



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:10 AM UTC

PDB ID : 1Q9M / pdb_00001q9m
Title : Three dimensional structures of PDE4D in complex with roliprams and implication on inhibitor selectivity
Authors : Huai, Q.; Wang, H.; Sun, Y.; Kim, H.Y.; Liu, Y.; Ke, H.
Deposited on : 2003-08-25
Resolution : 2.30 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) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

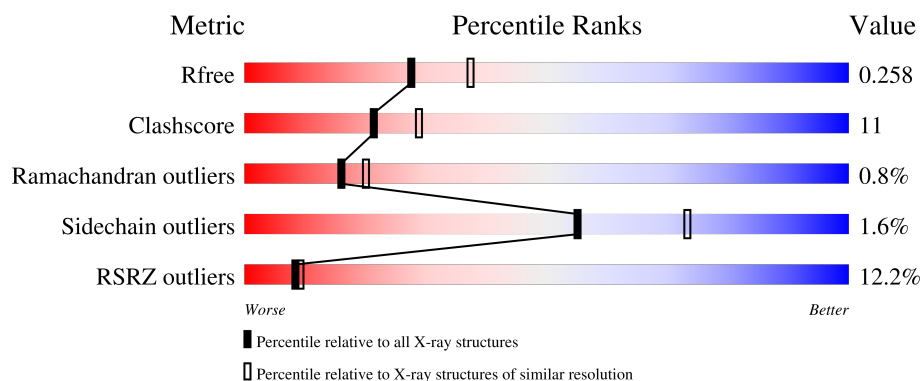
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	360	 8% 69% 22% • 7%
1	B	360	 20% 64% 26% • 9%
1	C	360	 11% 69% 20% • 9%
1	D	360	 5% 73% 19% • 7%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10972 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.

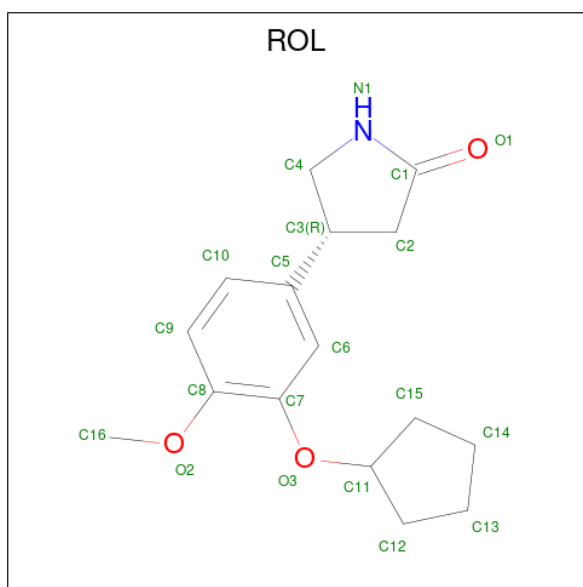
- Molecule 1 is a protein called cAMP-specific phosphodiesterase PDE4D2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	334	Total	C	N	O	S	0	0	0
			2704	1712	463	515	14			
1	B	327	Total	C	N	O	S	0	0	0
			2647	1673	452	508	14			
1	C	329	Total	C	N	O	S	0	0	0
			2663	1684	455	510	14			
1	D	334	Total	C	N	O	S	0	0	0
			2704	1712	463	515	14			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Zn	0	0
			2	2		
2	B	2	Total	Zn	0	0
			2	2		
2	C	2	Total	Zn	0	0
			2	2		
2	D	2	Total	Zn	0	0
			2	2		

- Molecule 3 is ROLIPRAM (CCD ID: ROL) (formula: C₁₆H₂₁NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			20	16	1	3		
3	B	1	Total	C	N	O	0	0
			20	16	1	3		
3	C	1	Total	C	N	O	0	0
			20	16	1	3		
3	D	1	Total	C	N	O	0	0
			20	16	1	3		

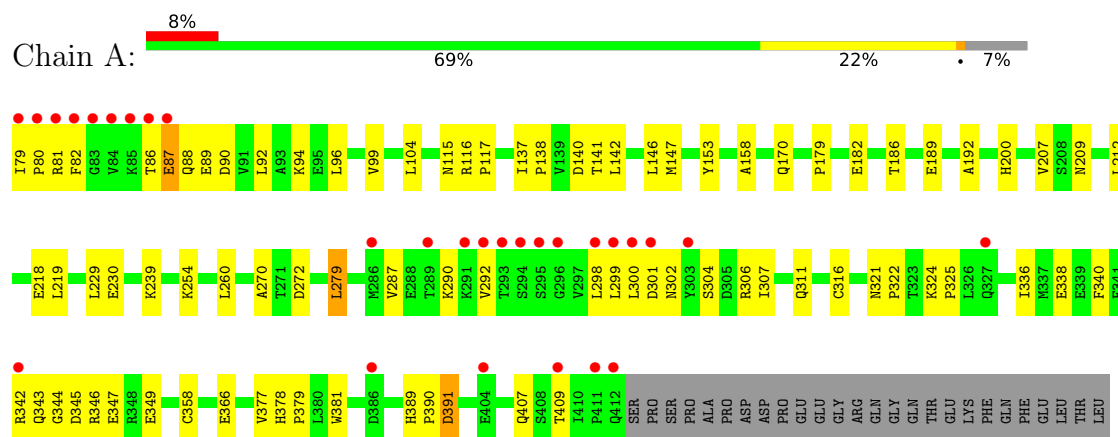
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	52	Total	O	0	0
			52	52		
4	B	36	Total	O	0	0
			36	36		
4	C	32	Total	O	0	0
			32	32		
4	D	46	Total	O	0	0
			46	46		

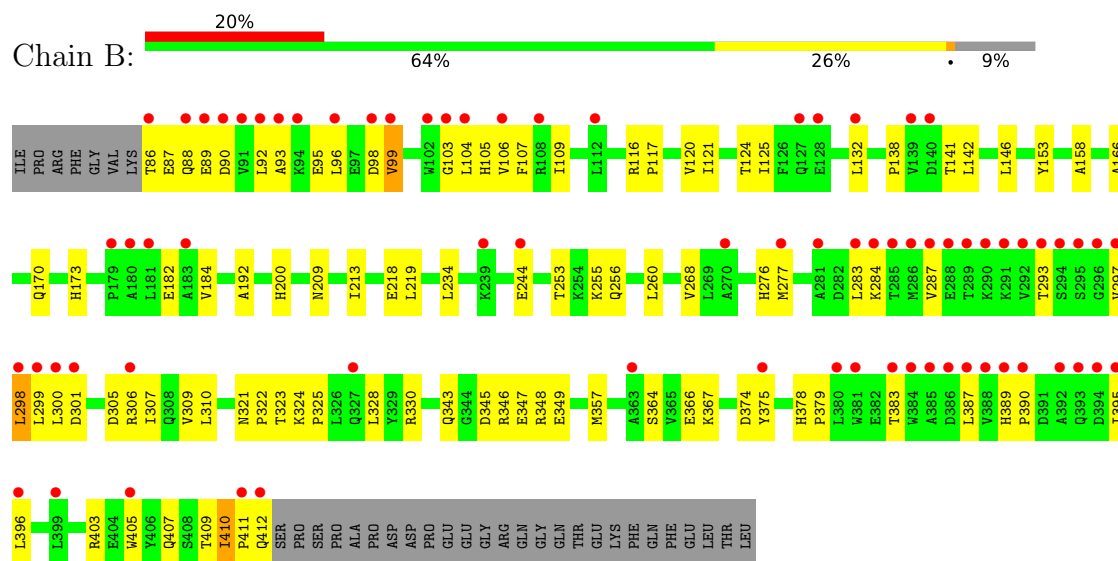
3 Residue-property plots [i](#)

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 and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: cAMP-specific phosphodiesterase PDE4D2

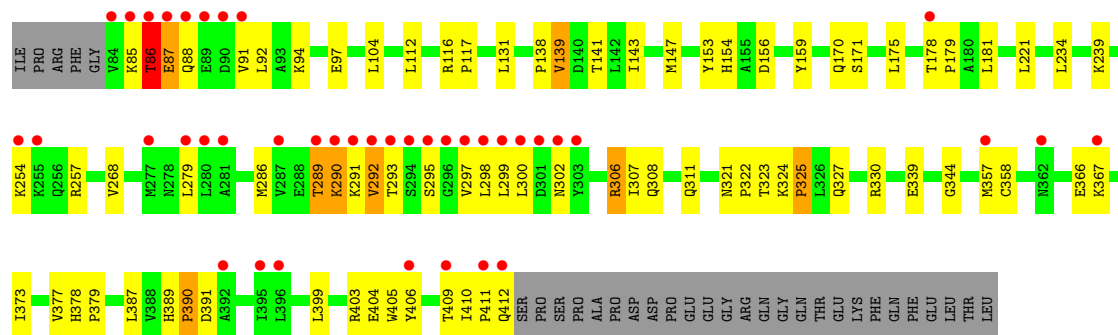


• Molecule 1: cAMP-specific phosphodiesterase PDE4D2

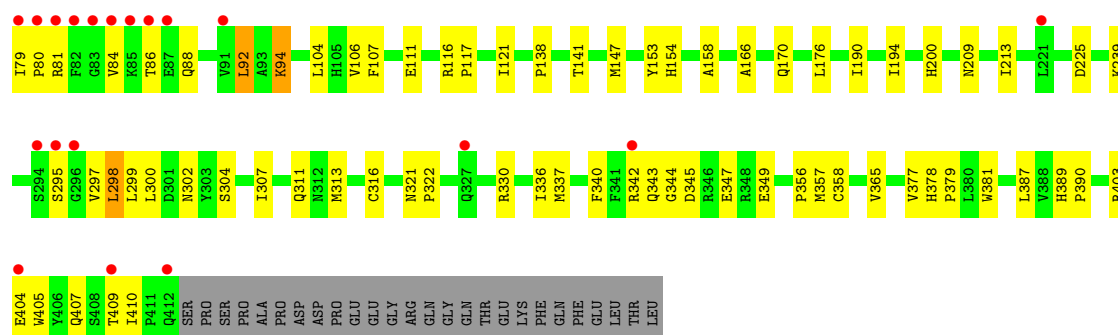


• Molecule 1: cAMP-specific phosphodiesterase PDE4D2





● Molecule 1: cAMP-specific phosphodiesterase PDE4D2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	99.27Å 112.49Å 160.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	99.00 – 2.30 92.20 – 2.30	Depositor EDS
% Data completeness (in resolution range)	93.9 (99.00-2.30) 93.9 (92.20-2.30)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.70 (at 2.29Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.224 , 0.259 0.223 , 0.258	Depositor DCC
R_{free} test set	7857 reflections (10.09%)	wwPDB-VP
Wilson B-factor (Å ²)	31.7	Xtriage
Anisotropy	0.077	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 32.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10972	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.42	0/2760	0.86	7/3749 (0.2%)
1	B	0.39	0/2701	0.87	7/3670 (0.2%)
1	C	0.39	0/2717	0.86	5/3691 (0.1%)
1	D	0.43	0/2760	0.90	9/3749 (0.2%)
All	All	0.41	0/10938	0.87	28/14859 (0.2%)

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	158	ALA	N-CA-C	7.52	119.48	111.28
1	D	153	TYR	N-CA-C	-6.92	98.33	109.96
1	C	153	TYR	N-CA-C	-6.44	99.14	109.96
1	A	377	VAL	N-CA-C	6.42	116.58	110.42
1	D	158	ALA	N-CA-C	6.26	118.10	111.28
1	B	153	TYR	N-CA-C	-6.08	99.32	109.24
1	D	81	ARG	N-CA-C	6.06	117.88	111.28
1	A	200	HIS	N-CA-C	6.00	120.14	112.34
1	D	377	VAL	N-CA-C	5.98	116.11	110.30
1	D	200	HIS	N-CA-C	5.85	119.95	112.34
1	B	374	ASP	N-CA-C	5.85	117.46	111.14
1	D	88	GLN	N-CA-C	-5.79	105.80	112.92
1	B	90	ASP	N-CA-C	-5.78	105.07	111.71
1	A	153	TYR	N-CA-C	-5.76	100.28	109.96
1	A	391	ASP	N-CA-C	5.75	118.29	111.33
1	C	131	LEU	N-CA-C	5.74	117.53	111.28
1	D	295	SER	N-CA-C	-5.54	102.56	110.59
1	D	297	VAL	N-CA-C	5.45	117.18	108.89
1	B	158	ALA	N-CA-C	5.37	117.13	111.28
1	C	139	VAL	N-CA-C	5.25	115.87	110.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	410	ILE	CA-C-N	5.21	126.36	119.84
1	B	410	ILE	C-N-CA	5.21	126.36	119.84
1	B	200	HIS	N-CA-C	5.11	118.61	112.38
1	A	137	ILE	CA-C-N	-5.06	114.55	119.76
1	A	137	ILE	C-N-CA	-5.06	114.55	119.76
1	C	178	THR	CA-C-N	5.05	125.08	119.32
1	C	178	THR	C-N-CA	5.05	125.08	119.32
1	D	106	VAL	N-CA-C	5.04	115.75	110.62

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	2704	0	2664	56	0
1	B	2647	0	2599	77	0
1	C	2663	0	2621	63	0
1	D	2704	0	2664	44	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	20	0	21	2	0
3	B	20	0	21	3	0
3	C	20	0	21	1	0
3	D	20	0	21	1	0
4	A	52	0	0	0	0
4	B	36	0	0	1	0
4	C	32	0	0	1	0
4	D	46	0	0	0	0
All	All	10972	0	10632	234	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 11.

All (234) close contacts within the same asymmetric unit are listed below, sorted by their clash

magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:255:LYS:H	1:B:255:LYS:HD2	1.29	0.95
1:C:293:THR:HG23	1:C:295:SER:H	1.46	0.80
1:B:364:SER:HB3	1:B:367:LYS:HD3	1.64	0.79
1:B:330:ARG:HD3	1:B:405:TRP:CH2	2.19	0.77
1:B:244:GLU:HG3	1:C:254:LYS:NZ	2.02	0.74
1:D:79:ILE:N	1:D:80:PRO:HD2	2.03	0.73
1:C:306:ARG:HB3	1:C:306:ARG:HH11	1.55	0.72
1:B:98:ASP:HB2	1:B:104:LEU:HD13	1.72	0.71
1:B:255:LYS:H	1:B:255:LYS:CD	2.04	0.69
1:C:293:THR:HG22	1:C:297:VAL:H	1.57	0.69
1:A:87:GLU:HG3	1:A:88:GLN:N	2.09	0.68
1:A:349:GLU:HB3	1:C:147:MET:HE1	1.74	0.68
1:B:96:LEU:O	1:B:99:VAL:HG23	1.94	0.68
1:D:302:ASN:H	1:D:302:ASN:ND2	1.90	0.67
1:A:96:LEU:O	1:A:99:VAL:HG23	1.94	0.67
1:C:85:LYS:O	1:C:86:THR:HG23	1.94	0.67
1:A:116:ARG:HE	1:A:147:MET:HE2	1.60	0.66
1:B:300:LEU:HB3	1:B:306:ARG:HG2	1.78	0.66
1:B:283:LEU:O	1:B:287:VAL:HG23	1.95	0.66
1:B:324:LYS:HB3	1:B:325:PRO:HD2	1.77	0.65
1:B:378:HIS:HB3	1:B:379:PRO:HD3	1.78	0.65
1:D:79:ILE:N	1:D:80:PRO:CD	2.60	0.65
1:D:138:PRO:HG2	1:D:141:THR:OG1	1.97	0.64
1:B:411:PRO:O	1:B:412:GLN:HB2	1.97	0.64
1:C:254:LYS:HG2	1:C:257:ARG:HH22	1.62	0.64
1:B:366:GLU:HG2	1:B:409:THR:HG22	1.79	0.63
1:A:344:GLY:HA3	1:A:358:CYS:O	1.99	0.63
1:C:181:LEU:HD21	1:C:298:LEU:HD12	1.80	0.63
1:D:407:GLN:NE2	1:D:410:ILE:HD12	2.13	0.63
1:B:389:HIS:CE1	1:B:390:PRO:HB3	2.34	0.62
1:D:79:ILE:N	1:D:79:ILE:HD12	2.15	0.62
1:A:116:ARG:NE	1:A:147:MET:HE2	2.14	0.62
1:B:284:LYS:HE2	1:B:383:THR:OG1	2.00	0.62
1:A:79:ILE:N	1:A:80:PRO:HD2	2.14	0.61
1:C:321:ASN:HB2	1:C:322:PRO:HD3	1.82	0.61
1:C:86:THR:O	1:C:88:GLN:N	2.34	0.61
1:D:337:MET:HE2	1:D:365:VAL:HG22	1.81	0.61
1:B:293:THR:HG22	1:B:299:LEU:HD23	1.82	0.61
1:B:321:ASN:HB2	1:B:322:PRO:HD3	1.83	0.61
1:D:302:ASN:OD1	1:D:304:SER:HB3	2.01	0.60
1:A:321:ASN:HB2	1:A:322:PRO:HD3	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:403:ARG:HD2	1:C:404:GLU:OE1	2.01	0.59
1:D:344:GLY:HA3	1:D:358:CYS:O	2.02	0.59
1:A:179:PRO:HD2	1:A:391:ASP:OD2	2.03	0.59
1:D:321:ASN:HB2	1:D:322:PRO:HD3	1.83	0.59
1:C:291:LYS:O	1:C:299:LEU:HB3	2.02	0.58
1:D:298:LEU:HD23	1:D:299:LEU:H	1.68	0.58
1:D:403:ARG:HD2	1:D:404:GLU:OE1	2.03	0.58
1:A:306:ARG:HH11	1:A:306:ARG:HG3	1.69	0.57
1:B:184:VAL:HG11	1:B:300:LEU:HD12	1.86	0.57
1:D:307:ILE:O	1:D:311:GLN:HG3	2.04	0.57
1:C:290:LYS:HG2	1:C:292:VAL:HG13	1.86	0.57
1:B:409:THR:HG22	1:B:409:THR:O	2.03	0.57
1:A:389:HIS:CE1	1:A:390:PRO:HB3	2.40	0.56
1:B:298:LEU:HD23	1:B:299:LEU:N	2.21	0.56
1:B:218:GLU:OE2	1:D:239:LYS:HE2	2.05	0.56
1:B:244:GLU:HG3	1:C:254:LYS:HZ2	1.71	0.56
1:C:254:LYS:CG	1:C:257:ARG:HH22	2.20	0.55
1:A:186:THR:OG1	1:A:189:GLU:HG3	2.06	0.55
1:B:192:ALA:HB2	1:B:260:LEU:HD12	1.87	0.55
1:A:307:ILE:O	1:A:311:GLN:HG3	2.06	0.55
1:B:293:THR:CG2	1:B:299:LEU:HD23	2.37	0.55
1:B:325:PRO:HG2	1:B:328:LEU:HD12	1.89	0.55
1:C:221:LEU:O	1:C:221:LEU:HD23	2.07	0.55
1:D:298:LEU:HD11	1:D:387:LEU:HD11	1.89	0.55
1:A:81:ARG:O	1:A:81:ARG:HG3	2.06	0.54
1:B:345:ASP:O	1:B:349:GLU:HG3	2.07	0.54
1:C:181:LEU:CD2	1:C:298:LEU:HD12	2.38	0.54
1:B:349:GLU:HB3	1:D:147:MET:HE1	1.88	0.54
1:D:84:VAL:HG11	1:D:116:ARG:HD2	1.90	0.54
1:C:330:ARG:HD3	1:C:405:TRP:CH2	2.43	0.53
1:C:293:THR:HG22	1:C:297:VAL:N	2.22	0.53
1:B:170:GLN:O	1:B:173:HIS:HB3	2.09	0.53
1:B:234:LEU:HD21	1:B:268:VAL:HB	1.90	0.53
1:C:307:ILE:O	1:C:311:GLN:HG3	2.09	0.53
1:C:234:LEU:HD21	1:C:268:VAL:HB	1.89	0.53
1:D:107:PHE:O	1:D:111:GLU:HG3	2.08	0.53
1:B:300:LEU:HD21	1:B:309:VAL:HG21	1.91	0.53
1:B:305:ASP:O	1:B:309:VAL:HG23	2.09	0.53
1:C:290:LYS:HE2	1:C:292:VAL:HG11	1.90	0.52
1:C:289:THR:O	1:C:289:THR:HG22	2.08	0.52
1:D:298:LEU:HD23	1:D:299:LEU:N	2.23	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:ARG:O	1:A:82:PHE:HB2	2.09	0.52
1:C:139:VAL:HG23	4:C:531:HOH:O	2.08	0.52
1:D:345:ASP:O	1:D:349:GLU:HG3	2.10	0.52
1:A:306:ARG:HG3	1:A:306:ARG:NH1	2.24	0.51
1:B:184:VAL:CG1	1:B:300:LEU:HD12	2.40	0.51
1:C:290:LYS:HE2	1:C:292:VAL:CG1	2.41	0.51
1:A:79:ILE:N	1:A:80:PRO:CD	2.74	0.51
1:A:138:PRO:HG2	1:A:141:THR:OG1	2.11	0.51
1:C:357:MET:SD	3:C:507:ROL:H131	2.51	0.51
1:B:87:GLU:HG3	1:B:88:GLN:HG3	1.93	0.51
1:D:104:LEU:HD22	1:D:170:GLN:HG3	1.91	0.51
1:B:138:PRO:HG2	1:B:141:THR:OG1	2.11	0.51
1:C:302:ASN:O	1:C:306:ARG:HG3	2.11	0.51
1:D:121:ILE:HD12	1:D:166:ALA:HB1	1.93	0.51
1:B:346:ARG:HD2	4:B:509:HOH:O	2.11	0.50
1:D:86:THR:HG21	1:D:92:LEU:HD12	1.94	0.50
1:B:99:VAL:HG21	1:B:124:THR:HG21	1.92	0.50
1:C:254:LYS:HG2	1:C:257:ARG:NH2	2.27	0.50
1:A:343:GLN:O	1:A:347:GLU:HG3	2.12	0.50
1:C:298:LEU:HD11	1:C:387:LEU:HG	1.94	0.49
1:A:292:VAL:HG12	1:A:298:LEU:HA	1.94	0.49
1:A:116:ARG:N	1:A:117:PRO:CD	2.75	0.49
1:C:389:HIS:CE1	1:C:390:PRO:HB3	2.47	0.49
1:D:79:ILE:O	1:D:79:ILE:HG22	2.11	0.49
1:C:221:LEU:HD23	1:C:221:LEU:C	2.37	0.49
1:C:293:THR:HG23	1:C:295:SER:N	2.22	0.49
1:A:179:PRO:HA	1:A:182:GLU:HG3	1.93	0.48
1:B:116:ARG:N	1:B:117:PRO:CD	2.77	0.48
1:B:375:TYR:CD1	1:B:375:TYR:N	2.81	0.48
1:A:212:LEU:HD11	1:A:229:LEU:HD21	1.95	0.48
1:C:291:LYS:HE3	1:C:299:LEU:O	2.13	0.48
1:C:324:LYS:O	1:C:325:PRO:C	2.56	0.48
1:C:411:PRO:O	1:C:412:GLN:HG3	2.12	0.48
1:A:142:LEU:O	1:A:146:LEU:HG	2.13	0.48
1:B:403:ARG:HG2	1:B:403:ARG:HH11	1.77	0.48
1:C:378:HIS:HB3	1:C:379:PRO:HD3	1.96	0.48
1:A:86:THR:HG22	1:A:87:GLU:N	2.28	0.48
1:A:192:ALA:HB2	1:A:260:LEU:HD12	1.95	0.48
1:A:207:VAL:HG11	1:A:346:ARG:NH2	2.28	0.48
1:B:121:ILE:HD12	1:B:166:ALA:HB1	1.96	0.48
1:C:366:GLU:HG2	1:C:409:THR:HG22	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:GLU:HG2	1:A:272:ASP:HB2	1.95	0.48
1:A:409:THR:HG22	1:A:409:THR:O	2.12	0.48
1:D:116:ARG:N	1:D:117:PRO:CD	2.77	0.48
1:B:411:PRO:O	1:B:412:GLN:CB	2.61	0.48
1:B:104:LEU:HD11	1:B:109:ILE:HD11	1.95	0.47
1:A:378:HIS:HB3	1:A:379:PRO:HD3	1.96	0.47
1:B:283:LEU:HD11	1:B:387:LEU:HD22	1.96	0.47
1:B:299:LEU:C	1:B:299:LEU:HD13	2.39	0.47
1:A:79:ILE:HG23	1:A:89:GLU:CD	2.40	0.47
1:A:340:PHE:CE2	3:A:503:ROL:H151	2.50	0.47
1:A:349:GLU:HB3	1:C:147:MET:CE	2.42	0.46
1:B:357:MET:SD	3:B:505:ROL:H131	2.56	0.46
1:C:87:GLU:HG3	1:C:88:GLN:H	1.80	0.46
1:B:244:GLU:HG3	1:C:254:LYS:HZ1	1.78	0.46
1:C:330:ARG:HD3	1:C:405:TRP:CZ3	2.50	0.46
1:D:336:ILE:HD11	1:D:340:PHE:CE1	2.51	0.46
1:B:132:LEU:H	1:B:132:LEU:HD12	1.81	0.45
1:A:298:LEU:HD21	1:A:300:LEU:HD21	1.98	0.45
1:B:182:GLU:O	1:B:297:VAL:HG11	2.16	0.45
1:B:330:ARG:HD3	1:B:405:TRP:CZ3	2.51	0.45
1:B:253:THR:OG1	1:B:256:GLN:HG3	2.17	0.45
1:C:154:HIS:HB3	1:C:156:ASP:OD1	2.16	0.45
1:C:344:GLY:HA3	1:C:358:CYS:O	2.17	0.45
1:D:299:LEU:O	1:D:300:LEU:HD23	2.16	0.45
1:A:79:ILE:HG13	1:A:89:GLU:OE2	2.17	0.45
1:B:104:LEU:HD12	1:B:105:HIS:H	1.81	0.45
1:B:298:LEU:HD23	1:B:299:LEU:H	1.82	0.45
1:A:209:ASN:OD1	1:A:229:LEU:HG	2.17	0.45
1:B:345:ASP:OD1	1:B:348:ARG:NH2	2.49	0.45
1:A:104:LEU:HD22	1:A:170:GLN:HG3	1.98	0.45
1:C:406:TYR:O	1:C:409:THR:HB	2.17	0.45
1:D:209:ASN:O	1:D:213:ILE:HG13	2.16	0.45
1:D:94:LYS:HB2	1:D:94:LYS:NZ	2.32	0.44
1:B:367:LYS:HE3	1:B:410:ILE:HG23	1.99	0.44
1:C:116:ARG:N	1:C:117:PRO:CD	2.80	0.44
1:D:330:ARG:HD3	1:D:405:TRP:CH2	2.53	0.44
1:C:91:VAL:HG12	1:C:112:LEU:HD11	2.00	0.44
1:D:302:ASN:H	1:D:302:ASN:HD22	1.61	0.44
3:D:509:ROL:H61	3:D:509:ROL:H21	1.78	0.44
1:B:284:LYS:HA	1:B:287:VAL:HG23	2.00	0.44
1:D:176:LEU:HG	1:D:313:MET:SD	2.58	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:ASP:O	1:A:94:LYS:HG3	2.17	0.44
1:C:143:ILE:O	1:C:147:MET:HG3	2.19	0.43
1:C:104:LEU:HD22	1:C:170:GLN:HG3	2.00	0.43
1:C:306:ARG:HH11	1:C:306:ARG:CB	2.26	0.43
1:B:89:GLU:HG2	1:B:93:ALA:HB2	1.99	0.43
1:C:300:LEU:HD13	1:C:306:ARG:HA	1.99	0.43
1:A:207:VAL:CG1	1:A:346:ARG:HH21	2.31	0.43
1:D:316:CYS:HB3	1:D:381:TRP:CZ2	2.53	0.43
1:A:239:LYS:NZ	1:D:239:LYS:NZ	2.66	0.43
1:B:96:LEU:HD11	1:B:120:VAL:CG1	2.49	0.43
1:D:154:HIS:CD2	1:D:154:HIS:N	2.86	0.43
1:A:336:ILE:HD11	1:A:340:PHE:CE1	2.54	0.42
1:A:366:GLU:HG2	1:A:409:THR:HG22	2.01	0.42
1:B:255:LYS:HD2	1:B:255:LYS:N	2.11	0.42
1:D:356:PRO:O	1:D:357:MET:HB2	2.19	0.42
1:B:142:LEU:O	1:B:146:LEU:HG	2.19	0.42
1:D:343:GLN:O	1:D:347:GLU:HG3	2.18	0.42
1:A:270:ALA:HB1	1:A:279:LEU:HD11	2.02	0.42
1:A:115:ASN:C	1:A:117:PRO:HD2	2.44	0.42
1:B:284:LYS:HG2	1:B:383:THR:CG2	2.49	0.42
1:A:86:THR:HG22	1:A:88:GLN:H	1.84	0.42
1:A:140:ASP:OD2	1:A:140:ASP:N	2.49	0.42
1:A:302:ASN:HD21	1:A:304:SER:HB2	1.84	0.42
1:C:138:PRO:HG2	1:C:141:THR:OG1	2.20	0.42
1:C:411:PRO:C	1:C:412:GLN:HG3	2.45	0.42
1:B:125:ILE:HD13	1:B:173:HIS:HB2	2.02	0.42
1:B:310:LEU:HD23	1:B:310:LEU:HA	1.94	0.42
1:B:396:LEU:HD23	1:B:396:LEU:O	2.20	0.42
1:A:316:CYS:HB3	1:A:381:TRP:CZ2	2.55	0.42
1:B:276:HIS:CD2	1:B:277:MET:HE2	2.54	0.42
1:B:389:HIS:ND1	1:B:390:PRO:HB3	2.35	0.42
1:D:378:HIS:HB3	1:D:379:PRO:HD3	2.01	0.42
1:B:92:LEU:O	1:B:95:GLU:N	2.53	0.42
3:B:505:ROL:H61	3:B:505:ROL:H111	1.69	0.42
1:A:287:VAL:O	1:A:287:VAL:HG12	2.19	0.42
1:C:94:LYS:O	1:C:97:GLU:HB2	2.20	0.42
3:A:503:ROL:H61	3:A:503:ROL:H111	1.59	0.41
1:B:323:THR:HB	1:B:395:ILE:HG23	2.01	0.41
1:A:324:LYS:O	1:A:325:PRO:C	2.63	0.41
1:B:86:THR:C	1:B:88:GLN:N	2.77	0.41
1:C:179:PRO:HD2	1:C:391:ASP:OD2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:106:VAL:HG13	1:B:107:PHE:N	2.35	0.41
1:D:213:ILE:HD13	1:D:225:ASP:HB3	2.01	0.41
1:D:299:LEU:HD23	1:D:300:LEU:N	2.35	0.41
1:B:409:THR:O	1:B:409:THR:CG2	2.67	0.41
1:C:171:SER:O	1:C:175:LEU:HG	2.21	0.41
1:D:389:HIS:HA	1:D:390:PRO:HA	1.82	0.41
1:C:367:LYS:HE2	1:C:410:ILE:HG23	2.02	0.41
3:B:505:ROL:H61	3:B:505:ROL:H21	1.79	0.41
1:A:207:VAL:CG1	1:A:346:ARG:NH2	2.84	0.41
1:A:218:GLU:HB3	1:C:239:LYS:HE3	2.03	0.41
1:A:299:LEU:C	1:A:299:LEU:HD23	2.45	0.41
1:B:367:LYS:HG3	1:B:410:ILE:HG23	2.02	0.41
1:C:323:THR:HG22	1:C:399:LEU:HD13	2.02	0.41
1:A:338:GLU:O	1:A:342:ARG:HG3	2.20	0.41
1:B:209:ASN:O	1:B:213:ILE:HG13	2.20	0.41
1:C:159:TYR:HB3	1:C:339:GLU:OE1	2.20	0.41
1:C:373:ILE:HA	1:C:377:VAL:HB	2.03	0.41
1:B:307:ILE:HD12	1:B:307:ILE:HA	1.96	0.41
1:B:92:LEU:CD2	1:B:96:LEU:HG	2.51	0.40
1:B:343:GLN:O	1:B:347:GLU:HG3	2.22	0.40
1:C:254:LYS:CG	1:C:257:ARG:NH2	2.84	0.40
1:A:219:LEU:HD12	1:A:219:LEU:HA	1.92	0.40
1:A:345:ASP:O	1:A:349:GLU:HG3	2.20	0.40
1:C:409:THR:HG22	1:C:409:THR:O	2.20	0.40
1:D:190:ILE:O	1:D:194:ILE:HG13	2.22	0.40
1:B:349:GLU:HB3	1:D:147:MET:CE	2.51	0.40
1:C:286:MET:HE3	1:C:308:GLN:OE1	2.20	0.40
1:B:389:HIS:HA	1:B:390:PRO:HA	1.84	0.40
1:D:409:THR:O	1:D:409:THR:HG22	2.21	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	332/360 (92%)	315 (95%)	15 (4%)	2 (1%)	21	27
1	B	325/360 (90%)	302 (93%)	20 (6%)	3 (1%)	14	17
1	C	327/360 (91%)	301 (92%)	21 (6%)	5 (2%)	8	8
1	D	332/360 (92%)	319 (96%)	13 (4%)	0	100	100
All	All	1316/1440 (91%)	1237 (94%)	69 (5%)	10 (1%)	16	20

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	301	ASP
1	C	86	THR
1	C	87	GLU
1	C	289	THR
1	C	290	LYS
1	A	290	LYS
1	A	301	ASP
1	B	103	GLY
1	C	292	VAL
1	B	99	VAL

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	305/328 (93%)	300 (98%)	5 (2%)	55	73
1	B	299/328 (91%)	296 (99%)	3 (1%)	68	82
1	C	301/328 (92%)	294 (98%)	7 (2%)	44	63
1	D	305/328 (93%)	301 (99%)	4 (1%)	61	77
All	All	1210/1312 (92%)	1191 (98%)	19 (2%)	55	73

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	87	GLU

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Mol	Chain	Res	Type
1	A	92	LEU
1	A	254	LYS
1	A	279	LEU
1	A	407	GLN
1	B	219	LEU
1	B	298	LEU
1	B	407	GLN
1	C	86	THR
1	C	92	LEU
1	C	279	LEU
1	C	306	ARG
1	C	325	PRO
1	C	327	GLN
1	C	390	PRO
1	D	92	LEU
1	D	94	LYS
1	D	298	LEU
1	D	342	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	123	HIS
1	A	245	ASN
1	A	250	GLN
1	A	258	GLN
1	A	369	GLN
1	A	407	GLN
1	B	170	GLN
1	B	210	GLN
1	B	245	ASN
1	B	378	HIS
1	C	123	HIS
1	C	210	GLN
1	C	242	GLN
1	C	245	ASN
1	C	250	GLN
1	C	402	ASN
1	C	407	GLN
1	D	127	GLN
1	D	210	GLN
1	D	245	ASN

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Mol	Chain	Res	Type
1	D	250	GLN
1	D	258	GLN
1	D	302	ASN
1	D	321	ASN
1	D	327	GLN
1	D	361	HIS
1	D	369	GLN
1	D	407	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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](#)

Of 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 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	$\# Z > 2$
3	ROL	B	505	-	22,22,22	0.60	0	30,30,30	1.48	4 (13%)
3	ROL	D	509	-	22,22,22	0.66	0	30,30,30	1.58	6 (20%)
3	ROL	A	503	-	22,22,22	0.76	0	30,30,30	1.47	6 (20%)
3	ROL	C	507	-	22,22,22	0.62	0	30,30,30	1.44	4 (13%)

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	ROL	B	505	-	-	2/10/26/26	0/3/3/3
3	ROL	D	509	-	-	2/10/26/26	0/3/3/3
3	ROL	A	503	-	-	4/10/26/26	0/3/3/3
3	ROL	C	507	-	-	2/10/26/26	0/3/3/3

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	509	ROL	C3-C2-C1	3.97	110.13	104.96
3	B	505	ROL	C3-C2-C1	3.71	109.80	104.96
3	A	503	ROL	C3-C2-C1	3.70	109.79	104.96
3	C	507	ROL	C3-C2-C1	3.65	109.72	104.96
3	D	509	ROL	C3-C4-N1	3.14	105.70	103.09
3	B	505	ROL	C16-O2-C8	-3.11	112.95	117.51
3	A	503	ROL	C16-O2-C8	-3.10	112.97	117.51
3	C	507	ROL	C4-N1-C1	3.00	117.04	114.13
3	D	509	ROL	C16-O2-C8	-2.95	113.18	117.51
3	D	509	ROL	C2-C1-N1	-2.93	105.89	108.68
3	B	505	ROL	C4-N1-C1	2.84	116.88	114.13
3	C	507	ROL	C16-O2-C8	-2.80	113.40	117.51
3	A	503	ROL	C4-N1-C1	2.68	116.73	114.13
3	D	509	ROL	C4-N1-C1	2.56	116.61	114.13
3	B	505	ROL	C2-C1-N1	-2.54	106.26	108.68
3	C	507	ROL	C2-C1-N1	-2.54	106.26	108.68
3	A	503	ROL	C2-C3-C5	-2.41	110.33	115.29
3	A	503	ROL	C2-C1-N1	-2.23	106.56	108.68
3	A	503	ROL	C7-O3-C11	-2.13	113.31	120.14
3	D	509	ROL	C4-C3-C2	-2.02	101.20	104.03

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	503	ROL	C15-C11-O3-C7
3	A	503	ROL	C12-C11-O3-C7
3	B	505	ROL	C15-C11-O3-C7

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Mol	Chain	Res	Type	Atoms
3	B	505	ROL	C12-C11-O3-C7
3	C	507	ROL	C15-C11-O3-C7
3	C	507	ROL	C12-C11-O3-C7
3	D	509	ROL	C15-C11-O3-C7
3	D	509	ROL	C12-C11-O3-C7
3	A	503	ROL	C7-C8-O2-C16
3	A	503	ROL	C9-C8-O2-C16

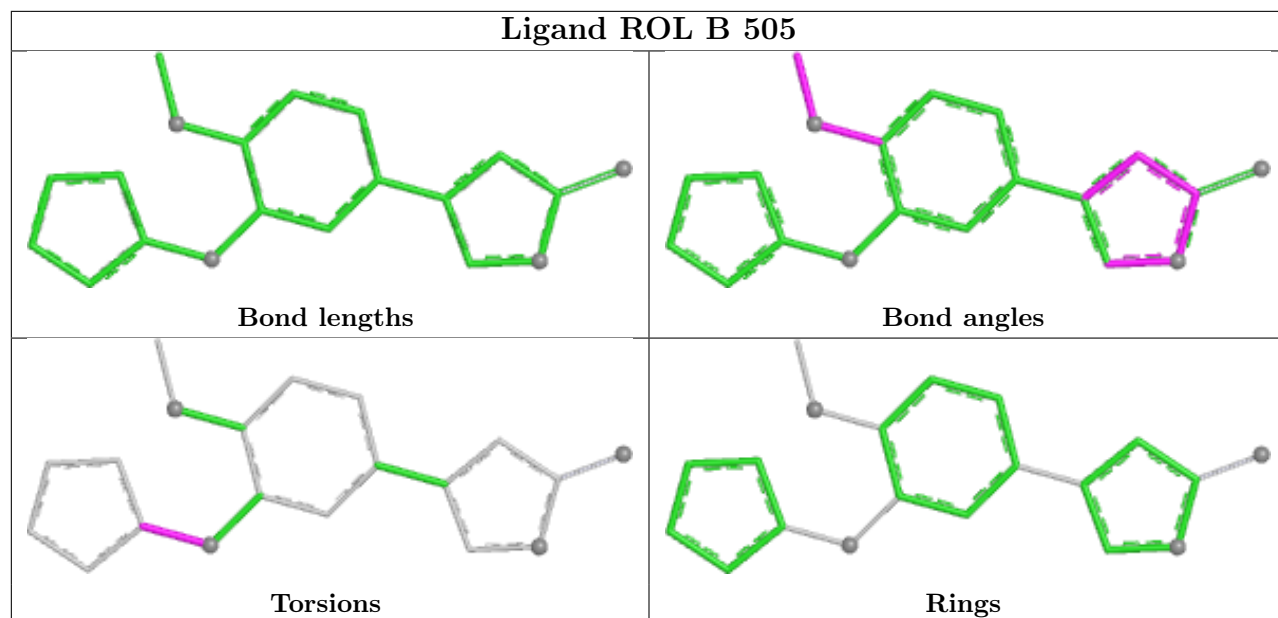
There are no ring outliers.

4 monomers are involved in 7 short contacts:

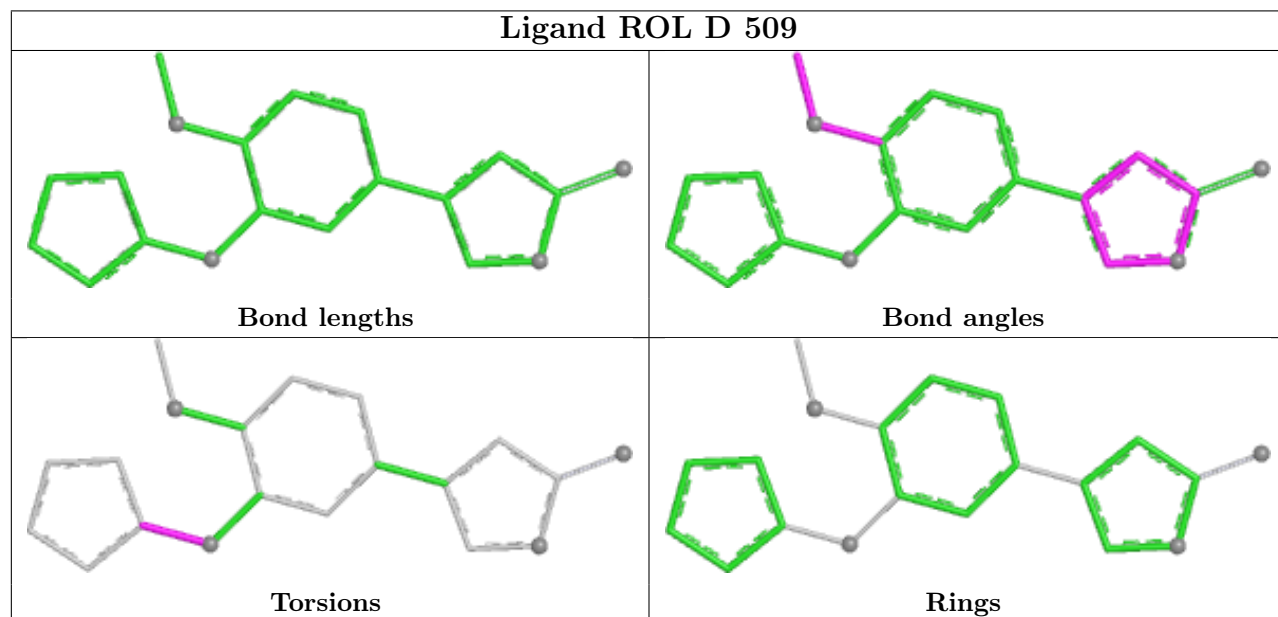
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	505	ROL	3	0
3	D	509	ROL	1	0
3	A	503	ROL	2	0
3	C	507	ROL	1	0

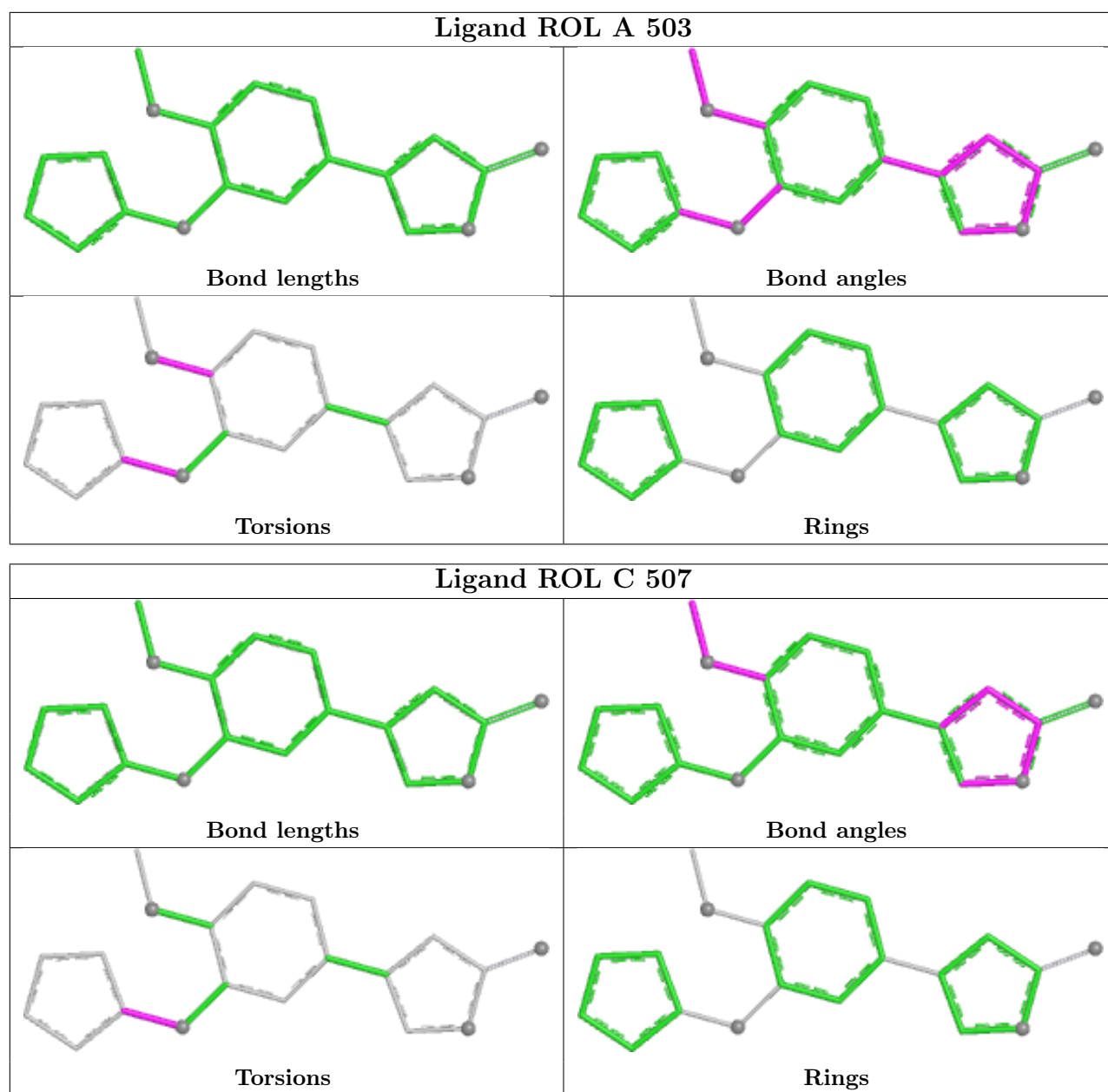
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 ROL B 505



Ligand ROL D 509





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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	334/360 (92%)	0.34	29 (8%) 16 17	16, 30, 71, 97	0
1	B	327/360 (90%)	1.11	73 (22%) 2 2	18, 42, 76, 100	0
1	C	329/360 (91%)	0.76	41 (12%) 8 9	17, 39, 79, 92	0
1	D	334/360 (92%)	0.28	19 (5%) 29 31	16, 30, 53, 76	0
All	All	1324/1440 (91%)	0.62	162 (12%) 8 9	16, 35, 73, 100	0

All (162) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	79	ILE	10.0
1	A	82	PHE	7.4
1	C	84	VAL	7.1
1	D	79	ILE	6.1
1	A	81	ARG	5.9
1	B	86	THR	5.7
1	C	86	THR	5.3
1	A	299	LEU	5.3
1	A	80	PRO	5.1
1	C	302	ASN	4.6
1	A	295	SER	4.5
1	D	412	GLN	4.5
1	B	244	GLU	4.5
1	B	299	LEU	4.4
1	C	293	THR	4.4
1	B	392	ALA	4.4
1	B	91	VAL	4.2
1	A	83	GLY	4.2
1	B	298	LEU	4.1
1	B	296	GLY	4.1
1	C	296	GLY	4.1

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Mol	Chain	Res	Type	RSRZ
1	C	301	ASP	4.1
1	D	86	THR	4.0
1	A	294	SER	4.0
1	C	297	VAL	4.0
1	B	292	VAL	3.9
1	C	298	LEU	3.9
1	D	84	VAL	3.8
1	A	293	THR	3.8
1	C	292	VAL	3.8
1	A	411	PRO	3.8
1	A	412	GLN	3.7
1	C	412	GLN	3.7
1	B	385	ALA	3.6
1	B	295	SER	3.6
1	D	296	GLY	3.6
1	B	92	LEU	3.6
1	B	297	VAL	3.6
1	C	287	VAL	3.6
1	A	298	LEU	3.5
1	B	327	GLN	3.5
1	B	89	GLU	3.4
1	B	88	GLN	3.4
1	B	381	TRP	3.4
1	C	299	LEU	3.4
1	A	291	LYS	3.4
1	B	294	SER	3.4
1	B	291	LYS	3.4
1	C	89	GLU	3.4
1	C	294	SER	3.4
1	A	409	THR	3.3
1	A	84	VAL	3.3
1	C	85	LYS	3.3
1	B	300	LEU	3.3
1	C	295	SER	3.3
1	B	395	ILE	3.3
1	C	411	PRO	3.3
1	B	285	THR	3.3
1	D	82	PHE	3.3
1	B	283	LEU	3.2
1	A	296	GLY	3.2
1	A	301	ASP	3.2
1	C	91	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	289	THR	3.2
1	D	85	LYS	3.1
1	B	180	ALA	3.1
1	B	93	ALA	3.1
1	C	300	LEU	3.1
1	A	286	MET	3.1
1	D	83	GLY	3.1
1	C	303	TYR	3.0
1	A	404	GLU	3.0
1	B	288	GLU	3.0
1	B	99	VAL	3.0
1	B	287	VAL	3.0
1	B	181	LEU	2.9
1	B	390	PRO	2.9
1	B	405	TRP	2.9
1	B	394	ASP	2.9
1	B	284	LYS	2.9
1	B	96	LEU	2.8
1	B	389	HIS	2.8
1	C	279	LEU	2.8
1	C	254	LYS	2.8
1	A	289	THR	2.8
1	D	80	PRO	2.8
1	A	342	ARG	2.8
1	B	102	TRP	2.8
1	C	88	GLN	2.8
1	C	367	LYS	2.8
1	C	392	ALA	2.8
1	D	409	THR	2.7
1	C	255	LYS	2.7
1	A	300	LEU	2.7
1	B	140	ASP	2.7
1	C	291	LYS	2.7
1	B	277	MET	2.7
1	B	387	LEU	2.7
1	B	270	ALA	2.6
1	C	406	TYR	2.6
1	C	87	GLU	2.6
1	A	86	THR	2.6
1	C	409	THR	2.6
1	C	277	MET	2.6
1	B	383	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	87	GLU	2.6
1	D	294	SER	2.6
1	B	132	LEU	2.5
1	B	396	LEU	2.5
1	D	404	GLU	2.5
1	C	281	ALA	2.5
1	B	384	TRP	2.5
1	A	292	VAL	2.5
1	C	395	ILE	2.5
1	B	399	LEU	2.5
1	B	411	PRO	2.5
1	B	293	THR	2.5
1	A	85	LYS	2.5
1	B	183	ALA	2.5
1	B	103	GLY	2.5
1	B	412	GLN	2.5
1	B	386	ASP	2.4
1	C	90	ASP	2.4
1	B	108	ARG	2.4
1	A	87	GLU	2.4
1	B	94	LYS	2.3
1	B	239	LYS	2.3
1	B	281	ALA	2.3
1	D	342	ARG	2.3
1	B	393	GLN	2.3
1	D	91	VAL	2.3
1	B	104	LEU	2.3
1	B	90	ASP	2.3
1	B	375	TYR	2.3
1	B	112	LEU	2.2
1	B	98	ASP	2.2
1	B	306	ARG	2.2
1	D	295	SER	2.2
1	B	380	LEU	2.2
1	B	106	VAL	2.2
1	B	179	PRO	2.2
1	B	290	LYS	2.2
1	C	290	LYS	2.2
1	B	363	ALA	2.1
1	C	178	THR	2.1
1	C	362	ASN	2.1
1	D	221	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	327	GLN	2.1
1	B	127	GLN	2.1
1	D	327	GLN	2.1
1	B	139	VAL	2.1
1	B	286	MET	2.1
1	C	357	MET	2.1
1	B	128	GLU	2.1
1	A	386	ASP	2.1
1	B	301	ASP	2.1
1	A	303	TYR	2.1
1	C	280	LEU	2.0
1	B	388	VAL	2.0
1	D	81	ARG	2.0
1	B	289	THR	2.0
1	C	396	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	ROL	A	503	20/20	0.89	0.12	34,36,37,40	0
3	ROL	C	507	20/20	0.89	0.12	38,39,46,48	0
3	ROL	D	509	20/20	0.91	0.11	32,37,41,44	0
3	ROL	B	505	20/20	0.92	0.11	34,37,41,43	0
2	ZN	A	502	1/1	0.93	0.20	61,61,61,61	0
2	ZN	B	504	1/1	0.93	0.17	64,64,64,64	0
2	ZN	D	508	1/1	0.94	0.14	61,61,61,61	0
2	ZN	C	506	1/1	0.96	0.17	59,59,59,59	0

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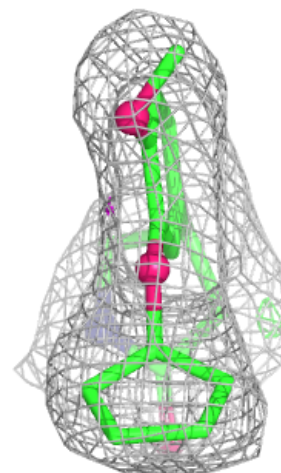
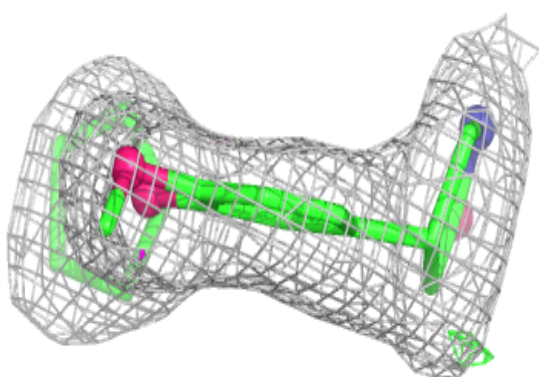
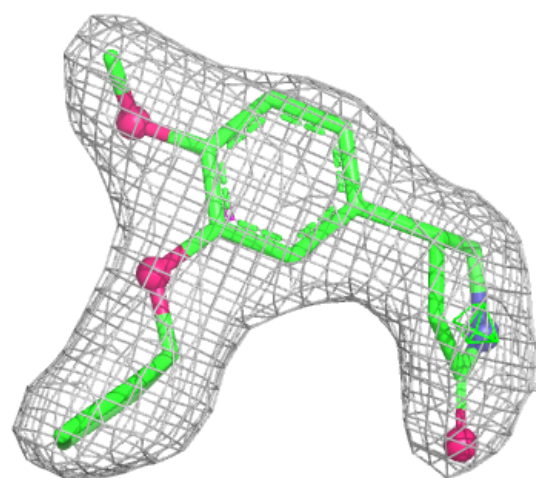
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	B	503	1/1	0.97	0.05	34,34,34,34	0
2	ZN	D	507	1/1	0.98	0.03	31,31,31,31	0
2	ZN	A	501	1/1	0.98	0.03	30,30,30,30	0
2	ZN	C	505	1/1	0.99	0.04	32,32,32,32	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

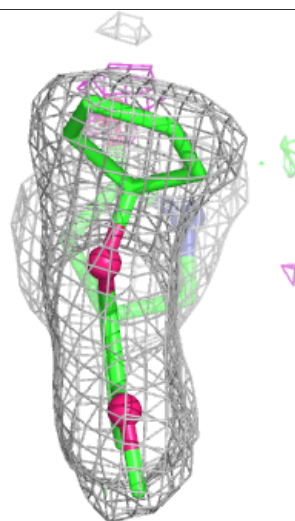
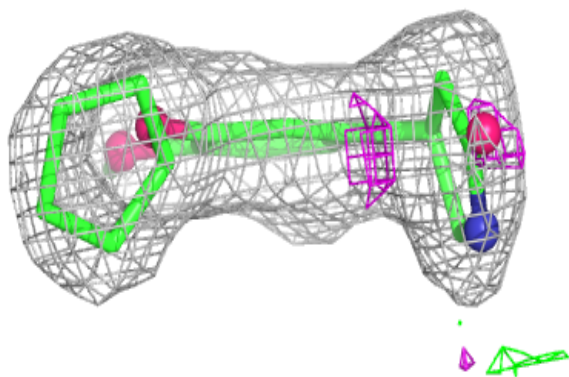
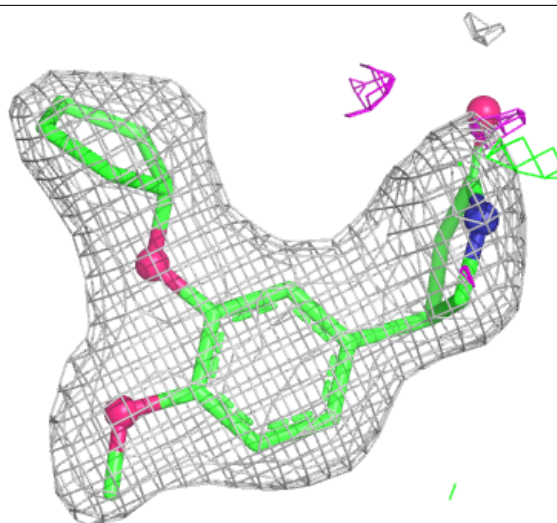
Electron density around ROL A 503:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



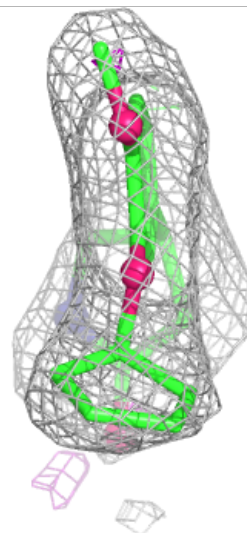
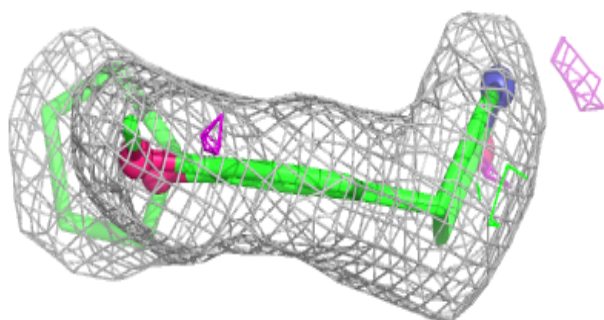
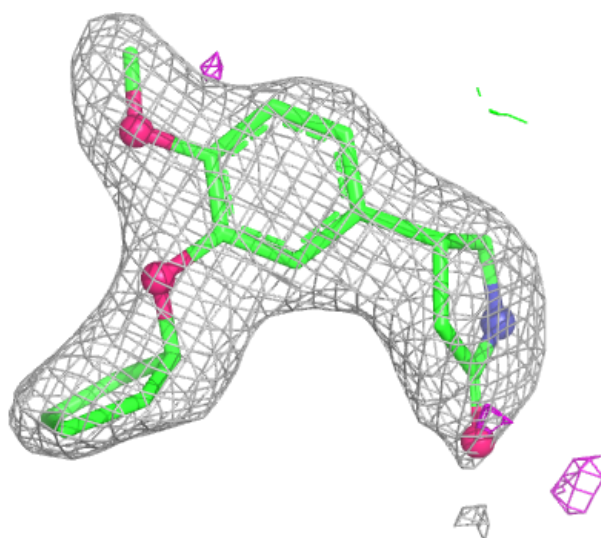
Electron density around ROL C 507:

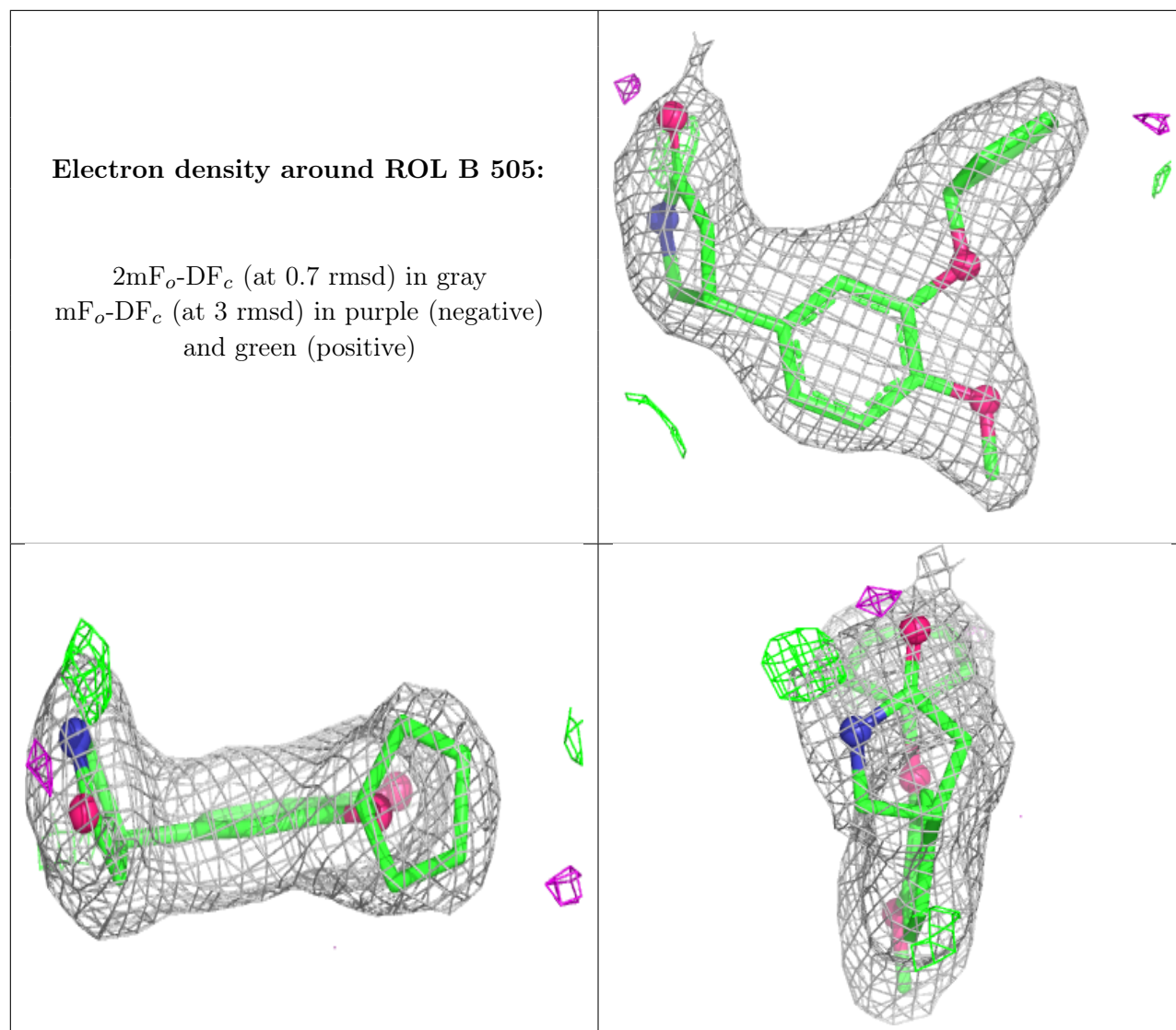
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ROL D 509:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers ⓘ

There are no such residues in this entry.