



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:32 PM UTC

PDB ID : 2PO2 / pdb_00002po2
Title : Crystal structure of the P. abyssi exosome RNase PH ring complexed with CDP
Authors : Navarro, M.V.A.S.; Guimaraes, B.G.
Deposited on : 2007-04-25
Resolution : 2.41 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

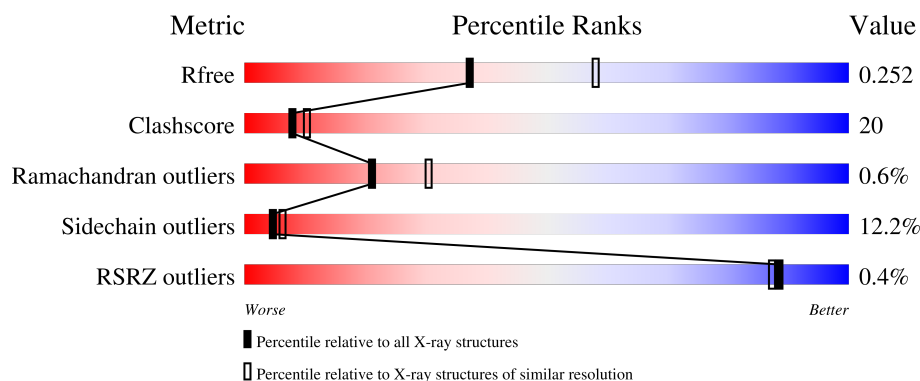
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 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	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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	249	 55% 34% . . 5%
2	B	277	 58% 31% 6% . .

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	MPD	A	251	X	-	X	-
4	MPD	A	283	X	-	-	-
4	MPD	B	282	X	-	-	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4144 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 Probable exosome complex exonuclease 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	236	Total	C	N	O	S	0	1	0
			1848	1176	326	339	7			

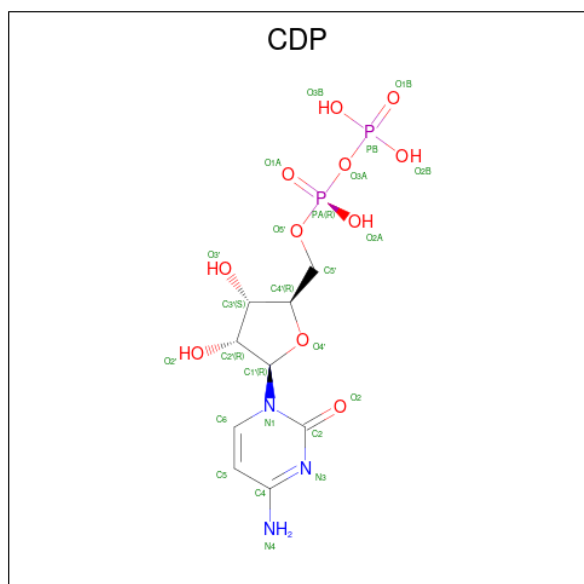
- Molecule 2 is a protein called Probable exosome complex exonuclease 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	267	Total	C	N	O	S	0	2	0
			2055	1306	342	400	7			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	GLY	-	expression tag	UNP Q9V118
B	-1	SER	-	expression tag	UNP Q9V118
B	0	HIS	-	expression tag	UNP Q9V118

- Molecule 3 is CYTIDINE-5'-DIPHOSPHATE (CCD ID: CDP) (formula: $C_9H_{15}N_3O_{11}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			25	9	3	11	2		

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			8	6	2		
4	A	1	Total	C	O	0	0
			8	6	2		
4	B	1	Total	C	O	0	0
			8	6	2		
4	B	1	Total	C	O	0	0
			8	6	2		
4	B	1	Total	C	O	0	0
			8	6	2		

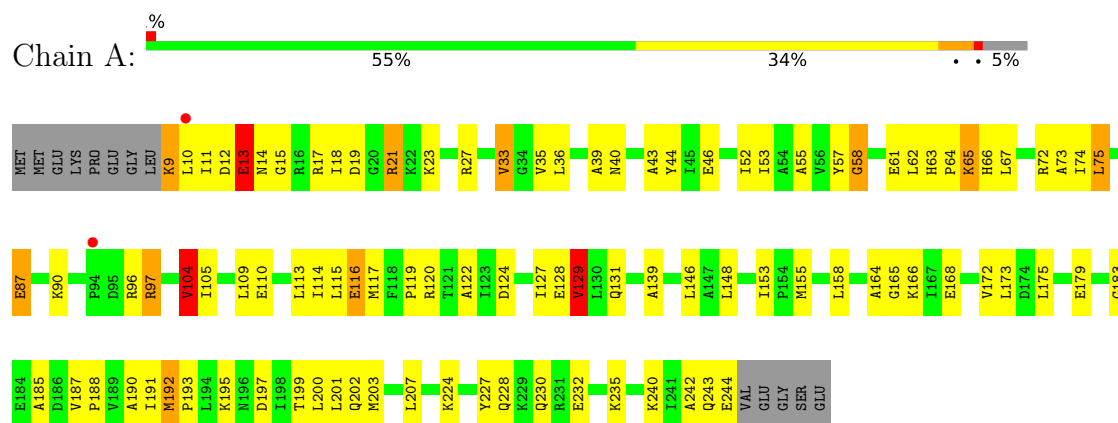
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	76	Total	O	0	0
			76	76		
5	B	100	Total	O	0	0
			100	100		

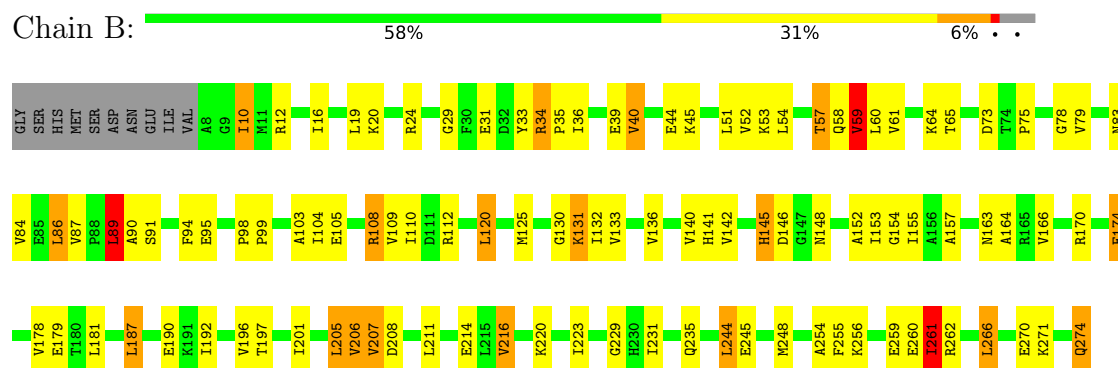
3 Residue-property plots [i](#)

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: Probable exosome complex exonuclease 1



- Molecule 2: Probable exosome complex exonuclease 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	93.50Å 93.50Å 125.71Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.80 – 2.41 19.80 – 2.41	Depositor EDS
% Data completeness (in resolution range)	99.3 (19.80-2.41) 99.3 (19.80-2.41)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.41Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.182 , 0.259 0.180 , 0.252	Depositor DCC
R_{free} test set	1273 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	37.0	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.053 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4144	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.28% 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: CDP, MPD

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	1.15	2/1878 (0.1%)	1.22	12/2528 (0.5%)
2	B	1.21	7/2090 (0.3%)	1.27	6/2832 (0.2%)
All	All	1.18	9/3968 (0.2%)	1.25	18/5360 (0.3%)

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

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	174	GLU	CD-OE1	9.64	1.43	1.25
2	B	174	GLU	CG-CD	9.41	1.75	1.52
2	B	79	VAL	CA-CB	7.02	1.63	1.54
1	A	240	LYS	CG-CD	6.88	1.73	1.52
2	B	174	GLU	CD-OE2	5.68	1.36	1.25
2	B	261	ILE	CA-CB	5.42	1.60	1.54
1	A	129	VAL	CA-C	-5.12	1.47	1.52
2	B	57	THR	CA-CB	5.10	1.61	1.53
2	B	59	VAL	CA-CB	5.02	1.61	1.54

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	95	GLU	CA-C-N	-6.48	113.08	119.76
2	B	95	GLU	C-N-CA	-6.48	113.08	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	129	VAL	CB-CA-C	6.37	119.34	111.25
1	A	192	MET	CA-C-N	6.17	125.90	119.05
1	A	192	MET	C-N-CA	6.17	125.90	119.05
1	A	173	LEU	CB-CA-C	5.59	118.87	109.48
2	B	89	LEU	CA-CB-CG	5.57	135.79	116.30
2	B	29	GLY	N-CA-C	-5.55	103.91	112.85
1	A	129	VAL	N-CA-CB	5.43	117.99	111.31
1	A	97	ARG	NE-CZ-NH2	-5.38	114.36	119.20
1	A	58	GLY	CA-C-N	5.38	139.90	127.00
1	A	58	GLY	C-N-CA	5.38	139.90	127.00
1	A	104	VAL	CB-CA-C	5.34	119.18	111.65
2	B	174	GLU	CB-CG-CD	-5.11	103.91	112.60
1	A	164	ALA	N-CA-C	-5.10	102.01	109.81
2	B	231	ILE	N-CA-C	-5.09	101.04	108.17
1	A	43	ALA	N-CA-C	-5.05	101.07	108.99
1	A	97	ARG	CD-NE-CZ	-5.02	117.37	124.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	9	LYS	Peptide

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	1848	0	1936	77	0
2	B	2055	0	2127	82	0
3	A	25	0	12	3	0
4	A	16	0	28	8	0
4	B	24	0	42	3	0
5	A	76	0	0	5	0
5	B	100	0	0	4	0
All	All	4144	0	4145	160	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 20.

All (160) 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:174:GLU:CD	2:B:174:GLU:CG	1.75	1.56
2:B:12:ARG:HE	2:B:16:ILE:HD11	1.16	1.06
4:A:251:MPD:HM2	4:A:251:MPD:H53	1.40	1.00
2:B:12:ARG:NE	2:B:16:ILE:HD11	1.78	0.97
1:A:19:ASP:OD2	1:A:21:ARG:HD3	1.70	0.89
2:B:270:GLU:HG2	5:B:348:HOH:O	1.73	0.88
1:A:183:GLY:O	4:A:251:MPD:HM3	1.79	0.82
1:A:97:ARG:HH22	3:A:250:CDP:PA	2.02	0.81
1:A:104:VAL:HG13	1:A:200:LEU:HD11	1.63	0.81
2:B:40:VAL:HG13	2:B:163:ASN:CG	2.05	0.81
1:A:193:PRO:HG2	1:A:227:TYR:CD1	2.17	0.80
1:A:12:ASP:O	1:A:13:GLU:HB2	1.82	0.78
4:A:251:MPD:H53	4:A:251:MPD:CM	2.16	0.74
1:A:97:ARG:NH2	3:A:250:CDP:O2A	2.22	0.73
1:A:104:VAL:HG13	1:A:200:LEU:CD1	2.18	0.73
1:A:53:ILE:HG23	2:B:89:LEU:HD21	1.72	0.71
1:A:227:TYR:HA	1:A:230:GLN:HE21	1.56	0.71
1:A:62:LEU:HD11	1:A:74:ILE:CD1	2.21	0.70
1:A:13:GLU:HG3	1:A:14:ASN:ND2	2.05	0.70
1:A:13:GLU:HG3	1:A:14:ASN:HD22	1.55	0.70
2:B:61:VAL:HG22	2:B:140:VAL:HG22	1.73	0.69
2:B:174:GLU:CD	2:B:174:GLU:CB	2.65	0.69
1:A:72:ARG:NH1	1:A:115:LEU:HD13	2.08	0.68
1:A:197:ASP:OD1	5:A:353:HOH:O	2.12	0.68
1:A:192:MET:HE2	1:A:199:THR:CG2	2.23	0.68
1:A:193:PRO:HG2	1:A:227:TYR:CG	2.30	0.66
2:B:34:ARG:NH1	2:B:146:ASP:O	2.29	0.66
1:A:192:MET:HE2	1:A:199:THR:HG21	1.78	0.65
4:A:283:MPD:HM2	2:B:58:GLN:OE1	1.97	0.65
4:B:280:MPD:H52	4:B:280:MPD:H11	1.78	0.64
1:A:97:ARG:NH2	3:A:250:CDP:PA	2.70	0.64
1:A:19:ASP:CG	1:A:21:ARG:HD3	2.25	0.62
2:B:105:GLU:O	2:B:109:VAL:HG23	1.99	0.62
1:A:228:GLN:HB3	5:A:338:HOH:O	2.01	0.61
2:B:16:ILE:HG23	2:B:20:LYS:HE3	1.83	0.61
2:B:40:VAL:HG13	2:B:163:ASN:ND2	2.16	0.61
2:B:205:LEU:H	4:B:282:MPD:H31	1.66	0.60
2:B:164:ALA:CB	2:B:187:LEU:HD22	2.32	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:LYS:HB2	1:A:172:VAL:HG12	1.83	0.60
1:A:72:ARG:HH12	1:A:115:LEU:HD13	1.64	0.60
2:B:214:GLU:OE2	2:B:220:LYS:NZ	2.30	0.59
2:B:271:LYS:O	2:B:274:GLN:HB2	2.02	0.59
2:B:164:ALA:HB3	2:B:187:LEU:HD22	1.83	0.59
2:B:229:GLY:HA2	2:B:255:PHE:CE2	2.39	0.58
2:B:87:VAL:HG22	2:B:142:VAL:O	2.03	0.58
2:B:40:VAL:CG1	2:B:163:ASN:ND2	2.68	0.57
2:B:31:GLU:HA	2:B:207:VAL:HG23	1.87	0.56
2:B:192:ILE:HB	2:B:262:ARG:HD2	1.87	0.56
2:B:78:GLY:HA3	2:B:125:MET:HE2	1.87	0.56
2:B:61:VAL:HG11	2:B:157:ALA:HA	1.86	0.56
2:B:86:LEU:HD22	2:B:103:ALA:HB2	1.86	0.56
2:B:104:ILE:O	2:B:108:ARG:HG3	2.06	0.56
1:A:62:LEU:HD11	1:A:74:ILE:HD13	1.88	0.55
1:A:64:PRO:O	1:A:66:HIS:N	2.39	0.55
1:A:166:LYS:NZ	5:A:339:HOH:O	2.38	0.55
2:B:65:THR:HG21	2:B:166:VAL:HG11	1.88	0.55
2:B:83:ASN:HB3	5:B:288:HOH:O	2.06	0.55
1:A:131:GLN:HE22	2:B:44:GLU:N	2.04	0.55
2:B:59:VAL:HG13	2:B:152:ALA:HB1	1.87	0.55
2:B:166:VAL:HG11	2:B:187:LEU:HD13	1.89	0.55
1:A:62:LEU:HG	1:A:122:ALA:HB2	1.89	0.55
1:A:165:GLY:HA3	1:A:175:LEU:HD11	1.89	0.54
4:A:251:MPD:HM2	4:A:251:MPD:C5	2.19	0.54
1:A:62:LEU:HD22	1:A:67:LEU:HB2	1.89	0.54
2:B:40:VAL:HG13	2:B:163:ASN:CB	2.37	0.53
2:B:256:LYS:O	2:B:260:GLU:HG3	2.08	0.53
1:A:114:ILE:O	1:A:117:MET:HB2	2.08	0.53
1:A:53:ILE:HG22	1:A:128:GLU:HB2	1.89	0.53
1:A:72:ARG:NH1	1:A:73:ALA:O	2.41	0.52
1:A:187:VAL:HG22	1:A:203:MET:HG2	1.91	0.52
1:A:17:ARG:HD2	1:A:179:GLU:OE1	2.10	0.52
1:A:185:ALA:CB	1:A:207:LEU:HD21	2.40	0.51
2:B:24:ARG:CZ	2:B:206:VAL:HG22	2.39	0.51
2:B:120:LEU:HD21	2:B:187:LEU:HG	1.92	0.51
4:A:251:MPD:CM	4:A:251:MPD:C5	2.73	0.51
1:A:62:LEU:HD11	1:A:74:ILE:HD11	1.92	0.51
2:B:40:VAL:HG13	2:B:163:ASN:HB2	1.92	0.51
2:B:90:ALA:HB3	2:B:94:PHE:HD1	1.76	0.51
1:A:64:PRO:C	1:A:66:HIS:H	2.20	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:78:GLY:HA2	2:B:133:VAL:HB	1.94	0.50
1:A:190:ALA:HB3	1:A:200:LEU:HB3	1.94	0.50
1:A:52:ILE:HG12	1:A:129:VAL:HB	1.94	0.49
1:A:185:ALA:HB2	1:A:207:LEU:HD21	1.93	0.49
2:B:34:ARG:HB3	2:B:35:PRO:CD	2.43	0.49
1:A:158:LEU:HB2	1:A:227:TYR:HE1	1.77	0.49
1:A:224:LYS:O	1:A:228:GLN:HG2	2.12	0.49
2:B:34:ARG:HB3	2:B:35:PRO:HD2	1.95	0.49
2:B:65:THR:HG22	2:B:136:VAL:HG22	1.95	0.49
2:B:166:VAL:HG11	2:B:187:LEU:CD1	2.43	0.49
1:A:12:ASP:O	1:A:13:GLU:CB	2.58	0.48
2:B:170:ARG:NH1	2:B:179:GLU:OE1	2.46	0.48
2:B:197:THR:HG21	5:B:316:HOH:O	2.12	0.48
2:B:190:GLU:C	4:B:280:MPD:HM3	2.37	0.48
1:A:75:LEU:HD11	1:A:109:LEU:HB2	1.96	0.48
2:B:155:ILE:HG12	2:B:261:ILE:HG21	1.95	0.48
1:A:188:PRO:HD2	1:A:202:GLN:O	2.15	0.47
4:A:283:MPD:HM1	2:B:51:LEU:HD21	1.97	0.47
2:B:52:VAL:HG21	2:B:155:ILE:HG22	1.96	0.47
2:B:90:ALA:HB3	2:B:94:PHE:CD1	2.50	0.47
2:B:108:ARG:O	2:B:112:ARG:HG3	2.14	0.47
1:A:19:ASP:OD2	1:A:21:ARG:CD	2.53	0.47
1:A:62:LEU:HD23	1:A:63:HIS:N	2.30	0.47
1:A:53:ILE:CG2	2:B:89:LEU:HD21	2.42	0.47
2:B:12:ARG:HD2	2:B:216:VAL:HB	1.97	0.47
2:B:244:LEU:O	2:B:248:MET:HG3	2.15	0.46
1:A:131:GLN:NE2	2:B:45:LYS:H	2.13	0.46
2:B:84:VAL:HA	2:B:140:VAL:HB	1.97	0.46
1:A:14:ASN:HA	1:A:15:GLY:HA2	1.82	0.46
1:A:113:LEU:HD23	1:A:155:MET:HG2	1.97	0.46
2:B:166:VAL:CG1	2:B:187:LEU:HD13	2.46	0.46
1:A:66:HIS:HB3	5:A:315:HOH:O	2.15	0.46
1:A:116:GLU:H	1:A:116:GLU:CD	2.24	0.46
2:B:65:THR:HG21	2:B:166:VAL:CG1	2.46	0.46
1:A:158:LEU:HB2	1:A:227:TYR:CE1	2.51	0.45
1:A:105:ILE:HD13	1:A:127:ILE:HD13	1.97	0.45
1:A:9:LYS:HA	1:A:10:LEU:HG	1.97	0.45
2:B:105:GLU:HG3	2:B:235:GLN:HE21	1.81	0.45
2:B:59:VAL:HG21	2:B:153:ILE:HG12	1.98	0.45
4:A:251:MPD:HM1	4:A:251:MPD:H4	1.55	0.44
2:B:10:ILE:H	2:B:10:ILE:HG12	1.64	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:91:SER:HB2	2:B:145:HIS:HB3	1.99	0.44
1:A:87:GLU:HB2	5:B:290:HOH:O	2.18	0.44
1:A:131:GLN:HE22	2:B:44:GLU:H	1.63	0.44
2:B:105:GLU:HG3	2:B:235:GLN:NE2	2.33	0.44
2:B:130:GLY:C	2:B:131:LYS:HG2	2.42	0.44
2:B:12:ARG:CZ	2:B:16:ILE:HD11	2.45	0.44
2:B:53:LYS:HA	2:B:57:THR:O	2.16	0.44
1:A:35:VAL:HG23	1:A:36:LEU:HG	2.00	0.44
1:A:188:PRO:HG2	1:A:202:GLN:HB2	1.98	0.43
1:A:52:ILE:HD12	1:A:139:ALA:HB3	2.00	0.43
2:B:33:TYR:OH	2:B:196:VAL:HG13	2.18	0.43
2:B:229:GLY:HA2	2:B:255:PHE:HE2	1.80	0.43
2:B:110:ILE:HD12	2:B:154:GLY:CA	2.48	0.43
2:B:40:VAL:CG1	2:B:163:ASN:HB2	2.48	0.43
1:A:44:TYR:CE2	1:A:46:GLU:HG3	2.53	0.43
2:B:75:PRO:CB	2:B:130:GLY:HA2	2.49	0.43
1:A:96:ARG:H	1:A:96:ARG:HG2	1.50	0.43
2:B:54:LEU:HD13	2:B:152:ALA:HA	1.99	0.43
2:B:130:GLY:O	2:B:131:LYS:HE3	2.19	0.43
1:A:155:MET:HE1	1:A:158:LEU:CD2	2.49	0.42
2:B:148:ASN:ND2	2:B:208:ASP:HA	2.33	0.42
2:B:84:VAL:HG12	2:B:86:LEU:HD13	2.02	0.42
2:B:223:ILE:HB	2:B:254:ALA:CB	2.49	0.42
1:A:40:ASN:OD1	1:A:58:GLY:HA2	2.20	0.42
1:A:13:GLU:HA	1:A:14:ASN:HA	1.56	0.42
1:A:39:ALA:HB2	1:A:55:ALA:HB1	2.02	0.42
2:B:19:LEU:HD23	2:B:19:LEU:HA	1.87	0.42
2:B:132:ILE:HD13	2:B:178:VAL:HG21	2.02	0.42
2:B:98:PRO:O	2:B:99:PRO:C	2.61	0.41
1:A:113:LEU:HD21	1:A:148:LEU:HD13	2.00	0.41
1:A:27:ARG:NH2	5:A:288:HOH:O	2.45	0.41
1:A:155:MET:CE	1:A:158:LEU:HD23	2.51	0.41
1:A:155:MET:HE1	1:A:158:LEU:HD23	2.02	0.41
1:A:232:GLU:OE1	1:A:235:LYS:HD3	2.20	0.41
2:B:19:LEU:HD11	2:B:201:ILE:HB	2.01	0.41
1:A:33:VAL:CG2	1:A:146:LEU:HB3	2.51	0.41
2:B:192:ILE:HD13	2:B:266:LEU:HD13	2.03	0.41
1:A:57:TYR:O	1:A:124:ASP:N	2.44	0.40
1:A:97:ARG:HH21	1:A:97:ARG:HD2	1.45	0.40
1:A:116:GLU:O	1:A:119:PRO:HD3	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	235/249 (94%)	220 (94%)	12 (5%)	3 (1%)	9	13
2	B	267/277 (96%)	259 (97%)	8 (3%)	0	100	100
All	All	502/526 (95%)	479 (95%)	20 (4%)	3 (1%)	21	30

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	13	GLU
1	A	65	LYS
1	A	242	ALA

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	192/202 (95%)	170 (88%)	22 (12%)	5	7
2	B	228/235 (97%)	199 (87%)	29 (13%)	4	5
All	All	420/437 (96%)	369 (88%)	51 (12%)	5	6

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ILE
1	A	13	GLU
1	A	18	ILE

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Mol	Chain	Res	Type
1	A	21	ARG
1	A	33	VAL
1	A	61	GLU
1	A	65	LYS
1	A	75	LEU
1	A	87	GLU
1	A	90	LYS
1	A	104	VAL
1	A	110	GLU
1	A	116	GLU
1	A	120	ARG
1	A	129	VAL
1	A	153	ILE
1	A	168	GLU
1	A	191	ILE
1	A	195	LYS
1	A	201	LEU
1	A	243	GLN
1	A	244	GLU
2	B	10	ILE
2	B	34	ARG
2	B	36	ILE
2	B	39	GLU
2	B	40	VAL
2	B	59	VAL
2	B	60	LEU
2	B	64	LYS
2	B	73	ASP
2	B	86	LEU
2	B	89	LEU
2	B	108	ARG
2	B	120	LEU
2	B	131	LYS
2	B	141	HIS
2	B	145	HIS
2	B	181	LEU
2	B	187	LEU
2	B	205	LEU
2	B	206	VAL
2	B	207	VAL
2	B	211	LEU
2	B	216	VAL

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Mol	Chain	Res	Type
2	B	244	LEU
2	B	245	GLU
2	B	259	GLU
2	B	261	ILE
2	B	266	LEU
2	B	274	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	131	GLN
1	A	196	ASN
1	A	230	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](#)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	CDP	A	250	-	25,26,26	1.43	2 (8%)	38,40,40	1.63	7 (18%)
4	MPD	A	251	-	7,7,7	0.77	0	9,10,10	1.53	2 (22%)
4	MPD	A	283	-	7,7,7	0.40	0	9,10,10	0.69	0
4	MPD	B	280	-	7,7,7	0.39	0	9,10,10	0.94	0
4	MPD	B	281	-	7,7,7	0.52	0	9,10,10	0.85	0
4	MPD	B	282	-	7,7,7	0.43	0	9,10,10	0.49	0

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	CDP	A	250	-	-	2/16/32/32	0/2/2/2
4	MPD	A	251	-	1/1/2/2	5/5/5/5	-
4	MPD	A	283	-	1/1/2/2	3/5/5/5	-
4	MPD	B	280	-	-	3/5/5/5	-
4	MPD	B	281	-	-	0/5/5/5	-
4	MPD	B	282	-	1/1/2/2	5/5/5/5	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	250	CDP	PA-O3A	5.41	1.65	1.59
3	A	250	CDP	C6-C5	2.32	1.40	1.35

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	250	CDP	O2-C2-N3	-4.37	115.44	122.33
3	A	250	CDP	O4'-C1'-N1	3.87	117.14	108.36
4	A	251	MPD	CM-C2-C1	-3.54	102.69	110.63
3	A	250	CDP	C2'-C1'-N1	-3.42	103.73	113.25
3	A	250	CDP	O2B-PB-O3B	3.26	120.04	107.80
3	A	250	CDP	O2-C2-N1	2.71	124.20	118.90
3	A	250	CDP	N4-C4-N3	-2.13	114.11	117.91
3	A	250	CDP	C5-C4-N4	2.01	124.15	120.63
4	A	251	MPD	O2-C2-C1	2.01	114.25	107.99

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	A	251	MPD	C4
4	A	283	MPD	C4
4	B	282	MPD	C4

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	250	CDP	PA-O3A-PB-O3B
4	A	251	MPD	O2-C2-C3-C4
4	A	251	MPD	CM-C2-C3-C4
4	A	251	MPD	C2-C3-C4-O4
4	A	251	MPD	C2-C3-C4-C5
4	A	283	MPD	C2-C3-C4-C5
4	B	280	MPD	C2-C3-C4-O4
4	B	280	MPD	C2-C3-C4-C5
4	B	282	MPD	C2-C3-C4-O4
4	B	282	MPD	C2-C3-C4-C5
4	B	282	MPD	O2-C2-C3-C4
4	B	282	MPD	C1-C2-C3-C4
4	B	282	MPD	CM-C2-C3-C4
3	A	250	CDP	PA-O3A-PB-O1B
4	A	283	MPD	O2-C2-C3-C4
4	A	251	MPD	C1-C2-C3-C4
4	A	283	MPD	CM-C2-C3-C4
4	B	280	MPD	CM-C2-C3-C4

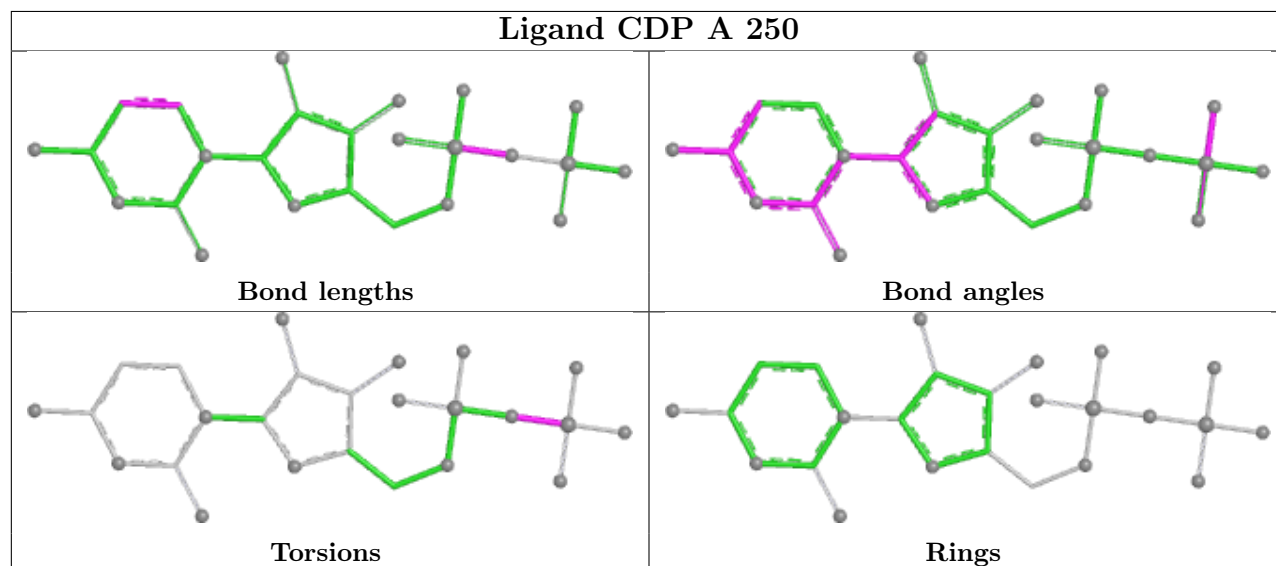
There are no ring outliers.

5 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	250	CDP	3	0
4	A	251	MPD	6	0
4	A	283	MPD	2	0
4	B	280	MPD	2	0
4	B	282	MPD	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.



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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	236/249 (94%)	-0.46	2 (0%) 82 80	15, 38, 76, 104	1 (0%)
2	B	267/277 (96%)	-0.64	0 100 100	18, 34, 65, 94	2 (0%)
All	All	503/526 (95%)	-0.56	2 (0%) 88 87	15, 36, 70, 104	3 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	94	PRO	2.3
1	A	10	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.

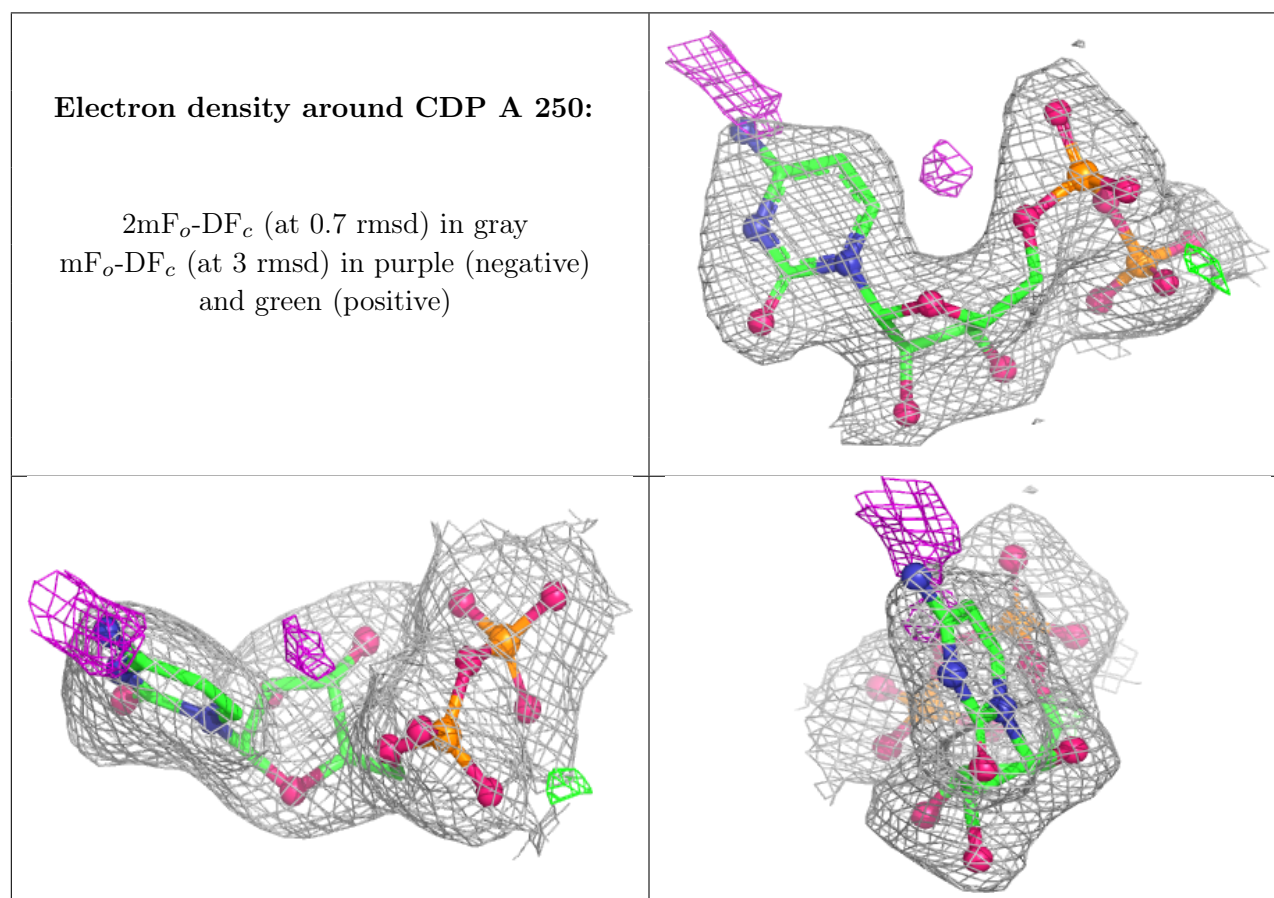
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MPD	A	251	8/8	0.79	0.17	44,49,53,54	0
4	MPD	B	282	8/8	0.86	0.13	53,59,62,64	0
4	MPD	B	281	8/8	0.87	0.13	58,65,68,68	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MPD	A	283	8/8	0.88	0.13	54,61,66,68	0
4	MPD	B	280	8/8	0.92	0.09	47,54,58,58	0
3	CDP	A	250	25/25	0.98	0.05	25,32,36,37	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.