



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 11:50 PM UTC

PDB ID : 6QIB / pdb_00006qib
Title : The crystal structure of Pol2CORE in complex with DNA and an incoming nucleotide, carrying an Fe-S cluster
Authors : Parkash, V.; Johansson, E.
Deposited on : 2019-01-18
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)	:	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

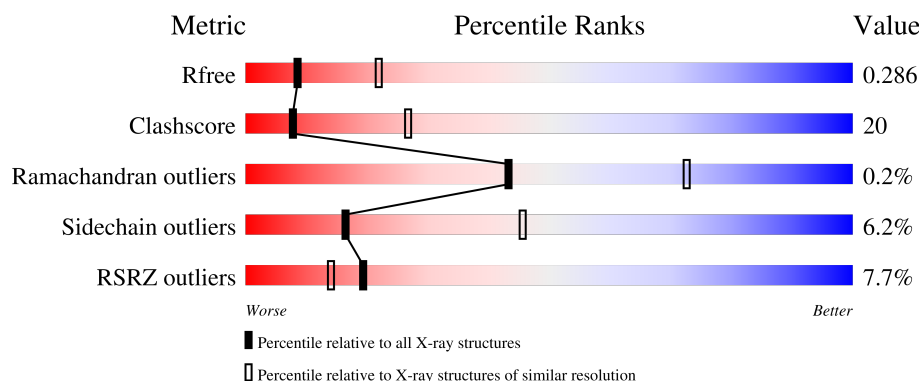
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

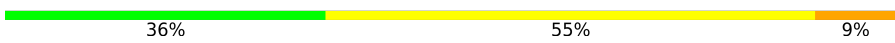
The reported resolution of this entry is 2.80 Å.

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))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (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 $\leq 5\%$. 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	1192	
2	P	11	
3	T	16	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9176 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 DNA polymerase epsilon catalytic subunit A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1123	Total	C	N	O	S	0	0	0
			8608	5494	1445	1626	43			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP P21951
A	-3	GLY	-	expression tag	UNP P21951
A	-2	ASP	-	expression tag	UNP P21951
A	-1	PRO	-	expression tag	UNP P21951
A	0	HIS	-	expression tag	UNP P21951
A	290	ALA	ASP	engineered mutation	UNP P21951
A	292	ALA	GLU	engineered mutation	UNP P21951

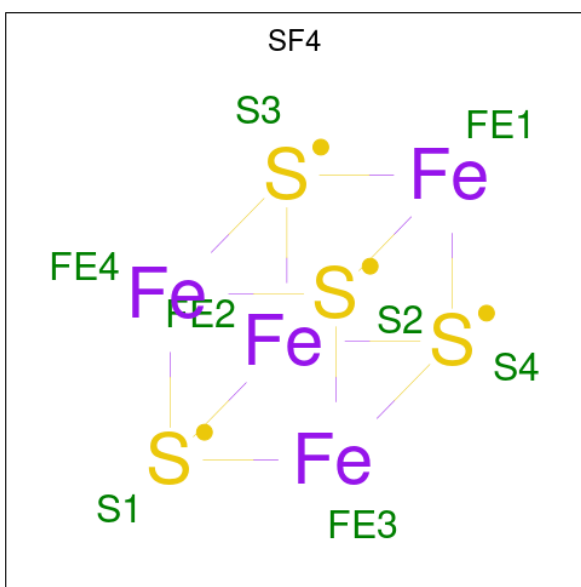
- Molecule 2 is a DNA chain called Primer_11ddC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	11	Total	C	N	O	P	0	0	0
			221	106	38	66	11			

- Molecule 3 is a DNA chain called Template16.

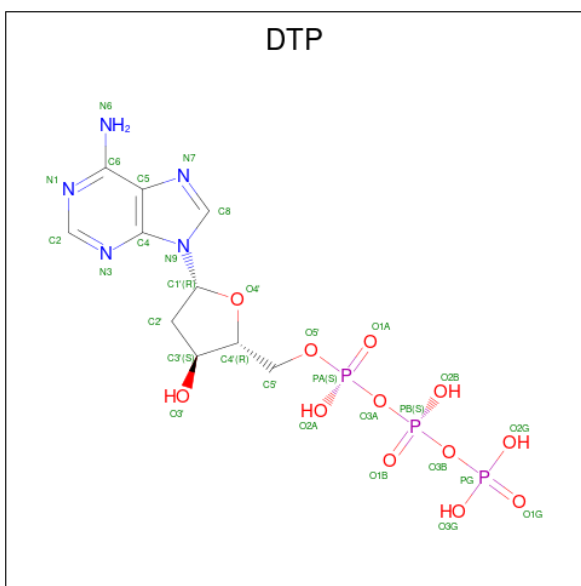
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	T	15	Total	C	N	O	P	0	0	0
			308	147	54	92	15			

- Molecule 4 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	Fe	S		0	0
			8	4	4			

- Molecule 5 is 2'-DEOXYADENOSINE 5'-TRIPHOSPHATE (CCD ID: DTP) (formula: $C_{10}H_{16}N_5O_{12}P_3$)



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	
			30	10	5	12	3	

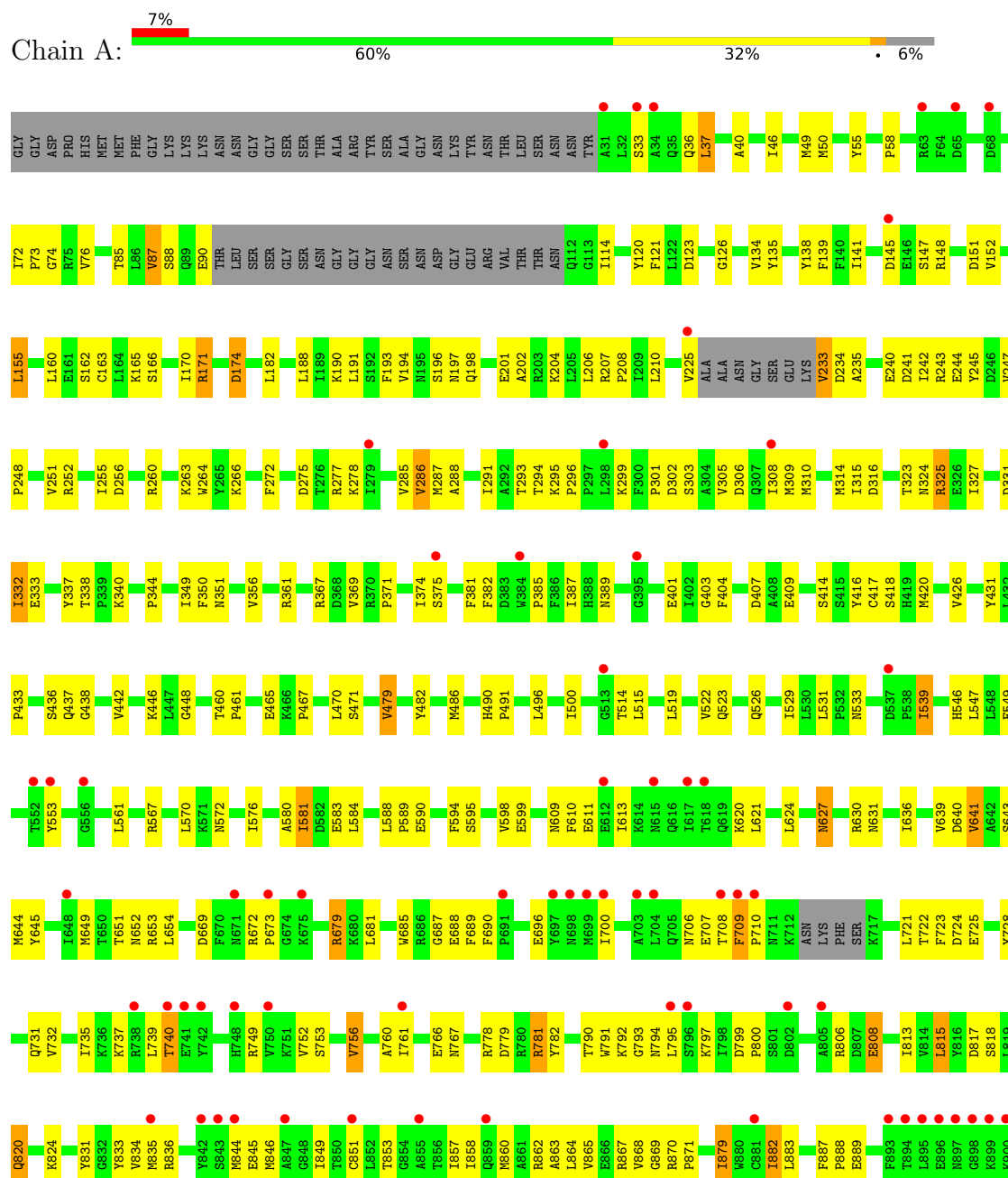
- Molecule 6 is CALCIUM ION (CCD ID: CA) (formula: Ca).

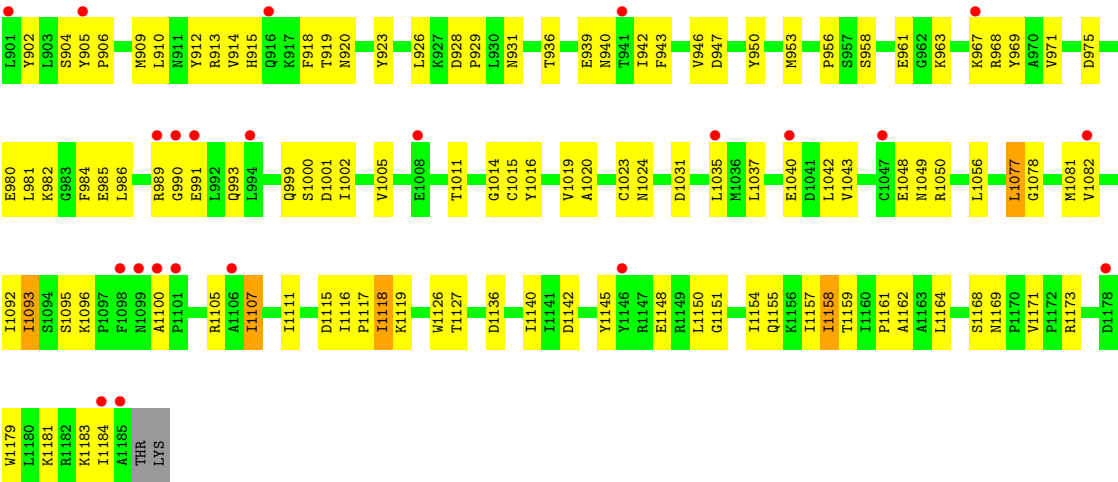
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total 1	Ca 1	0	0

3 Residue-property plots

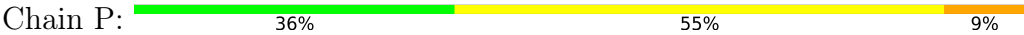
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: DNA polymerase epsilon catalytic subunit A





● Molecule 2: Primer_11ddC



● Molecule 3: Template16



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	158.36Å 70.40Å 154.24Å 90.00° 112.99° 90.00°	Depositor
Resolution (Å)	78.48 – 2.80 78.48 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.3 (78.48-2.80) 98.3 (78.48-2.80)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.82Å)	Xtriage
Refinement program	REFMAC 5.8.0232	Depositor
R, R_{free}	0.223 , 0.286 0.228 , 0.286	Depositor DCC
R_{free} test set	1872 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å ²)	65.2	Xtriage
Anisotropy	0.557	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 75.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9176	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.40% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SF4, CA, DOC, DTP

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	0.45	0/8806	0.99	2/11978 (0.0%)
2	P	0.54	0/226	0.90	2/346 (0.6%)
3	T	0.39	0/344	0.85	1/529 (0.2%)
All	All	0.45	0/9376	0.98	5/12853 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	P	6	DG	C2'-C3'-O3'	5.63	119.94	111.50
1	A	381	PHE	CB-CA-C	5.40	118.45	109.38
1	A	975	ASP	CA-CB-CG	5.33	117.93	112.60
3	T	15	DT	C2'-C3'-O3'	5.33	119.49	111.50
2	P	6	DG	C4'-C3'-O3'	5.15	117.72	110.00

There are no chirality outliers.

There are no planarity outliers.

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	8608	0	7997	331	0
2	P	221	0	125	7	0
3	T	308	0	171	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	8	0	0	0	0
5	A	30	0	12	4	0
6	A	1	0	0	0	0
All	All	9176	0	8305	340	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 (340) 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:553:TYR:CE1	1:A:835:MET:HE3	1.82	1.12
1:A:553:TYR:HE1	1:A:835:MET:HE3	1.06	1.10
1:A:155:LEU:HD22	1:A:235:ALA:HB1	1.45	0.98
1:A:652:ASN:HD22	1:A:681:LEU:HD11	1.27	0.96
1:A:576:ILE:HG21	1:A:624:LEU:HD21	1.49	0.94
1:A:863:ALA:O	1:A:867:ARG:HG3	1.69	0.92
1:A:296:PRO:HG2	1:A:299:LYS:HB2	1.52	0.89
1:A:332:ILE:HG23	1:A:349:ILE:HG21	1.54	0.88
1:A:853:THR:O	1:A:857:ILE:HG13	1.74	0.88
1:A:1056:LEU:HB2	1:A:1082:VAL:HG11	1.55	0.88
1:A:860:MET:HE2	1:A:918:PHE:HE2	1.43	0.84
1:A:760:ALA:HB3	1:A:849:ILE:HD11	1.59	0.83
1:A:553:TYR:CE1	1:A:835:MET:CE	2.61	0.83
1:A:1002:ILE:O	1:A:1005:VAL:HG12	1.79	0.82
1:A:285:VAL:HG23	1:A:371:PRO:HA	1.61	0.81
1:A:581:ILE:HD13	1:A:624:LEU:HD22	1.60	0.80
1:A:906:PRO:O	1:A:910:LEU:HD12	1.82	0.80
1:A:685:TRP:CE3	1:A:756:VAL:HG12	2.19	0.77
1:A:76:VAL:HG11	1:A:278:LYS:HE3	1.67	0.77
1:A:182:LEU:HD11	1:A:669:ASP:HB2	1.66	0.77
1:A:953:MET:HG3	1:A:971:VAL:HG22	1.66	0.77
1:A:114:ILE:HD12	1:A:134:VAL:HG11	1.67	0.76
1:A:696:GLU:O	1:A:700:ILE:HG12	1.87	0.74
1:A:539:ILE:HG21	1:A:728:TYR:CE2	2.22	0.74
1:A:749:ARG:HD2	1:A:752:VAL:HG22	1.68	0.74
1:A:553:TYR:HE2	1:A:851:CYS:HG	1.35	0.73
1:A:887:PHE:CD1	1:A:888:PRO:HD2	2.23	0.73
1:A:1181:LYS:O	1:A:1184:ILE:HG13	1.88	0.73
1:A:679:ARG:HH21	1:A:920:ASN:HA	1.54	0.72
1:A:293:THR:HG22	1:A:294:THR:O	1.88	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:653:ARG:O	1:A:653:ARG:HG2	1.89	0.71
1:A:1019:VAL:HG12	1:A:1158:ILE:HD12	1.72	0.71
1:A:1056:LEU:CB	1:A:1082:VAL:HG11	2.21	0.71
1:A:834:VAL:HG21	1:A:846:MET:HE2	1.72	0.70
1:A:956:PRO:HG3	1:A:984:PHE:HZ	1.56	0.70
1:A:653:ARG:NH1	1:A:766:GLU:O	2.24	0.69
1:A:709:PHE:CD1	1:A:710:PRO:HD3	2.27	0.69
1:A:588:LEU:O	1:A:588:LEU:HD23	1.93	0.68
1:A:332:ILE:CG2	1:A:349:ILE:HG21	2.24	0.67
1:A:496:LEU:HG	1:A:500:ILE:HD12	1.75	0.67
1:A:860:MET:CE	1:A:918:PHE:HE2	2.08	0.67
1:A:315:ILE:HG21	1:A:369:VAL:HG21	1.77	0.67
1:A:547:LEU:HD22	1:A:690:PHE:HE1	1.60	0.67
1:A:293:THR:HG21	1:A:301:PRO:HB3	1.74	0.67
1:A:1019:VAL:CG1	1:A:1158:ILE:HD12	2.24	0.66
1:A:792:LYS:HA	1:A:813:ILE:HD11	1.77	0.66
1:A:1093:ILE:HD11	1:A:1145:TYR:CE1	2.31	0.65
1:A:709:PHE:HB3	1:A:710:PRO:CD	2.27	0.64
1:A:171:ARG:HB2	1:A:171:ARG:HH21	1.61	0.64
1:A:709:PHE:CG	1:A:710:PRO:HD3	2.32	0.64
1:A:1096:LYS:O	1:A:1127:THR:HG22	1.98	0.64
1:A:576:ILE:HG21	1:A:624:LEU:CD2	2.28	0.63
1:A:1031:ASP:OD2	1:A:1173:ARG:NH2	2.30	0.63
1:A:337:TYR:CE2	1:A:479:VAL:HG11	2.34	0.63
1:A:630:ARG:HG2	1:A:631:ASN:H	1.64	0.63
1:A:295:LYS:HD2	1:A:299:LYS:O	1.99	0.62
1:A:76:VAL:CG1	1:A:278:LYS:HE3	2.29	0.62
1:A:862:ARG:HA	1:A:865:VAL:HG22	1.81	0.62
1:A:731:GLN:O	1:A:735:ILE:HG13	2.00	0.61
1:A:641:VAL:HG22	1:A:644:MET:HB2	1.81	0.61
1:A:139:PHE:CE2	1:A:242:ILE:HD11	2.36	0.61
1:A:1116:ILE:N	1:A:1117:PRO:HD2	2.15	0.61
1:A:1020:ALA:HB1	1:A:1171:VAL:HG22	1.83	0.61
1:A:685:TRP:HE3	1:A:756:VAL:HG12	1.64	0.60
1:A:709:PHE:CB	1:A:710:PRO:HD3	2.31	0.60
1:A:707:GLU:HG2	1:A:708:THR:H	1.67	0.60
1:A:781:ARG:NH1	1:A:820:GLN:OE1	2.34	0.60
1:A:990:GLY:CA	1:A:1049:ASN:O	2.49	0.60
1:A:337:TYR:HE2	1:A:479:VAL:HG11	1.66	0.60
2:P:10:DT:H2''	2:P:11:DOC:H5'	1.84	0.59
1:A:707:GLU:HG3	1:A:721:LEU:H	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:639:VAL:HG12	1:A:946:VAL:HG12	1.84	0.59
1:A:652:ASN:ND2	1:A:681:LEU:HD11	2.10	0.59
1:A:989:ARG:NH2	2:P:8:DG:OP2	2.35	0.58
1:A:286:VAL:CG2	1:A:482:TYR:HE1	2.17	0.58
1:A:990:GLY:HA3	1:A:1049:ASN:O	2.03	0.58
1:A:374:ILE:HG13	1:A:417:CYS:SG	2.44	0.58
1:A:887:PHE:CG	1:A:888:PRO:HD2	2.37	0.58
1:A:845:GLU:O	1:A:849:ILE:HG13	2.04	0.58
1:A:968:ARG:HE	3:T:10:DG:H4'	1.68	0.58
1:A:791:TRP:HA	1:A:794:ASN:HB2	1.85	0.58
1:A:233:VAL:HG12	1:A:234:ASP:H	1.69	0.57
1:A:999:GLN:HA	1:A:1002:ILE:HG22	1.86	0.57
1:A:956:PRO:HG3	1:A:984:PHE:CZ	2.37	0.57
1:A:824:LYS:NZ	5:A:1202:DTP:O3B	2.38	0.57
1:A:310:MET:SD	1:A:470:LEU:HD23	2.44	0.57
1:A:332:ILE:HG22	1:A:351:ASN:OD1	2.04	0.57
1:A:722:THR:H	1:A:725:GLU:HB2	1.68	0.57
1:A:611:GLU:C	1:A:613:ILE:H	2.11	0.57
1:A:141:ILE:HD11	1:A:191:LEU:HD21	1.86	0.57
1:A:942:ILE:C	1:A:943:PHE:HD1	2.13	0.57
1:A:969:TYR:CZ	1:A:982:LYS:HG3	2.40	0.57
1:A:910:LEU:O	1:A:914:VAL:HG23	2.05	0.56
3:T:8:DA:H1'	3:T:9:DC:H5''	1.87	0.56
1:A:350:PHE:HE2	1:A:361:ARG:HE	1.52	0.56
1:A:123:ASP:OD1	1:A:126:GLY:N	2.39	0.56
1:A:989:ARG:HH21	2:P:8:DG:P	2.29	0.56
1:A:382:PHE:O	1:A:385:PRO:HD2	2.06	0.55
1:A:1159:THR:O	1:A:1162:ALA:HB3	2.07	0.55
1:A:340:LYS:HD2	1:A:448:GLY:O	2.07	0.55
1:A:296:PRO:HG2	1:A:299:LYS:CB	2.33	0.55
1:A:709:PHE:CB	1:A:710:PRO:CD	2.85	0.55
1:A:547:LEU:HD22	1:A:690:PHE:CE1	2.39	0.55
1:A:288:ALA:HB2	1:A:375:SER:HB3	1.88	0.55
1:A:947:ASP:HB3	1:A:950:TYR:OH	2.07	0.55
1:A:171:ARG:HB2	1:A:171:ARG:NH2	2.23	0.54
1:A:679:ARG:HH11	1:A:679:ARG:CG	2.20	0.54
1:A:1095:SER:HA	1:A:1142:ASP:HB2	1.90	0.54
1:A:580:ALA:HA	1:A:867:ARG:HH11	1.73	0.54
1:A:870:ARG:HB2	1:A:882:ILE:HD11	1.89	0.54
1:A:170:ILE:HD12	1:A:190:LYS:HD2	1.90	0.54
1:A:778:ARG:NH1	1:A:779:ASP:OD1	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:433:PRO:HD2	1:A:436:SER:HB2	1.90	0.53
1:A:961:GLU:HG3	1:A:1179:TRP:CE3	2.44	0.53
1:A:1150:LEU:O	1:A:1154:ILE:HG12	2.06	0.53
1:A:286:VAL:HG23	1:A:482:TYR:HE1	1.74	0.53
1:A:858:ILE:HD11	1:A:879:ILE:HG23	1.90	0.53
1:A:138:TYR:HA	1:A:191:LEU:O	2.09	0.53
1:A:685:TRP:CE3	1:A:756:VAL:CG1	2.91	0.53
1:A:316:ASP:HA	1:A:482:TYR:OH	2.08	0.53
1:A:333:GLU:C	1:A:349:ILE:HD12	2.34	0.53
1:A:553:TYR:HE2	1:A:851:CYS:SG	2.30	0.53
1:A:595:SER:O	1:A:599:GLU:HB2	2.09	0.53
1:A:813:ILE:O	1:A:817:ASP:HB2	2.08	0.53
1:A:163:CYS:CB	1:A:198:GLN:HE21	2.21	0.53
1:A:323:THR:HB	1:A:351:ASN:HA	1.91	0.53
1:A:1081:MET:HA	1:A:1081:MET:HE2	1.90	0.53
1:A:519:LEU:O	1:A:522:VAL:HG22	2.09	0.53
1:A:1107:ILE:HG12	1:A:1126:TRP:CE3	2.44	0.53
1:A:539:ILE:CG2	1:A:728:TYR:CE2	2.92	0.52
1:A:722:THR:HG22	1:A:723:PHE:N	2.25	0.52
1:A:581:ILE:HD11	1:A:621:LEU:O	2.09	0.52
1:A:1107:ILE:CG1	1:A:1126:TRP:CE3	2.92	0.52
1:A:460:THR:HB	1:A:461:PRO:HD3	1.92	0.52
1:A:793:GLY:HA2	1:A:797:LYS:CB	2.40	0.51
1:A:636:ILE:HG12	1:A:882:ILE:HG22	1.92	0.51
1:A:594:PHE:CE2	1:A:915:HIS:CD2	2.98	0.51
1:A:291:ILE:HG21	1:A:387:ILE:HD11	1.93	0.51
1:A:145:ASP:C	1:A:147:SER:H	2.19	0.51
1:A:722:THR:HG22	1:A:723:PHE:H	1.75	0.50
1:A:1092:ILE:C	1:A:1092:ILE:HD12	2.37	0.50
1:A:170:ILE:HG22	1:A:171:ARG:N	2.26	0.50
1:A:835:MET:HE1	1:A:844:MET:HG3	1.94	0.50
1:A:1136:ASP:O	1:A:1140:ILE:HD12	2.11	0.50
1:A:245:TYR:HE1	1:A:533:ASN:OD1	1.94	0.50
1:A:251:VAL:O	1:A:255:ILE:HG13	2.11	0.50
1:A:919:THR:HA	1:A:940:ASN:HD22	1.76	0.50
1:A:1181:LYS:C	1:A:1183:LYS:H	2.19	0.50
1:A:382:PHE:C	1:A:385:PRO:HD2	2.37	0.50
1:A:88:SER:O	1:A:90:GLU:HG2	2.12	0.49
1:A:287:MET:HB2	1:A:315:ILE:HG12	1.93	0.49
1:A:1024:ASN:OD1	1:A:1173:ARG:HD3	2.11	0.49
1:A:1100:ALA:O	1:A:1105:ARG:NH1	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1169:ASN:OD1	1:A:1169:ASN:C	2.55	0.49
1:A:287:MET:HA	1:A:314:MET:O	2.13	0.49
1:A:309:MET:HG2	1:A:327:ILE:HG21	1.94	0.49
1:A:707:GLU:CD	1:A:707:GLU:H	2.20	0.49
1:A:1023:CYS:HB3	1:A:1154:ILE:HD12	1.95	0.49
1:A:350:PHE:HD2	1:A:361:ARG:HG2	1.77	0.49
1:A:141:ILE:HD12	1:A:191:LEU:HD11	1.95	0.49
1:A:490:HIS:HB3	1:A:491:PRO:CD	2.43	0.49
1:A:594:PHE:HE2	1:A:915:HIS:CD2	2.31	0.49
1:A:928:ASP:OD2	1:A:931:ASN:HB2	2.13	0.49
1:A:72:ILE:O	1:A:266:LYS:NZ	2.28	0.49
1:A:152:VAL:HA	1:A:155:LEU:HD23	1.94	0.49
1:A:553:TYR:CD1	1:A:835:MET:CE	2.96	0.49
1:A:1037:LEU:HD23	1:A:1042:LEU:HD12	1.94	0.49
1:A:264:TRP:CD2	1:A:278:LYS:HD2	2.48	0.48
1:A:795:LEU:HA	1:A:799:ASP:CB	2.43	0.48
1:A:906:PRO:O	1:A:910:LEU:CD1	2.57	0.48
1:A:1040:GLU:O	1:A:1043:VAL:HG22	2.14	0.48
1:A:286:VAL:HG21	1:A:486:MET:SD	2.53	0.48
1:A:332:ILE:HD12	1:A:471:SER:CB	2.43	0.48
1:A:968:ARG:O	1:A:969:TYR:HB3	2.14	0.48
1:A:490:HIS:HB3	1:A:491:PRO:HD3	1.96	0.48
1:A:799:ASP:N	1:A:800:PRO:CD	2.77	0.48
1:A:672:ARG:CB	1:A:673:PRO:HD2	2.44	0.48
1:A:1011:THR:HG23	1:A:1014:GLY:H	1.79	0.48
1:A:210:LEU:HD21	1:A:241:ASP:HA	1.94	0.48
1:A:915:HIS:CE1	1:A:940:ASN:OD1	2.67	0.48
2:P:4:DC:H2'	2:P:5:DC:C6	2.49	0.47
1:A:986:LEU:HD13	1:A:1000:SER:HG	1.78	0.47
1:A:1011:THR:O	1:A:1015:CYS:N	2.47	0.47
1:A:549:GLU:HB3	1:A:688:GLU:O	2.14	0.47
1:A:981:LEU:CD2	1:A:986:LEU:HD23	2.45	0.47
1:A:438:GLY:O	1:A:442:VAL:HG23	2.14	0.47
1:A:572:ASN:HB3	1:A:870:ARG:HH21	1.78	0.47
1:A:865:VAL:HA	1:A:868:VAL:HG12	1.97	0.47
1:A:912:TYR:C	1:A:912:TYR:CD2	2.93	0.47
1:A:155:LEU:HD22	1:A:235:ALA:CB	2.30	0.47
1:A:324:ASN:ND2	1:A:327:ILE:HD13	2.29	0.47
1:A:709:PHE:HB3	1:A:710:PRO:HD3	1.95	0.47
1:A:869:GLY:HA3	1:A:883:LEU:CD2	2.45	0.47
1:A:522:VAL:O	1:A:526:GLN:HG3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:LEU:HD23	1:A:40:ALA:HB3	1.97	0.46
1:A:302:ASP:O	1:A:306:ASP:HB2	2.15	0.46
1:A:651:THR:HG23	1:A:939:GLU:O	2.15	0.46
1:A:790:THR:O	1:A:794:ASN:HB2	2.15	0.46
1:A:645:TYR:CG	5:A:1202:DTP:H2'1	2.50	0.46
1:A:649:MET:HA	1:A:654:LEU:HD12	1.97	0.46
1:A:679:ARG:HH11	1:A:679:ARG:HG2	1.80	0.46
1:A:1048:GLU:OE2	1:A:1050:ARG:NH2	2.47	0.46
1:A:151:ASP:O	1:A:155:LEU:HD23	2.16	0.46
1:A:860:MET:HE2	1:A:918:PHE:CE2	2.34	0.46
1:A:989:ARG:NH2	2:P:8:DG:P	2.88	0.46
1:A:990:GLY:HA2	1:A:1049:ASN:O	2.13	0.46
1:A:332:ILE:HD12	1:A:471:SER:HB3	1.96	0.46
1:A:465:GLU:O	1:A:467:PRO:HD3	2.15	0.46
1:A:567:ARG:HE	1:A:570:LEU:HD22	1.79	0.46
1:A:588:LEU:N	1:A:589:PRO:CD	2.78	0.46
1:A:1155:GLN:HA	1:A:1159:THR:HB	1.97	0.46
3:T:2:DT:H2''	3:T:3:DC:O5'	2.15	0.46
1:A:737:LYS:O	1:A:740:THR:HG22	2.15	0.46
1:A:1078:GLY:O	1:A:1081:MET:HB2	2.16	0.46
1:A:243:ARG:NH1	1:A:531:LEU:O	2.48	0.46
1:A:1179:TRP:O	1:A:1183:LYS:HB2	2.16	0.46
1:A:55:TYR:OH	1:A:74:GLY:HA2	2.15	0.45
1:A:170:ILE:CG2	1:A:171:ARG:N	2.80	0.45
1:A:191:LEU:N	1:A:191:LEU:HD12	2.31	0.45
1:A:1116:ILE:HA	1:A:1119:LYS:HB3	1.98	0.45
1:A:689:PHE:CE2	1:A:752:VAL:HB	2.51	0.45
1:A:135:TYR:OH	1:A:256:ASP:OD2	2.33	0.45
1:A:308:ILE:O	1:A:308:ILE:HG13	2.16	0.45
1:A:645:TYR:HD2	1:A:831:TYR:CD2	2.35	0.45
1:A:905:TYR:N	1:A:906:PRO:CD	2.80	0.45
1:A:46:ILE:O	1:A:49:MET:HB3	2.16	0.45
1:A:114:ILE:HG22	1:A:194:VAL:HB	1.98	0.45
1:A:645:TYR:CD2	5:A:1202:DTP:H2'1	2.52	0.45
1:A:165:LYS:HD2	1:A:194:VAL:HG12	1.99	0.45
1:A:260:ARG:HB2	1:A:263:LYS:HD2	1.99	0.45
1:A:806:ARG:C	1:A:808:GLU:H	2.25	0.45
1:A:332:ILE:HD12	1:A:471:SER:OG	2.17	0.45
1:A:833:TYR:C	1:A:835:MET:H	2.24	0.45
1:A:561:LEU:HD13	1:A:871:PRO:HG2	1.99	0.44
1:A:1115:ASP:O	1:A:1118:ILE:HG22	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:244:GLU:HG2	1:A:531:LEU:HB2	1.98	0.44
1:A:332:ILE:HD13	1:A:349:ILE:HD13	1.98	0.44
1:A:906:PRO:O	1:A:909:MET:HB3	2.17	0.44
1:A:583:GLU:OE1	1:A:913:ARG:NH2	2.45	0.44
1:A:1169:ASN:OD1	1:A:1171:VAL:N	2.45	0.44
1:A:46:ILE:HD13	1:A:403:GLY:HA2	1.99	0.44
1:A:138:TYR:CD1	1:A:138:TYR:C	2.95	0.44
1:A:148:ARG:HG2	1:A:151:ASP:HB2	1.99	0.44
1:A:611:GLU:C	1:A:613:ILE:N	2.75	0.44
1:A:163:CYS:HB3	1:A:198:GLN:HE21	1.82	0.44
1:A:182:LEU:CD1	1:A:669:ASP:HB2	2.43	0.44
1:A:865:VAL:O	1:A:869:GLY:N	2.50	0.44
1:A:193:PHE:HB3	1:A:198:GLN:HB3	1.99	0.44
1:A:288:ALA:CB	1:A:375:SER:HB3	2.47	0.44
1:A:609:ASN:C	1:A:611:GLU:N	2.76	0.44
1:A:791:TRP:HA	1:A:794:ASN:CB	2.47	0.44
1:A:160:LEU:HD22	1:A:163:CYS:SG	2.57	0.44
1:A:598:VAL:HG23	1:A:599:GLU:N	2.33	0.44
1:A:539:ILE:HG21	1:A:728:TYR:HE2	1.79	0.44
3:T:2:DT:H2"	3:T:3:DC:C6	2.53	0.43
1:A:985:GLU:HG3	1:A:991:GLU:OE1	2.18	0.43
1:A:275:ASP:OD2	1:A:278:LYS:NZ	2.49	0.43
1:A:303:SER:HB3	1:A:389:ASN:HB3	2.00	0.43
1:A:643:SER:N	5:A:1202:DTP:O1G	2.42	0.43
1:A:202:ALA:O	1:A:206:LEU:CD1	2.67	0.43
1:A:700:ILE:HG22	1:A:700:ILE:O	2.18	0.43
1:A:1168:SER:O	1:A:1169:ASN:C	2.61	0.43
1:A:367:ARG:HH12	1:A:401:GLU:CD	2.26	0.43
1:A:549:GLU:N	1:A:688:GLU:O	2.41	0.43
1:A:490:HIS:N	1:A:491:PRO:HD2	2.34	0.43
1:A:609:ASN:C	1:A:611:GLU:H	2.25	0.43
1:A:639:VAL:HG22	1:A:879:ILE:HG13	2.00	0.43
1:A:293:THR:O	1:A:309:MET:HE2	2.19	0.43
1:A:338:THR:HG23	1:A:344:PRO:HA	2.00	0.43
1:A:700:ILE:HG21	1:A:739:LEU:HA	1.99	0.43
1:A:404:PHE:CD2	1:A:414:SER:HB3	2.54	0.43
1:A:446:LYS:HA	1:A:446:LYS:HD3	1.85	0.43
1:A:580:ALA:HB2	1:A:867:ARG:HE	1.83	0.43
1:A:609:ASN:O	1:A:611:GLU:N	2.51	0.43
1:A:1157:ILE:O	1:A:1161:PRO:HG2	2.19	0.43
1:A:687:GLY:O	1:A:753:SER:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:T:15:DT:H2''	3:T:16:DA:OP2	2.19	0.43
1:A:36:GLN:OE1	1:A:197:ASN:ND2	2.46	0.42
1:A:252:ARG:HD3	1:A:252:ARG:HA	1.72	0.42
1:A:201:GLU:O	1:A:204:LYS:HB2	2.18	0.42
1:A:514:THR:OG1	1:A:836:ARG:NH2	2.51	0.42
1:A:864:LEU:HD13	1:A:913:ARG:HH12	1.83	0.42
1:A:993:GLN:O	1:A:993:GLN:HG3	2.19	0.42
1:A:50:MET:HE2	1:A:416:TYR:CE2	2.54	0.42
1:A:926:LEU:HD21	1:A:929:PRO:HA	2.01	0.42
1:A:590:GLU:CB	1:A:912:TYR:CE1	3.02	0.42
1:A:958:SER:HB3	1:A:963:LYS:O	2.20	0.42
1:A:500:ILE:HD13	1:A:515:LEU:HD22	2.02	0.42
1:A:174:ASP:OD1	1:A:174:ASP:C	2.62	0.42
1:A:431:TYR:CE1	1:A:824:LYS:HD3	2.54	0.42
1:A:1077:LEU:HB3	1:A:1081:MET:HG3	2.01	0.42
1:A:864:LEU:HD13	1:A:913:ARG:NH1	2.34	0.42
1:A:1151:GLY:O	1:A:1155:GLN:HG3	2.19	0.42
1:A:121:PHE:HZ	1:A:272:PHE:CE2	2.37	0.42
1:A:202:ALA:O	1:A:206:LEU:HD12	2.20	0.42
1:A:275:ASP:C	1:A:277:ARG:H	2.28	0.42
1:A:653:ARG:HB2	1:A:923:TYR:CE2	2.54	0.42
1:A:1016:TYR:CZ	1:A:1162:ALA:HA	2.54	0.42
1:A:707:GLU:HG2	1:A:708:THR:N	2.33	0.42
3:T:9:DC:H2''	3:T:10:DG:C8	2.54	0.42
1:A:778:ARG:HD3	1:A:782:TYR:CZ	2.55	0.41
1:A:882:ILE:H	1:A:882:ILE:HG13	1.63	0.41
1:A:87:VAL:O	1:A:114:ILE:HG12	2.21	0.41
1:A:620:LYS:O	1:A:624:LEU:HB2	2.20	0.41
1:A:241:ASP:OD1	1:A:242:ILE:N	2.45	0.41
1:A:407:ASP:C	1:A:409:GLU:H	2.29	0.41
1:A:426:VAL:HG12	1:A:437:GLN:HG2	2.02	0.41
1:A:1115:ASP:O	1:A:1119:LYS:HB2	2.20	0.41
1:A:247:VAL:O	1:A:248:PRO:C	2.62	0.41
1:A:584:LEU:HD11	1:A:864:LEU:HD11	2.01	0.41
1:A:33:SER:O	1:A:37:LEU:HB2	2.21	0.41
1:A:420:MET:HE2	1:A:420:MET:HA	2.02	0.41
1:A:627:ASN:OD1	1:A:627:ASN:N	2.53	0.41
1:A:1107:ILE:HD11	1:A:1126:TRP:HE3	1.85	0.41
3:T:2:DT:O2	3:T:2:DT:H2'	2.20	0.41
1:A:160:LEU:HD11	1:A:206:LEU:HD11	2.01	0.41
1:A:588:LEU:HD23	1:A:588:LEU:C	2.45	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:709:PHE:HD1	1:A:709:PHE:HA	1.76	0.41
1:A:58:PRO:HG2	1:A:73:PRO:HB2	2.02	0.41
1:A:433:PRO:HB3	2:P:9:DT:H5'	2.01	0.41
3:T:14:DT:H2'	3:T:15:DT:H72	2.03	0.41
1:A:207:ARG:N	1:A:208:PRO:CD	2.83	0.41
1:A:546:HIS:HB3	1:A:690:PHE:O	2.21	0.41
1:A:404:PHE:CE2	1:A:414:SER:HB3	2.56	0.41
1:A:986:LEU:HD13	1:A:1000:SER:OG	2.21	0.41
1:A:1181:LYS:C	1:A:1183:LYS:N	2.79	0.41
1:A:325:ARG:HA	1:A:325:ARG:HD2	1.81	0.41
1:A:529:ILE:HD11	1:A:767:ASN:ND2	2.36	0.41
1:A:967:LYS:NZ	2:P:10:DT:O2	2.52	0.41
1:A:1092:ILE:HG13	1:A:1107:ILE:HG22	2.02	0.41
1:A:685:TRP:CZ3	1:A:756:VAL:HG12	2.55	0.40
1:A:724:ASP:HA	1:A:731:GLN:NE2	2.36	0.40
1:A:806:ARG:HD3	1:A:808:GLU:HB2	2.03	0.40
1:A:707:GLU:CG	1:A:708:THR:H	2.31	0.40
1:A:761:ILE:HD12	1:A:761:ILE:N	2.36	0.40
1:A:120:TYR:CD1	1:A:120:TYR:N	2.90	0.40
1:A:815:LEU:HD22	1:A:815:LEU:HA	1.88	0.40
1:A:860:MET:CE	1:A:918:PHE:CE2	2.97	0.40
1:A:580:ALA:CA	1:A:867:ARG:HH11	2.34	0.40
1:A:36:GLN:CD	1:A:197:ASN:HD21	2.26	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	1115/1192 (94%)	998 (90%)	115 (10%)	2 (0%)	43 72

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	889	GLU
1	A	610	PHE

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	869/1066 (82%)	815 (94%)	54 (6%)	16 45

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	LEU
1	A	85	THR
1	A	87	VAL
1	A	155	LEU
1	A	162	SER
1	A	166	SER
1	A	171	ARG
1	A	174	ASP
1	A	188	LEU
1	A	196	SER
1	A	225	VAL
1	A	233	VAL
1	A	240	GLU
1	A	286	VAL
1	A	305	VAL
1	A	325	ARG
1	A	331	ASP
1	A	332	ILE
1	A	356	VAL
1	A	418	SER
1	A	479	VAL
1	A	523	GLN
1	A	539	ILE
1	A	581	ILE
1	A	627	ASN

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Mol	Chain	Res	Type
1	A	640	ASP
1	A	641	VAL
1	A	679	ARG
1	A	706	ASN
1	A	709	PHE
1	A	732	VAL
1	A	740	THR
1	A	756	VAL
1	A	781	ARG
1	A	808	GLU
1	A	815	LEU
1	A	818	SER
1	A	820	GLN
1	A	879	ILE
1	A	882	ILE
1	A	902	TYR
1	A	904	SER
1	A	936	THR
1	A	980	GLU
1	A	1001	ASP
1	A	1035	LEU
1	A	1077	LEU
1	A	1093	ILE
1	A	1107	ILE
1	A	1111	ILE
1	A	1118	ILE
1	A	1148	GLU
1	A	1158	ILE
1	A	1164	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	144	ASN
1	A	185	GLN
1	A	198	GLN
1	A	217	ASN
1	A	389	ASN
1	A	394	HIS
1	A	468	GLN
1	A	546	HIS
1	A	652	ASN

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Mol	Chain	Res	Type
1	A	767	ASN
1	A	859	GLN
1	A	915	HIS
1	A	922	GLN
1	A	999	GLN
1	A	1033	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	DOC	P	11	2,3	16,19,20	1.56	3 (18%)	20,26,29	1.44	2 (10%)

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	DOC	P	11	2,3	-	2/7/18/19	0/2/2/2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	P	11	DOC	C4-N3	3.11	1.40	1.34
2	P	11	DOC	C2-N1	-2.52	1.34	1.40
2	P	11	DOC	C2-N3	2.38	1.41	1.36

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	P	11	DOC	O4'-C4'-C5'	-4.41	101.38	109.34
2	P	11	DOC	C3'-C2'-C1'	2.23	105.44	102.87

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	P	11	DOC	C3'-C4'-C5'-O5'
2	P	11	DOC	O4'-C4'-C5'-O5'

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	P	11	DOC	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	DTP	A	1202	6	31,32,32	1.73	7 (22%)	46,50,50	2.11	15 (32%)
4	SF4	A	1201	1	0,12,12	-	-	-		

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
5	DTP	A	1202	6	-	4/22/34/34	0/3/3/3
4	SF4	A	1201	1	-	-	0/6/5/5

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1202	DTP	C2-N1	4.50	1.41	1.33
5	A	1202	DTP	C2-N3	3.40	1.40	1.33
5	A	1202	DTP	PB-O3B	3.17	1.62	1.59
5	A	1202	DTP	C5-C6	-2.91	1.32	1.41
5	A	1202	DTP	C5-N7	-2.76	1.34	1.39
5	A	1202	DTP	C5-C4	-2.45	1.34	1.39
5	A	1202	DTP	C4-N9	-2.28	1.33	1.37

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1202	DTP	N3-C2-N1	-7.11	117.83	128.58
5	A	1202	DTP	N9-C8-N7	-3.99	108.27	113.94
5	A	1202	DTP	O4'-C1'-N9	-3.78	101.18	107.96
5	A	1202	DTP	C5-C4-N3	-3.62	121.73	126.72
5	A	1202	DTP	O3B-PG-O1G	-3.38	93.23	111.04
5	A	1202	DTP	C5-C6-N6	-3.16	115.46	123.29
5	A	1202	DTP	C2-N3-C4	3.04	119.25	111.83
5	A	1202	DTP	C4-N9-C8	2.76	108.64	105.74
5	A	1202	DTP	C2'-C1'-N9	-2.62	108.48	114.63
5	A	1202	DTP	O2B-PB-O3B	2.51	114.05	107.27
5	A	1202	DTP	C2'-C3'-C4'	-2.40	97.92	102.80
5	A	1202	DTP	N3-C4-N9	2.39	131.24	127.17
5	A	1202	DTP	C6-C5-C4	2.21	120.20	117.18
5	A	1202	DTP	C5-N7-C8	2.16	106.85	103.45
5	A	1202	DTP	O3G-PG-O2G	2.08	115.60	107.80

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1202	DTP	O4'-C4'-C5'-O5'
5	A	1202	DTP	C3'-C4'-C5'-O5'
5	A	1202	DTP	PB-O3A-PA-O1A

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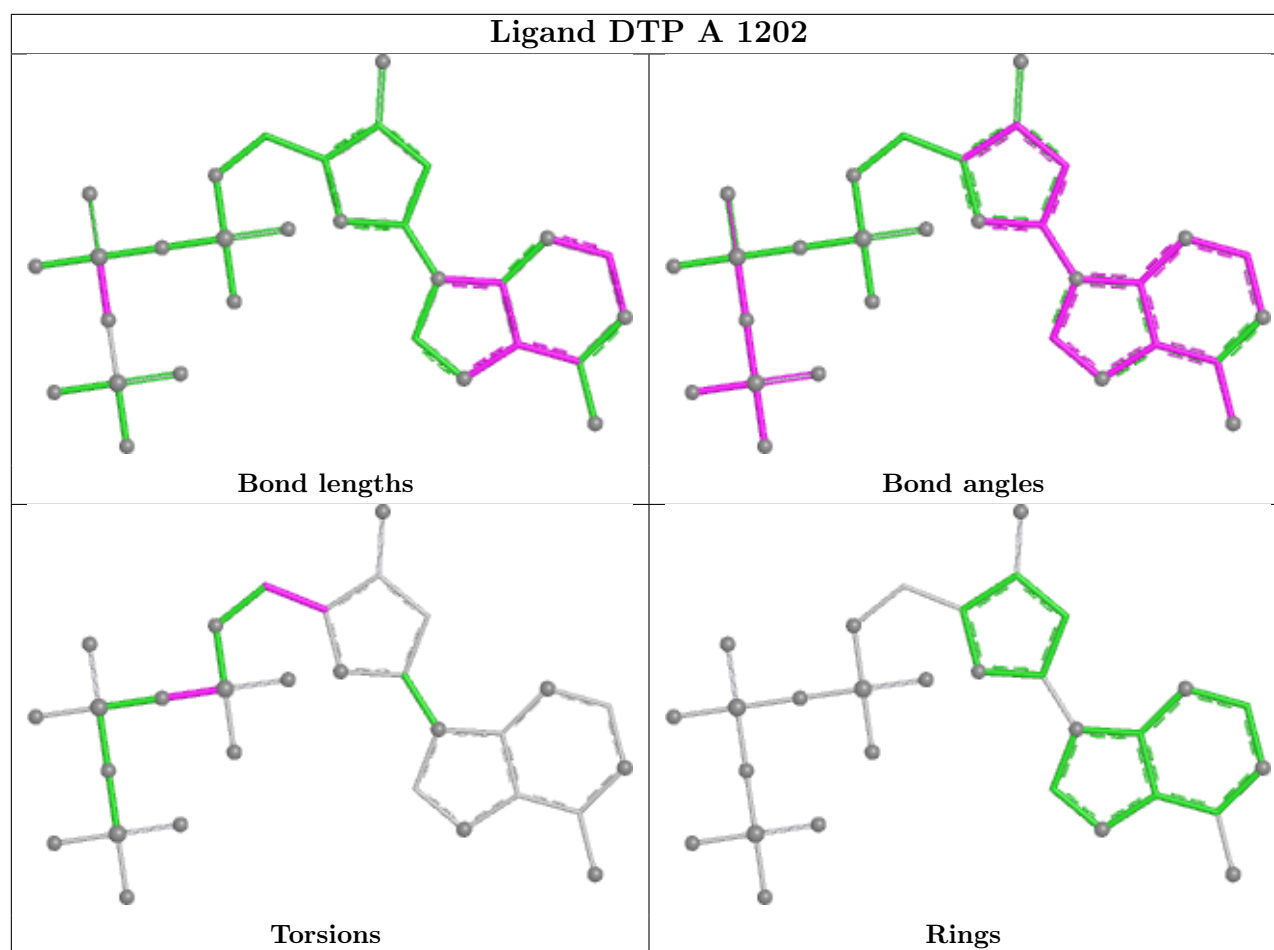
Mol	Chain	Res	Type	Atoms
5	A	1202	DTP	PB-O3A-PA-O2A

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1202	DTP	4	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

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	1123/1192 (94%)	0.70	88 (7%) 19 14	32, 74, 112, 145	0
2	P	10/11 (90%)	-0.22	0 100 100	39, 56, 80, 98	0
3	T	15/16 (93%)	-0.16	0 100 100	34, 54, 86, 103	0
All	All	1148/1219 (94%)	0.68	88 (7%) 19 14	32, 74, 112, 145	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	893	PHE	5.0
1	A	1106	ALA	4.7
1	A	1185	ALA	4.5
1	A	537	ASP	4.4
1	A	847	ALA	4.4
1	A	1099	ASN	3.8
1	A	742	TYR	3.6
1	A	750	VAL	3.6
1	A	899	LYS	3.5
1	A	740	THR	3.5
1	A	710	PRO	3.4
1	A	615	ASN	3.4
1	A	709	PHE	3.4
1	A	897	ASN	3.4
1	A	881	CYS	3.4
1	A	894	THR	3.4
1	A	65	ASP	3.3
1	A	898	GLY	3.3
1	A	31	ALA	3.2
1	A	895	LEU	3.2
1	A	851	CYS	3.2
1	A	63	ARG	3.1
1	A	698	ASN	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	748	HIS	3.1
1	A	671	ASN	3.1
1	A	896	GLU	3.1
1	A	1184	ILE	3.0
1	A	1101	PRO	3.0
1	A	553	TYR	2.9
1	A	802	ASP	2.9
1	A	761	ILE	2.9
1	A	552	THR	2.8
1	A	298	LEU	2.7
1	A	68	ASP	2.7
1	A	1100	ALA	2.7
1	A	279	ILE	2.7
1	A	703	ALA	2.7
1	A	1146	TYR	2.7
1	A	617	ILE	2.6
1	A	844	MET	2.6
1	A	556	GLY	2.6
1	A	990	GLY	2.6
1	A	704	LEU	2.6
1	A	375	SER	2.5
1	A	708	THR	2.5
1	A	1178	ASP	2.5
1	A	675	LYS	2.5
1	A	673	PRO	2.4
1	A	805	ALA	2.4
1	A	795	LEU	2.4
1	A	699	MET	2.4
1	A	901	LEU	2.4
1	A	34	ALA	2.4
1	A	994	LEU	2.4
1	A	900	LYS	2.3
1	A	145	ASP	2.3
1	A	1008	GLU	2.3
1	A	941	THR	2.3
1	A	991	GLU	2.3
1	A	835	MET	2.3
1	A	1035	LEU	2.2
1	A	738	ARG	2.2
1	A	1040	GLU	2.2
1	A	967	LYS	2.2
1	A	395	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	513	GLY	2.2
1	A	741	GLU	2.2
1	A	700	ILE	2.2
1	A	989	ARG	2.2
1	A	842	TYR	2.2
1	A	905	TYR	2.2
1	A	33	SER	2.2
1	A	1098	PHE	2.2
1	A	855	ALA	2.2
1	A	916	GLN	2.1
1	A	225	VAL	2.1
1	A	697	TYR	2.1
1	A	843	SER	2.1
1	A	612	GLU	2.1
1	A	308	ILE	2.1
1	A	859	GLN	2.1
1	A	1082	VAL	2.1
1	A	618	THR	2.1
1	A	1047	CYS	2.1
1	A	796	SER	2.0
1	A	691	PRO	2.0
1	A	384	TRP	2.0
1	A	648	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
2	DOC	P	11	18/19	0.97	0.08	33,37,47,52	0

6.3 Carbohydrates [i](#)

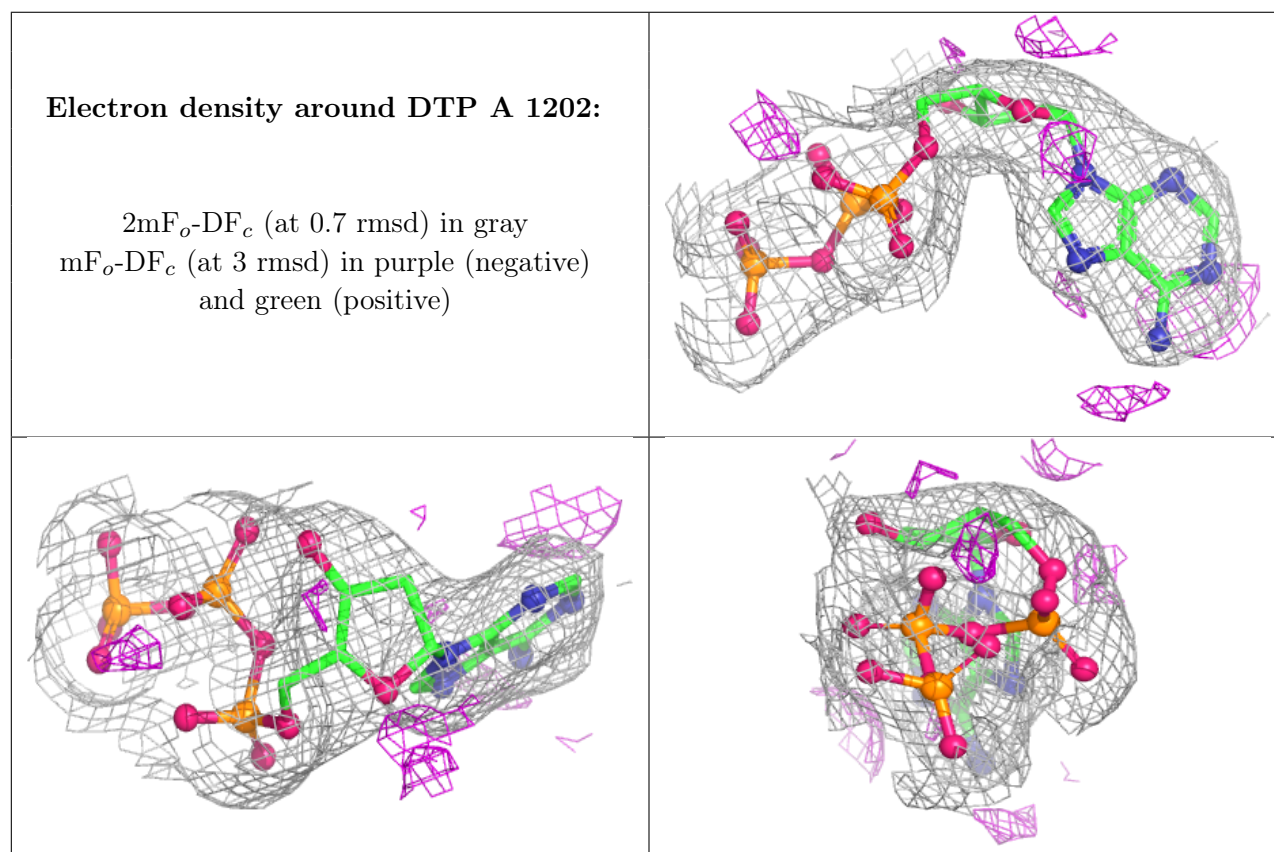
There are no oligosaccharides in this entry.

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($\text{\AA}^2$)	Q<0.9
4	SF4	A	1201	8/8	0.97	0.05	68,77,85,85	8
5	DTP	A	1202	30/30	0.97	0.06	35,40,48,49	0
6	CA	A	1203	1/1	0.98	0.07	46,46,46,46	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.