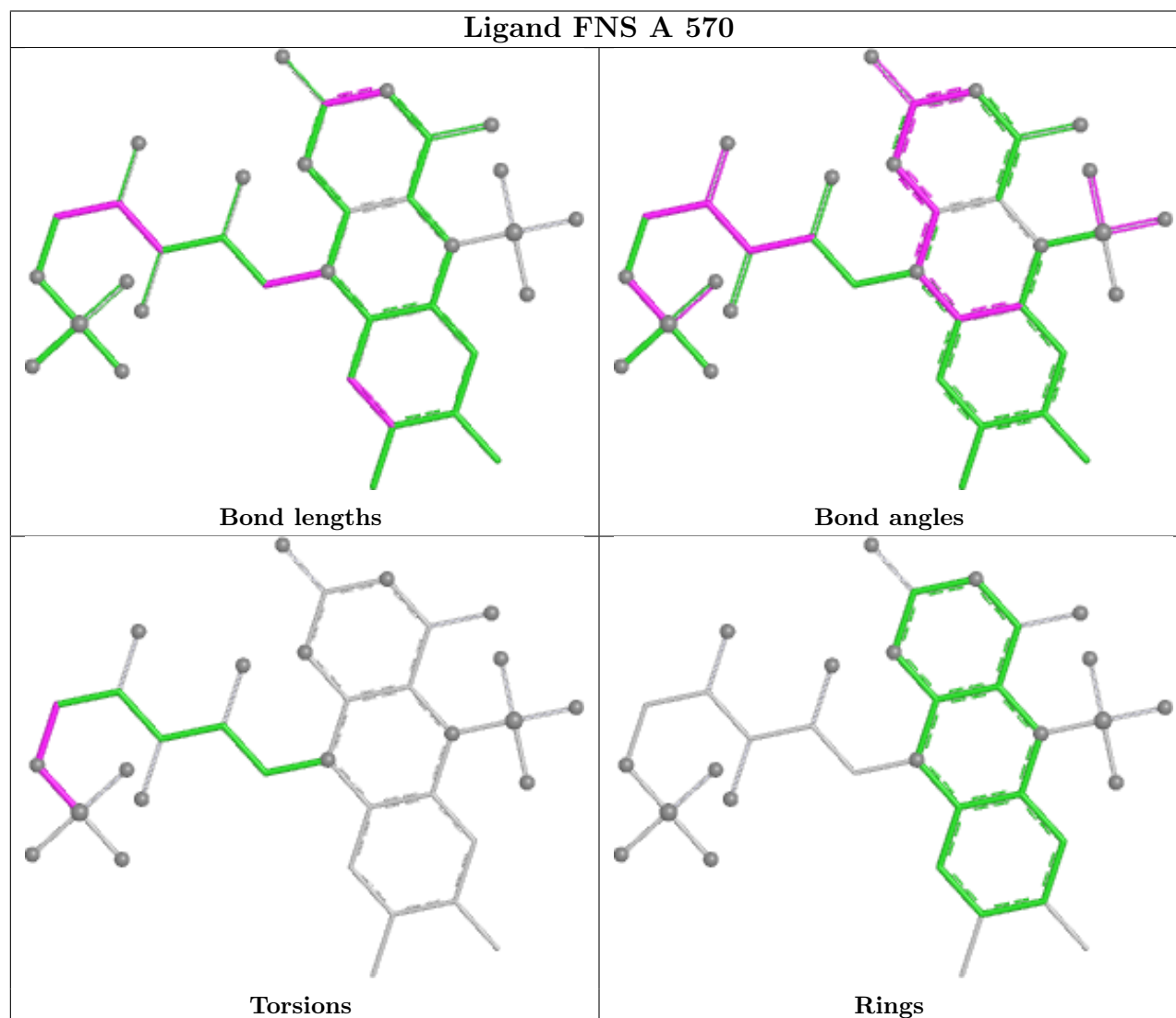
There are no ring outliers.

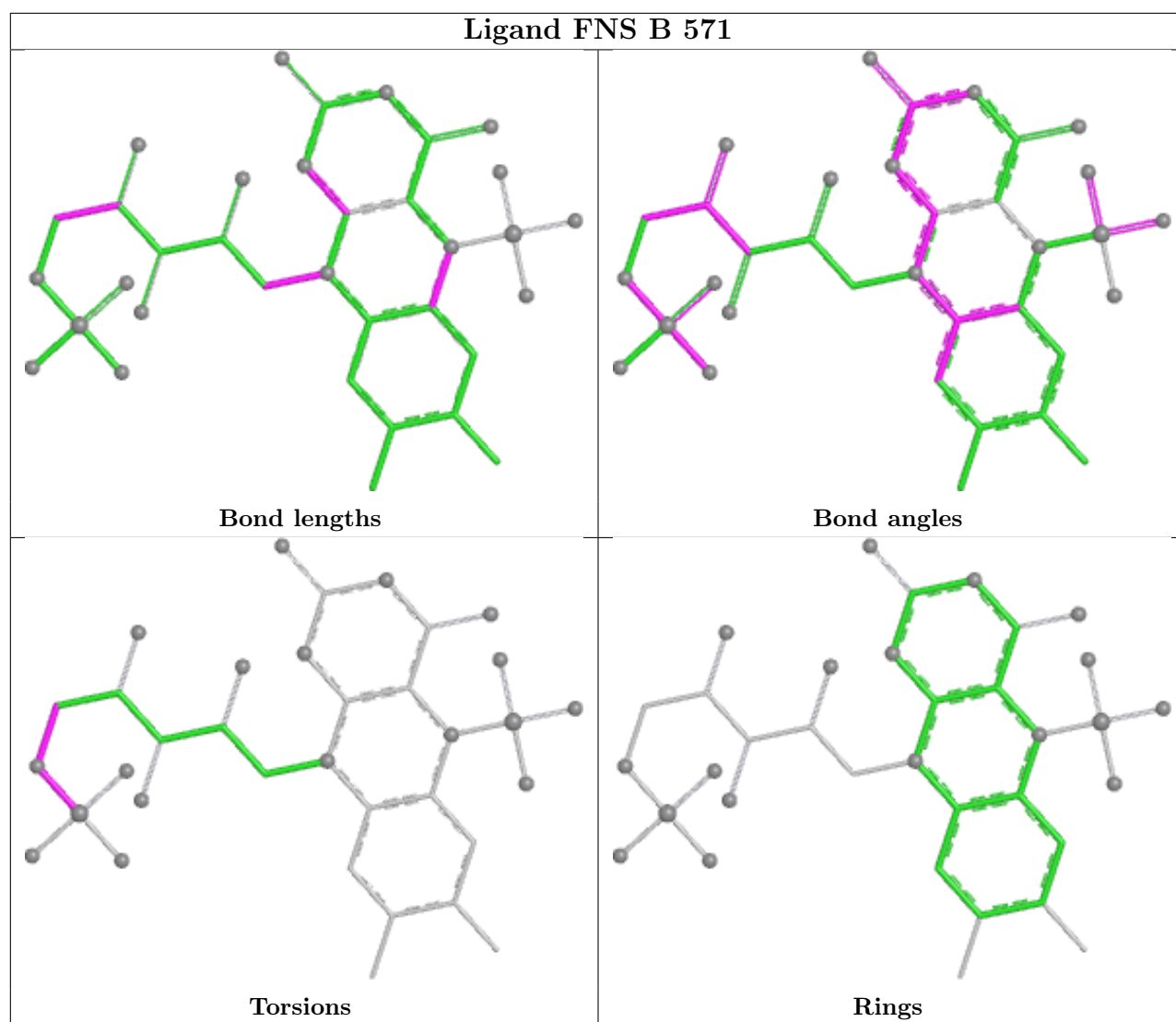
2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	570	FNS	1	0
2	B	571	FNS	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 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.