



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:27 PM UTC

PDB ID : 1G3M / pdb_00001g3m
Title : CRYSTAL STRUCTURE OF HUMAN ESTROGEN SULFOTRANSFERASE IN COMPLEX WITH IN-ACTIVE COFACTOR PAP AND 3,5,3',5'-TETRACHLORO-BIPHENYL-4,4'-DIOL
Authors : Shevtsov, S.; Petrochenko, E.V.; Negishi, M.; Pedersen, L.C.
Deposited on : 2000-10-24
Resolution : 1.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

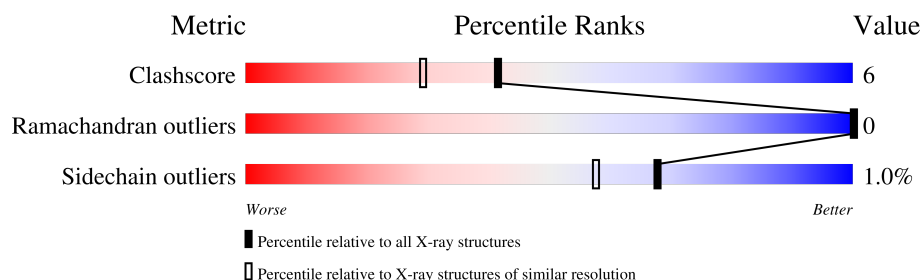
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	294	
1	B	294	

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	EDO	A	801	-	X	-	-
4	EDO	A	802	-	X	-	-
4	EDO	A	804	-	X	-	-

2 Entry composition [i](#)

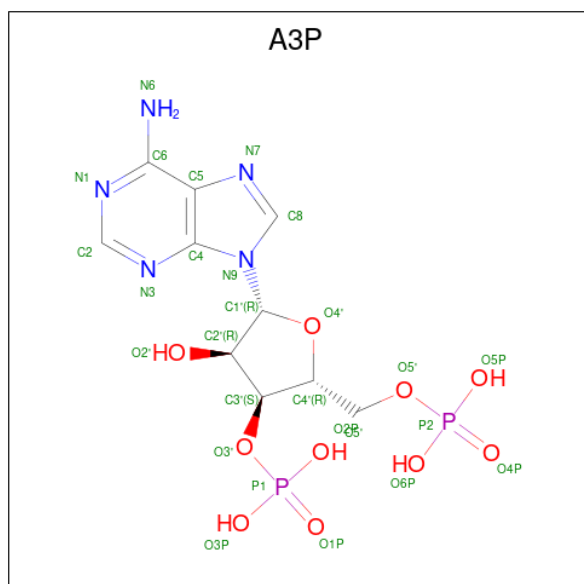
There are 5 unique types of molecules in this entry. The entry contains 5484 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ESTROGEN SULFOTRANSFERASE.

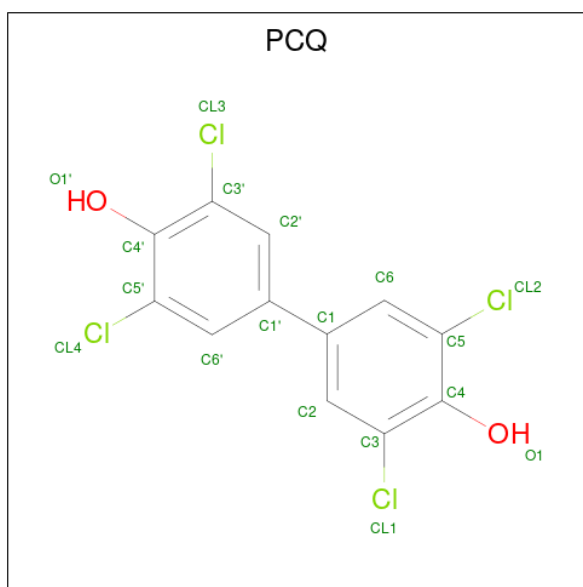
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	290	Total	C	N	O	S	0	0	0
			2422	1577	397	435	13			
1	B	285	Total	C	N	O	S	0	0	0
			2353	1535	387	418	13			

- Molecule 2 is ADENOSINE-3'-5'-DIPHOSPHATE (CCD ID: A3P) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 3 is 3,5,3',5'-TETRACHLORO-BIPHENYL-4,4'-DIOL (CCD ID: PCQ) (formula: $C_{12}H_6Cl_4O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Cl	O	0	0
			18	12	4	2		
3	B	1	Total	C	Cl	O	0	0
			18	12	4	2		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

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

- Molecule 5 is water.

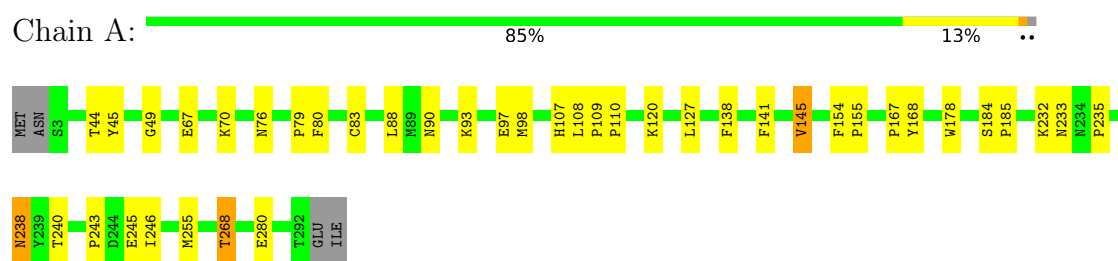
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	317	Total	O	0	0
			317	317		
5	B	286	Total	O	0	0
			286	286		

3 Residue-property plots [i](#)

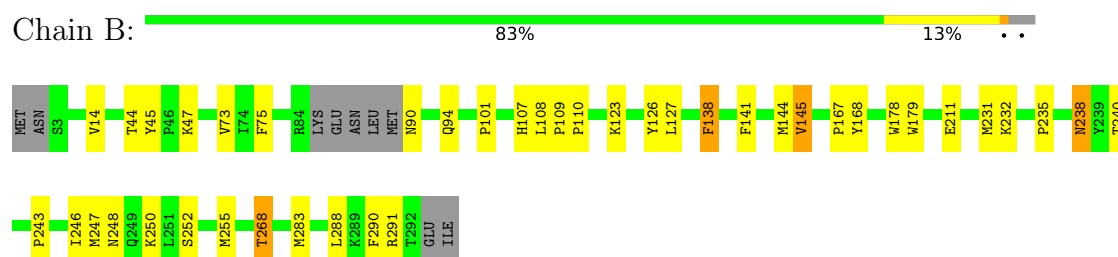
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

• Molecule 1: ESTROGEN SULFOTRANSFERASE



• Molecule 1: ESTROGEN SULFOTRANSFERASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	62.69Å 96.89Å 61.72Å 90.00° 92.58° 90.00°	Depositor
Resolution (Å)	24.53 – 1.70	Depositor
% Data completeness (in resolution range)	91.5 (24.53-1.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.193 , 0.219	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5484	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

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: EDO, PCQ, A3P

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	0.36	0/2495	0.86	6/3374 (0.2%)
1	B	0.37	0/2425	0.87	7/3283 (0.2%)
All	All	0.37	0/4920	0.87	13/6657 (0.2%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	45	TYR	N-CA-C	-8.86	97.31	110.24
1	B	45	TYR	N-CA-C	-8.38	98.01	110.24
1	B	109	PRO	N-CA-C	-6.63	102.61	110.70
1	A	145	VAL	N-CA-C	6.62	117.82	108.36
1	A	268	THR	N-CA-C	-6.00	102.79	110.53
1	B	145	VAL	N-CA-C	5.80	116.66	108.36
1	A	109	PRO	N-CA-C	-5.74	103.70	110.70
1	A	76	ASN	N-CA-C	-5.60	106.58	113.41
1	A	108	LEU	N-CA-C	5.51	116.83	109.83
1	B	283	MET	N-CA-C	5.21	121.36	114.12
1	B	268	THR	N-CA-C	-5.20	103.83	110.53
1	B	250	LYS	N-CA-C	-5.08	105.83	112.23
1	B	108	LEU	N-CA-C	5.08	116.28	109.83

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	2422	0	2354	30	0
1	B	2353	0	2264	31	0
2	A	27	0	11	0	0
2	B	27	0	11	0	0
3	A	18	0	6	1	0
3	B	18	0	5	0	0
4	A	16	0	24	2	0
5	A	317	0	0	3	0
5	B	286	0	0	1	0
All	All	5484	0	4675	61	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 6.

All (61) 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:144:MET:SD	1:B:247:MET:HG2	2.29	0.72
1:A:67:GLU:HA	1:A:70:LYS:HE2	1.72	0.70
1:B:232:LYS:HA	1:B:255:MET:HG3	1.76	0.68
1:B:144:MET:HG3	1:B:252:SER:O	1.97	0.64
1:B:110:PRO:HG3	1:B:178:TRP:CD2	2.35	0.61
1:B:144:MET:HG2	1:B:248:ASN:O	2.00	0.61
1:A:90:ASN:C	1:A:90:ASN:HD22	2.09	0.61
1:A:232:LYS:HA	1:A:255:MET:HG3	1.83	0.61
1:B:44:THR:O	1:B:107:HIS:HA	2.03	0.58
1:B:243:PRO:HD2	1:B:246:ILE:HD12	1.84	0.58
1:A:120:LYS:HE2	5:A:1076:HOH:O	2.03	0.57
1:A:238:ASN:C	1:A:238:ASN:HD22	2.11	0.57
1:A:44:THR:O	1:A:107:HIS:HA	2.06	0.56
1:B:238:ASN:C	1:B:238:ASN:HD22	2.12	0.56
1:B:123:LYS:NZ	1:B:211:GLU:HG2	2.21	0.56
1:B:123:LYS:HZ1	1:B:211:GLU:HG2	1.72	0.55
1:A:88:LEU:HD12	1:A:88:LEU:N	2.22	0.55
1:B:107:HIS:HB2	1:B:168:TYR:CE2	2.42	0.55
1:B:141:PHE:CE1	1:B:145:VAL:HG21	2.41	0.55
1:A:184:SER:OG	1:A:185:PRO:HD2	2.09	0.53
1:B:90:ASN:HD22	1:B:90:ASN:C	2.16	0.52
1:B:238:ASN:ND2	1:B:240:THR:H	2.07	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:ASN:O	1:B:94:GLN:HG3	2.11	0.51
1:A:235:PRO:HA	1:A:238:ASN:ND2	2.26	0.51
1:A:107:HIS:HB2	1:A:168:TYR:CE2	2.45	0.51
1:A:79:PRO:HD3	1:A:98:MET:HE1	1.93	0.51
1:A:107:HIS:HB2	1:A:168:TYR:CZ	2.46	0.50
1:B:235:PRO:HA	1:B:238:ASN:HD21	1.77	0.50
1:B:235:PRO:HA	1:B:238:ASN:ND2	2.27	0.49
1:A:246:ILE:HG23	3:A:712:PCQ:CL3	2.50	0.49
1:A:110:PRO:HG3	1:A:178:TRP:CD2	2.49	0.48
1:B:101:PRO:HB2	5:B:834:HOH:O	2.13	0.48
1:B:126:TYR:CE1	1:B:178:TRP:CD1	3.02	0.48
1:A:154:PHE:HB3	1:A:155:PRO:HD3	1.95	0.48
1:B:167:PRO:O	1:B:168:TYR:HB2	2.14	0.47
1:A:235:PRO:HA	1:A:238:ASN:HD21	1.80	0.47
1:A:238:ASN:ND2	1:A:240:THR:H	2.14	0.46
1:A:93:LYS:O	1:A:97:GLU:HG2	2.16	0.46
4:A:802:EDO:H22	5:A:891:HOH:O	2.16	0.45
1:B:14:VAL:HG13	1:B:14:VAL:O	2.15	0.45
1:A:268:THR:HA	1:B:268:THR:HA	1.98	0.45
1:B:231:MET:HB3	1:B:255:MET:HE3	1.98	0.45
1:A:80:PHE:CD2	1:A:83:CYS:HB2	2.52	0.45
1:B:107:HIS:HB2	1:B:168:TYR:CZ	2.52	0.44
1:A:238:ASN:C	1:A:238:ASN:ND2	2.75	0.44
1:A:243:PRO:HD2	1:A:246:ILE:HD12	1.98	0.44
1:A:233:ASN:ND2	4:A:804:EDO:H12	2.33	0.43
1:B:238:ASN:CG	1:B:240:THR:HG23	2.44	0.43
1:A:44:THR:HG22	1:A:127:LEU:HB3	2.01	0.42
1:B:44:THR:HG22	1:B:127:LEU:HB3	2.01	0.42
1:A:167:PRO:O	1:A:168:TYR:HB2	2.20	0.42
1:A:88:LEU:N	1:A:88:LEU:CD1	2.84	0.41
1:A:141:PHE:CE1	1:A:145:VAL:HG21	2.56	0.41
1:B:290:PHE:HB3	1:B:291:ARG:H	1.46	0.41
1:A:90:ASN:ND2	1:A:93:LYS:H	2.18	0.41
1:B:179:TRP:CE3	1:B:288:LEU:HD13	2.56	0.41
1:A:49:GLY:HA2	5:A:868:HOH:O	2.21	0.41
1:B:73:VAL:HG21	1:B:75:PHE:CZ	2.56	0.41
1:B:290:PHE:O	1:B:291:ARG:HG3	2.21	0.41
1:B:47:LYS:HG3	1:B:138:PHE:HA	2.04	0.40
1:A:245:GLU:CD	1:A:245:GLU:H	2.28	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	268/294 (91%)	257 (96%)	11 (4%)	0	100	100
1	B	281/294 (96%)	273 (97%)	8 (3%)	0	100	100
All	All	549/588 (93%)	530 (96%)	19 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	265/275 (96%)	262 (99%)	3 (1%)	65	54
1	B	253/275 (92%)	251 (99%)	2 (1%)	73	65
All	All	518/550 (94%)	513 (99%)	5 (1%)	68	58

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	138	PHE
1	A	238	ASN
1	A	280	GLU
1	B	138	PHE
1	B	238	ASN

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	90	ASN
1	A	94	GLN
1	A	99	ASN
1	A	229	GLN
1	A	233	ASN
1	A	238	ASN
1	A	281	GLN
1	B	90	ASN
1	B	99	ASN
1	B	165	GLN
1	B	208	HIS
1	B	224	HIS
1	B	233	ASN
1	B	238	ASN
1	B	265	ASN

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](#)

8 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	PCQ	A	712	-	19,19,19	3.08	10 (52%)	28,28,28	0.84	0
4	EDO	A	804	-	3,3,3	2.28	2 (66%)	2,2,2	0.47	0
4	EDO	A	803	-	3,3,3	2.28	2 (66%)	2,2,2	0.42	0
4	EDO	A	801	-	3,3,3	2.26	2 (66%)	2,2,2	0.46	0
4	EDO	A	802	-	3,3,3	2.24	2 (66%)	2,2,2	0.40	0
2	A3P	B	702	-	29,29,29	0.95	0	44,45,45	0.70	1 (2%)
3	PCQ	B	711	-	19,19,19	3.12	12 (63%)	28,28,28	0.87	0
2	A3P	A	701	-	29,29,29	0.96	0	44,45,45	0.67	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	PCQ	A	712	-	-	0/4/4/4	0/2/2/2
4	EDO	A	804	-	-	1/1/1/1	-
4	EDO	A	803	-	-	0/1/1/1	-
4	EDO	A	801	-	-	1/1/1/1	-
4	EDO	A	802	-	-	1/1/1/1	-
2	A3P	B	702	-	-	0/15/31/31	0/3/3/3
3	PCQ	B	711	-	-	0/4/4/4	0/2/2/2
2	A3P	A	701	-	-	0/15/31/31	0/3/3/3

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	712	PCQ	C2-C1	8.37	1.54	1.39
3	B	711	PCQ	C2-C1	8.31	1.54	1.39
3	A	712	PCQ	C4-C5	4.39	1.46	1.39
3	B	711	PCQ	C6'-C1'	4.36	1.47	1.39
3	A	712	PCQ	C6'-C1'	4.21	1.47	1.39
3	B	711	PCQ	C4-C5	4.16	1.45	1.39
3	B	711	PCQ	C4-C3	4.05	1.45	1.39
3	A	712	PCQ	C4-C3	3.72	1.45	1.39
3	B	711	PCQ	C4'-C5'	3.52	1.45	1.39
3	A	712	PCQ	C4'-C5'	3.35	1.44	1.39
3	B	711	PCQ	C6-C1	3.13	1.45	1.39
3	A	712	PCQ	C2'-C1'	3.06	1.45	1.39
4	A	803	EDO	O2-C2	3.05	1.57	1.42
4	A	804	EDO	O2-C2	3.00	1.57	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	802	EDO	O2-C2	3.00	1.57	1.42
4	A	801	EDO	O2-C2	3.00	1.57	1.42
3	A	712	PCQ	C6'-C5'	2.98	1.43	1.38
3	B	711	PCQ	C2'-C1'	2.94	1.44	1.39
3	B	711	PCQ	C6'-C5'	2.85	1.43	1.38
3	A	712	PCQ	C6-C1	2.81	1.44	1.39
3	B	711	PCQ	C4'-C3'	2.78	1.43	1.39
3	A	712	PCQ	C4'-C3'	2.66	1.43	1.39
4	A	804	EDO	O1-C1	2.43	1.54	1.42
4	A	803	EDO	O1-C1	2.38	1.54	1.42
4	A	801	EDO	O1-C1	2.38	1.54	1.42
4	A	802	EDO	O1-C1	2.36	1.54	1.42
3	B	711	PCQ	C6-C5	2.34	1.42	1.38
3	A	712	PCQ	C2'-C3'	2.28	1.42	1.38
3	B	711	PCQ	C2'-C3'	2.07	1.42	1.38
3	B	711	PCQ	C2-C3	2.03	1.41	1.38

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	702	A3P	O2'-C2'-C3'	2.25	117.48	111.19

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	802	EDO	O1-C1-C2-O2
4	A	801	EDO	O1-C1-C2-O2
4	A	804	EDO	O1-C1-C2-O2

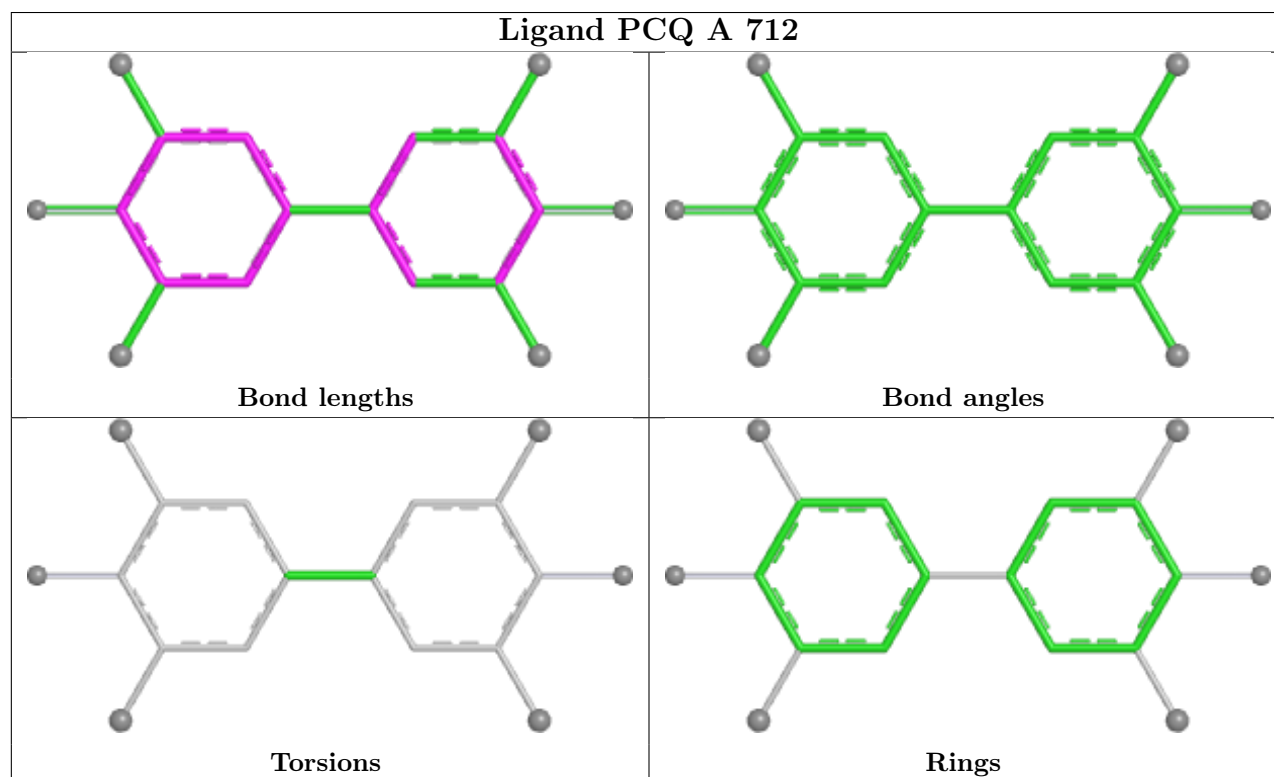
There are no ring outliers.

3 monomers are involved in 3 short contacts:

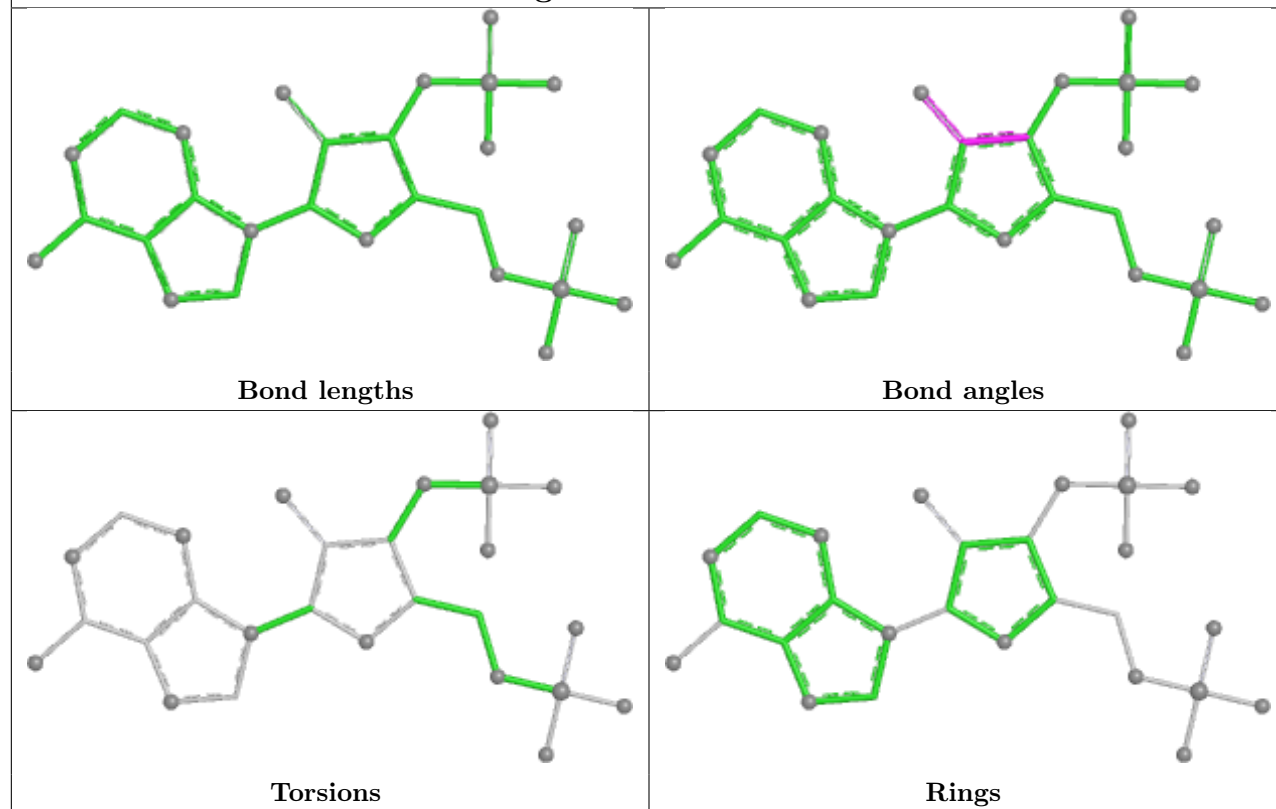
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	712	PCQ	1	0
4	A	804	EDO	1	0
4	A	802	EDO	1	0

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

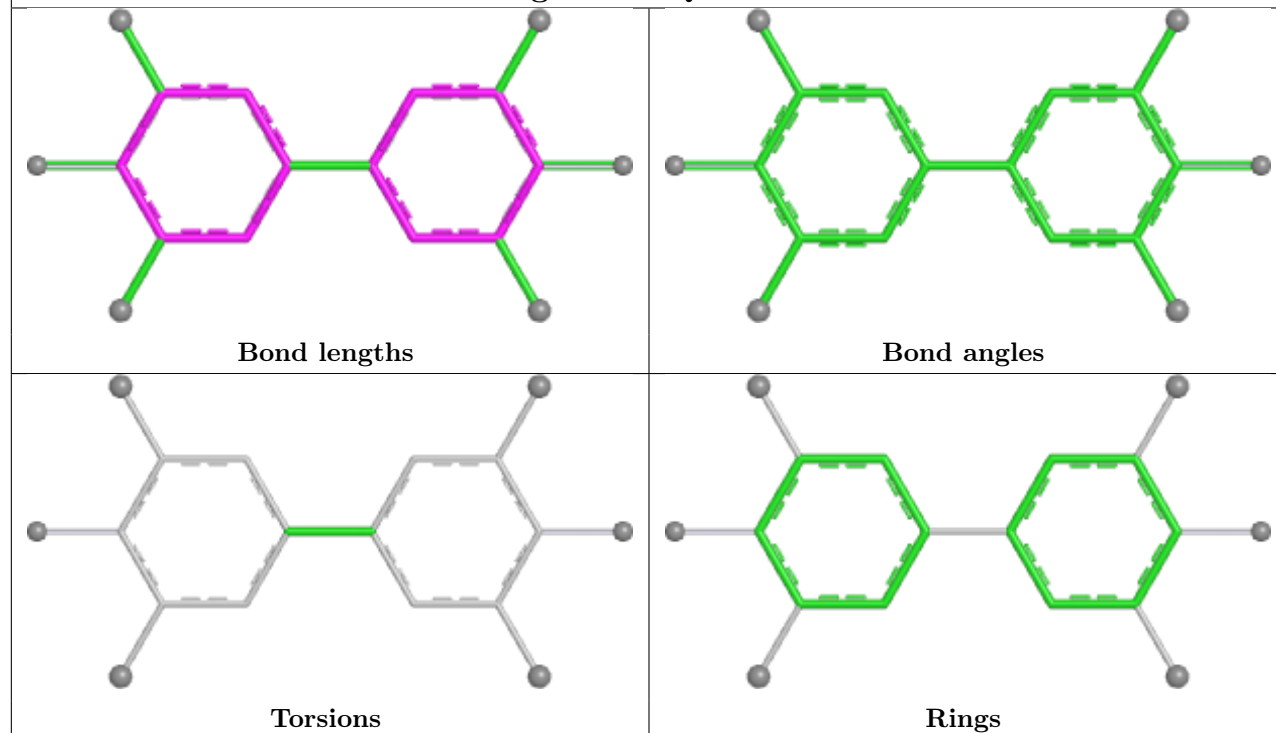
also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted 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 A3P B 702



Ligand PCQ B 711



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.