Full wwPDB X-ray Structure Validation Report (i)

Jun 13, 2024 – 12:42 AM EDT

PDB ID	:	3CAS
Title	:	Crystal structure of 5beta-reductase (AKR1D1) in complex with NADP+ and 4-androstenedione
Authors	:	Faucher, F.; Cantin, L.; Breton, R.
Deposited on	:	2008-02-20
Resolution	:	2.00 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 (i) 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 (i)) were used in the production of this report:

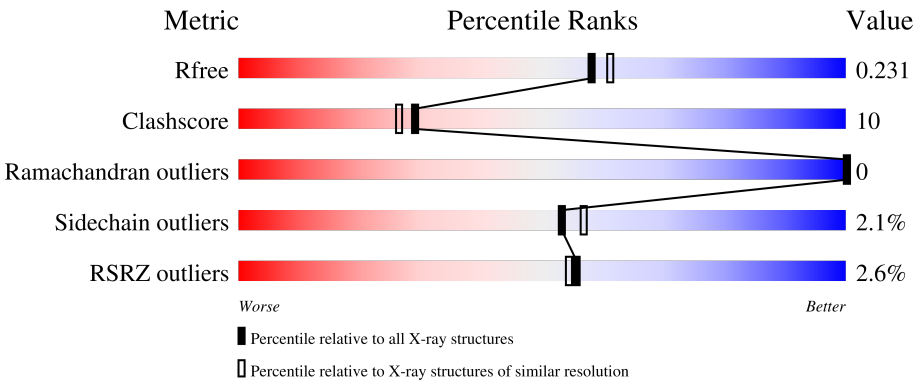
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 <=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	326	86% 13% •
1	B	326	5% 79% 20% •

2 Entry composition (i)

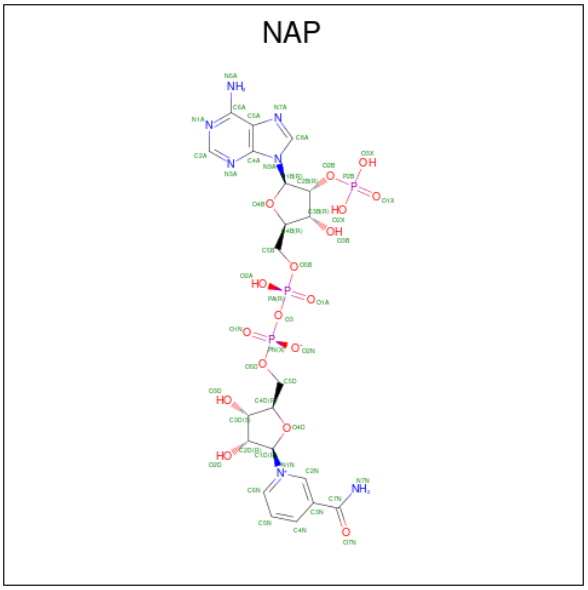
There are 6 unique types of molecules in this entry. The entry contains 5923 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 3-oxo-5-beta-steroid 4-dehydrogenase.

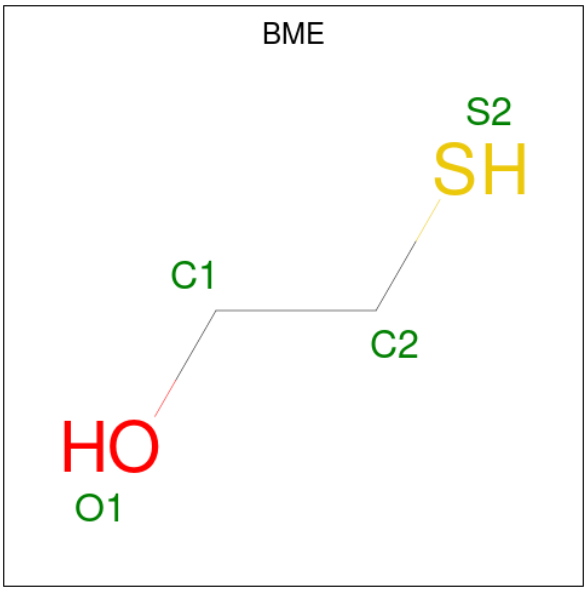
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	325	2642	1689	458	484	11	0	2	0
1	B	325	2654	1696	462	485	11	0	3	0

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



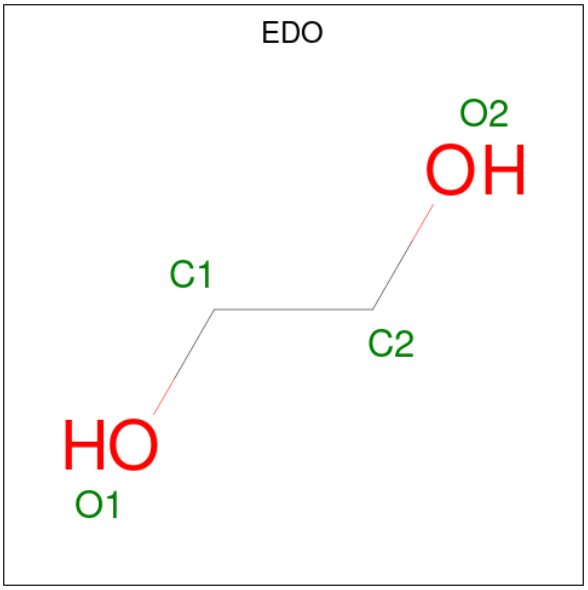
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	48	21	7	17	3	0	0
2	B	1	48	21	7	17	3	0	0

- Molecule 3 is BETA-MERCAPTOETHANOL (three-letter code: BME) (formula: C_2H_6OS).



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

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

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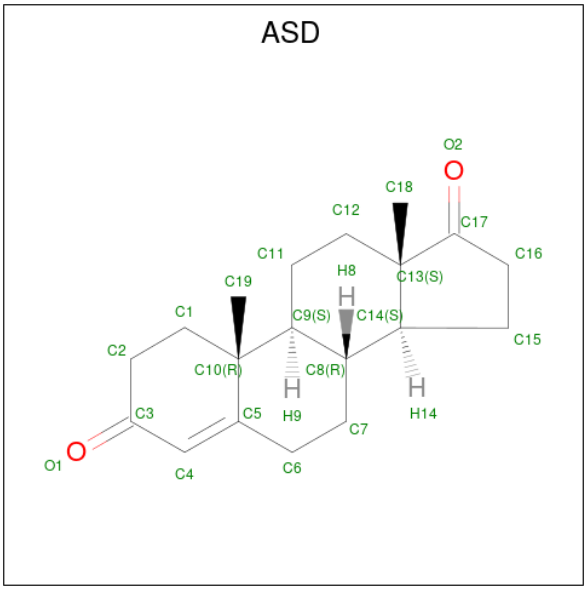
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0

- Molecule 5 is 4-ANDROSTENE-3-17-DIONE (three-letter code: ASD) (formula: $C_{19}H_{26}O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 21 19 2	0	0
5	B	1	Total C O 21 19 2	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	238	Total O 238 238	0	0
6	B	223	Total O 223 223	0	0

3 Residue-property plots (i)

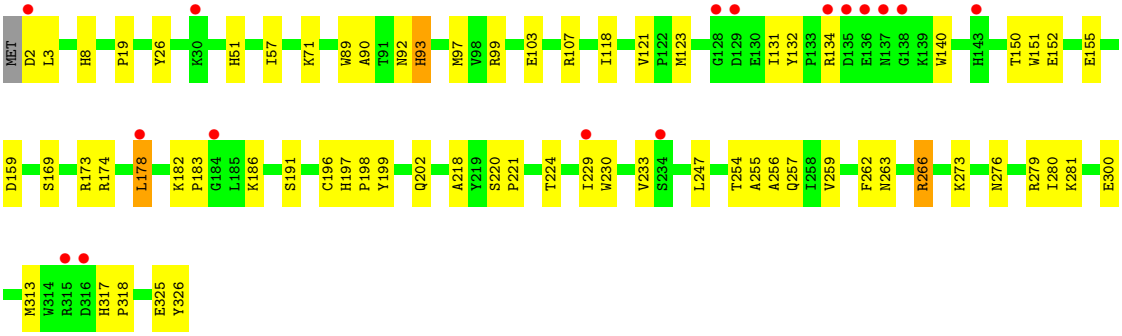
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: 3-oxo-5-beta-steroid 4-dehydrogenase

Chain A:

• Molecule 1: 3-oxo-5-beta-steroid 4-dehydrogenase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	50.41Å 110.87Å 130.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.63 – 2.00 19.63 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.8 (19.63-2.00) 96.0 (19.63-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.12	Depositor
$< I/\sigma(I) >$ [1]	4.12 (at 2.01Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.208 , 0.238 0.201 , 0.231	Depositor DCC
R_{free} test set	2435 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	12.7	Xtriage
Anisotropy	0.406	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 55.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.44$, $< L^2 > = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	5923	wwPDB-VP
Average B, all atoms (Å^2)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.00% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: NAP, BME, EDO, ASD

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	0/2707	0.58	0/3670
1	B	0.34	0/2719	0.56	0/3684
All	All	0.35	0/5426	0.57	0/7354

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	219	TYR	Sidechain

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	2642	0	2635	50	0
1	B	2654	0	2652	56	0
2	A	48	0	25	5	0
2	B	48	0	25	3	0
3	A	4	0	6	1	0
4	A	12	0	15	4	0
4	B	12	0	15	2	0
5	A	21	0	26	3	0
5	B	21	0	26	3	0
6	A	238	0	0	2	0
6	B	223	0	0	5	0
All	All	5923	0	5425	105	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 10.

All (105) 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:148:CYS:SG	3:A:328:BME:S2	2.36	1.14
1:A:12:LEU:HB3	1:A:190[B]:VAL:HG21	1.55	0.87
1:A:276:ASN:HD22	1:A:279:ARG:H	1.25	0.82
1:B:276:ASN:HD22	1:B:279:ARG:H	1.23	0.82
1:B:93:HIS:O	1:B:150[A]:THR:HG22	1.82	0.78
1:A:12:LEU:HB3	1:A:190[A]:VAL:HG11	1.68	0.74
1:A:254:THR:H	1:A:257:GLN:HE21	1.35	0.73
1:A:12:LEU:HD22	1:A:190[A]:VAL:HG12	1.71	0.72
1:B:254:THR:H	1:B:257:GLN:HE21	1.37	0.72
1:A:197:HIS:HD2	1:A:199:TYR:H	1.38	0.71
1:A:254:THR:H	1:A:257:GLN:NE2	1.89	0.70
1:A:152:GLU:HG3	1:A:182:LYS:HE3	1.74	0.69
1:A:12:LEU:HB3	1:A:190[B]:VAL:CG2	2.25	0.65
1:A:8:HIS:H	4:A:329:EDO:H21	1.60	0.65
1:B:191:SER:HB2	6:B:656:HOH:O	1.97	0.63
1:B:8:HIS:H	4:B:330:EDO:H12	1.63	0.63
1:A:298:ASP:OD1	1:B:107[A]:ARG:HD3	1.99	0.62
1:A:147:LEU:HB2	4:A:330:EDO:H22	1.82	0.61
1:B:92:ASN:HA	1:B:97:MET:HE3	1.83	0.60
1:B:197:HIS:HD2	1:B:199:TYR:H	1.50	0.59
1:B:159:ASP:OD2	1:B:186[B]:LYS:HE2	2.03	0.59
1:A:12:LEU:HD22	1:A:190[A]:VAL:CG1	2.33	0.59
1:B:134:ARG:HG2	1:B:140:TRP:HA	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:273:LYS:HD3	1:B:273:LYS:C	2.24	0.58
1:A:191:SER:HB2	6:A:431:HOH:O	2.03	0.58
1:B:276:ASN:ND2	1:B:279:ARG:H	1.99	0.56
1:B:262:PHE:O	1:B:266:ARG:HG2	2.04	0.55
1:B:2:ASP:HA	6:B:665:HOH:O	2.05	0.55
1:B:281:LYS:HG3	6:B:673:HOH:O	2.07	0.55
1:B:256:ALA:O	1:B:259:VAL:HG22	2.06	0.55
1:B:276:ASN:O	1:B:280:ILE:HG13	2.07	0.55
1:A:166:LEU:O	1:A:190[B]:VAL:HG12	2.08	0.54
1:A:195:GLU:OE2	1:A:197:HIS:HE1	1.90	0.54
2:B:327:NAP:H5N	5:B:331:ASD:H62	1.90	0.54
1:A:12:LEU:HB3	1:A:190[A]:VAL:CG1	2.32	0.54
1:B:152:GLU:HG3	1:B:182:LYS:HE3	1.90	0.53
1:B:92:ASN:HA	1:B:97:MET:CE	2.39	0.53
1:A:26:TYR:CE2	5:A:332:ASD:H12	2.44	0.53
1:A:273:LYS:C	1:A:273:LYS:HD3	2.29	0.53
1:B:8:HIS:H	4:B:330:EDO:C1	2.21	0.52
1:B:132:TYR:HB3	1:B:140:TRP:CZ3	2.44	0.52
1:A:99:ARG:HB3	1:A:100:PRO:HD3	1.93	0.51
1:A:12:LEU:HD22	1:A:190[B]:VAL:HG22	1.92	0.51
1:B:99:ARG:O	1:B:103:GLU:HG3	2.10	0.51
1:B:118:ILE:HD12	1:B:169:SER:HB2	1.92	0.51
1:A:187:HIS:HD2	6:A:457:HOH:O	1.95	0.50
1:B:254:THR:H	1:B:257:GLN:NE2	2.06	0.50
1:A:263:ASN:HA	1:A:266:ARG:HD2	1.94	0.50
1:A:118:ILE:HD12	1:A:169:SER:HB2	1.94	0.49
1:A:134:ARG:HG2	1:A:140:TRP:HA	1.94	0.49
1:A:26:TYR:CD2	5:A:332:ASD:H193	2.48	0.49
1:A:224:THR:C	2:A:327:NAP:H51A	2.32	0.49
2:A:327:NAP:H5N	5:A:332:ASD:H62	1.95	0.49
1:B:132:TYR:HB3	1:B:140:TRP:HZ3	1.76	0.48
1:A:16[A]:ASN:HD22	1:A:17:SER:N	2.12	0.48
1:A:16[A]:ASN:ND2	1:A:17:SER:N	2.60	0.48
1:A:262:PHE:O	1:A:266:ARG:HG2	2.12	0.48
1:B:152:GLU:CG	1:B:182:LYS:HE3	2.43	0.47
1:A:263:ASN:O	1:A:266:ARG:HG3	2.14	0.47
1:B:134:ARG:HD3	1:B:140:TRP:CE2	2.49	0.47
1:B:174:ARG:O	1:B:178:LEU:HD12	2.15	0.46
1:B:230:TRP:CD1	5:B:331:ASD:H11	2.50	0.46
1:A:254:THR:N	1:A:257:GLN:HE21	2.09	0.46
1:A:279:ARG:HG2	2:A:327:NAP:C5A	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:19:PRO:HB2	1:B:51:HIS:HB2	1.97	0.46
1:B:263:ASN:HA	1:B:266:ARG:HD2	1.98	0.45
1:B:57:ILE:HG22	1:B:131:ILE:HD12	2.00	0.44
1:B:2:ASP:OD1	1:B:3:LEU:N	2.50	0.44
1:A:8:HIS:H	4:A:329:EDO:C2	2.27	0.44
1:B:121:VAL:HB	1:B:123:MET:CE	2.48	0.44
1:B:220:SER:N	1:B:221:PRO:CD	2.81	0.44
1:B:173:ARG:HD2	1:B:326:TYR:CE2	2.53	0.43
1:B:263:ASN:O	1:B:266:ARG:HG3	2.18	0.43
1:A:273:LYS:O	2:A:327:NAP:H8A	2.18	0.43
1:B:197:HIS:HB2	1:B:198:PRO:CD	2.48	0.43
1:B:151:TRP:O	1:B:155:GLU:HG3	2.18	0.43
1:A:298:ASP:HA	1:B:107[B]:ARG:HD3	2.00	0.43
1:B:26:TYR:CD2	5:B:331:ASD:H193	2.54	0.43
1:A:262:PHE:O	1:A:266:ARG:CG	2.67	0.43
1:B:233:VAL:O	1:B:233:VAL:HG12	2.19	0.43
1:B:197:HIS:HD2	1:B:199:TYR:N	2.17	0.42
1:A:276:ASN:HD21	1:A:278:GLU:HB2	1.84	0.42
1:A:279:ARG:HG2	2:A:327:NAP:C6A	2.49	0.42
1:B:273:LYS:O	2:B:327:NAP:H8A	2.18	0.42
1:A:297:LYS:HD2	1:B:107[A]:ARG:NH2	2.34	0.42
1:B:196:CYS:HB3	1:B:218:ALA:CB	2.50	0.42
1:B:255:ALA:O	1:B:259:VAL:HG13	2.20	0.42
1:A:204:LYS:HE3	1:A:325:GLU:OE2	2.20	0.42
1:B:89:TRP:CG	1:B:90:ALA:N	2.88	0.41
1:A:263:ASN:HA	1:A:266:ARG:CD	2.50	0.41
1:B:93:HIS:HD2	6:B:674:HOH:O	2.03	0.41
1:B:71:LYS:HD3	1:B:71:LYS:HA	1.91	0.41
1:A:9:ARG:HG3	4:A:329:EDO:O2	2.20	0.41
1:A:220:SER:N	1:A:221:PRO:CD	2.84	0.41
1:A:256:ALA:O	1:A:259:VAL:HG22	2.20	0.41
1:A:152:GLU:CG	1:A:182:LYS:HE3	2.49	0.41
1:B:182:LYS:HA	1:B:183:PRO:HD3	1.94	0.41
1:B:317:HIS:HA	1:B:318:PRO:HD3	1.98	0.41
1:B:90:ALA:HA	1:B:93:HIS:CE1	2.56	0.40
1:A:276:ASN:ND2	1:A:279:ARG:H	2.06	0.40
1:B:229:ILE:HG12	6:B:531:HOH:O	2.21	0.40
1:B:224:THR:C	2:B:327:NAP:H51A	2.41	0.40
1:B:325:GLU:O	1:B:326:TYR:HB3	2.21	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	325/326 (100%)	317 (98%)	8 (2%)	0	100	100
1	B	326/326 (100%)	317 (97%)	9 (3%)	0	100	100
All	All	651/652 (100%)	634 (97%)	17 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	291/290 (100%)	285 (98%)	6 (2%)	53	57
1	B	292/290 (101%)	285 (98%)	7 (2%)	49	51
All	All	583/580 (100%)	570 (98%)	13 (2%)	53	55

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16[A]	ASN
1	A	16[B]	ASN
1	A	202	GLN
1	A	266	ARG
1	A	300	GLU
1	A	316	ASP
1	B	93	HIS
1	B	178	LEU

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Mol	Chain	Res	Type
1	B	202	GLN
1	B	247	LEU
1	B	266	ARG
1	B	300	GLU
1	B	313	MET

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

Mol	Chain	Res	Type
1	A	51	HIS
1	A	59	GLN
1	A	92	ASN
1	A	110	GLN
1	A	187	HIS
1	A	197	HIS
1	A	202	GLN
1	A	211	GLN
1	A	257	GLN
1	A	276	ASN
1	B	59	GLN
1	B	92	ASN
1	B	170	ASN
1	B	187	HIS
1	B	197	HIS
1	B	202	GLN
1	B	210	GLN
1	B	257	GLN
1	B	276	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

11 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	BME	A	328	-	3,3,3	0.28	0	2,2,2	0.34	0
4	EDO	A	329	-	3,3,3	2.39	1 (33%)	2,2,2	0.35	0
2	NAP	B	327	-	46,52,52	1.76	8 (17%)	61,80,80	1.03	4 (6%)
4	EDO	B	329	-	3,3,3	2.29	1 (33%)	2,2,2	0.37	0
4	EDO	A	331	-	3,3,3	2.29	1 (33%)	2,2,2	0.38	0
5	ASD	A	332	-	24,24,24	3.44	10 (41%)	39,39,39	2.00	12 (30%)
2	NAP	A	327	-	46,52,52	1.66	8 (17%)	61,80,80	1.04	3 (4%)
4	EDO	B	330	-	3,3,3	2.38	1 (33%)	2,2,2	0.39	0
5	ASD	B	331	-	24,24,24	3.45	11 (45%)	39,39,39	2.00	14 (35%)
4	EDO	A	330	-	3,3,3	2.29	1 (33%)	2,2,2	0.34	0
4	EDO	B	328	-	3,3,3	2.32	1 (33%)	2,2,2	0.40	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	BME	A	328	-	-	0/1/1/1	-
4	EDO	A	329	-	-	1/1/1/1	-
2	NAP	B	327	-	-	9/31/67/67	0/5/5/5
4	EDO	B	329	-	-	0/1/1/1	-
4	EDO	A	331	-	-	1/1/1/1	-
5	ASD	A	332	-	-	-	0/4/4/4
2	NAP	A	327	-	-	6/31/67/67	0/5/5/5
4	EDO	B	330	-	-	0/1/1/1	-
5	ASD	B	331	-	-	-	0/4/4/4
4	EDO	A	330	-	-	1/1/1/1	-
4	EDO	B	328	-	-	1/1/1/1	-

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	331	ASD	C4-C5	14.28	1.54	1.34
5	A	332	ASD	C4-C5	14.24	1.54	1.34
2	B	327	NAP	C2N-N1N	4.95	1.40	1.35
2	B	327	NAP	C4N-C3N	4.61	1.46	1.39
2	B	327	NAP	C6N-N1N	4.37	1.45	1.35
2	A	327	NAP	C6N-N1N	4.16	1.44	1.35
2	A	327	NAP	C2N-N1N	4.16	1.39	1.35
4	A	329	EDO	O2-C2	-4.08	1.21	1.42
4	B	330	EDO	O2-C2	-4.06	1.21	1.42
4	B	328	EDO	O2-C2	-3.96	1.21	1.42
2	B	327	NAP	C2A-N3A	3.93	1.38	1.32
2	A	327	NAP	C4N-C3N	3.91	1.45	1.39
4	B	329	EDO	O2-C2	-3.90	1.22	1.42
4	A	331	EDO	O2-C2	-3.90	1.22	1.42
4	A	330	EDO	O2-C2	-3.88	1.22	1.42
2	A	327	NAP	C2A-N3A	3.75	1.37	1.32
5	A	332	ASD	C16-C17	3.59	1.57	1.51
2	A	327	NAP	C4A-N3A	3.55	1.40	1.35
2	B	327	NAP	C4A-N3A	3.42	1.40	1.35
5	B	331	ASD	C16-C17	3.38	1.57	1.51
5	A	332	ASD	C8-C9	3.14	1.59	1.53
2	A	327	NAP	P2B-O2B	2.98	1.64	1.59
5	A	332	ASD	C2-C3	2.90	1.56	1.50
5	B	331	ASD	C7-C8	2.89	1.58	1.53
5	B	331	ASD	C2-C3	2.89	1.56	1.50
5	B	331	ASD	C8-C9	2.80	1.58	1.53
2	B	327	NAP	C5N-C4N	2.71	1.43	1.38
2	A	327	NAP	C5N-C4N	2.59	1.43	1.38
2	B	327	NAP	C2A-N1A	2.53	1.38	1.33
5	B	331	ASD	C7-C6	2.52	1.58	1.53
5	A	332	ASD	C7-C8	2.45	1.57	1.53
5	A	332	ASD	C7-C6	2.44	1.58	1.53
5	A	332	ASD	C1-C10	2.33	1.58	1.54
5	B	331	ASD	C1-C10	2.30	1.58	1.54
2	A	327	NAP	C2A-N1A	2.29	1.38	1.33
5	A	332	ASD	C11-C9	2.28	1.57	1.53
5	B	331	ASD	C11-C9	2.21	1.57	1.53
5	B	331	ASD	C6-C5	2.18	1.54	1.50
5	B	331	ASD	C12-C11	2.17	1.57	1.53
5	A	332	ASD	C6-C5	2.13	1.54	1.50
5	B	331	ASD	C4-C3	-2.10	1.41	1.45
2	B	327	NAP	P2B-O2B	2.08	1.63	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	332	ASD	C12-C11	2.03	1.57	1.53

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	A	332	ASD	C19-C10-C9	4.99	117.26	111.66
5	B	331	ASD	C19-C10-C9	4.72	116.96	111.66
5	B	331	ASD	C14-C13-C17	-4.53	95.29	100.61
5	A	332	ASD	C14-C13-C17	-4.43	95.40	100.61
5	A	332	ASD	C16-C15-C14	-3.94	97.03	103.00
5	B	331	ASD	C16-C15-C14	-3.89	97.10	103.00
5	A	332	ASD	C2-C3-C4	3.44	121.96	116.73
5	B	331	ASD	C18-C13-C14	3.33	117.78	112.99
5	B	331	ASD	C2-C3-C4	3.31	121.75	116.73
5	B	331	ASD	C15-C14-C8	-3.20	114.00	119.10
5	A	332	ASD	C18-C13-C14	3.14	117.50	112.99
5	A	332	ASD	C15-C14-C8	-3.06	114.22	119.10
5	A	332	ASD	C18-C13-C17	3.01	110.73	105.18
5	B	331	ASD	C18-C13-C17	3.01	110.72	105.18
5	A	332	ASD	C2-C1-C10	2.89	118.45	113.42
2	A	327	NAP	N3A-C2A-N1A	-2.57	125.18	128.67
2	B	327	NAP	N3A-C2A-N1A	-2.42	125.38	128.67
5	B	331	ASD	C2-C1-C10	2.42	117.63	113.42
5	B	331	ASD	C18-C13-C12	-2.33	108.44	111.10
5	A	332	ASD	C5-C4-C3	-2.32	120.08	123.67
5	B	331	ASD	C5-C4-C3	-2.30	120.11	123.67
5	B	331	ASD	C6-C5-C4	-2.29	117.20	120.86
2	A	327	NAP	C6N-N1N-C1D	2.23	124.10	119.73
5	B	331	ASD	C12-C13-C14	-2.20	105.47	108.88
5	B	331	ASD	C19-C10-C5	-2.16	105.08	108.38
5	A	332	ASD	C18-C13-C12	-2.16	108.64	111.10
2	B	327	NAP	C6N-N1N-C1D	2.15	123.96	119.73
5	A	332	ASD	C19-C10-C5	-2.14	105.12	108.38
5	B	331	ASD	C13-C14-C8	2.12	115.36	113.13
2	B	327	NAP	C4A-C5A-N7A	2.06	111.52	109.34
5	A	332	ASD	C6-C5-C4	-2.06	117.57	120.86
2	B	327	NAP	O3D-C3D-C4D	2.05	116.96	111.08
2	A	327	NAP	C4A-C5A-N7A	2.02	111.47	109.34

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	327	NAP	C5B-O5B-PA-O1A
2	B	327	NAP	C5B-O5B-PA-O1A
2	B	327	NAP	C5B-O5B-PA-O2A
2	B	327	NAP	C5B-O5B-PA-O3
2	B	327	NAP	O4B-C4B-C5B-O5B
2	B	327	NAP	C3B-C4B-C5B-O5B
4	A	329	EDO	O1-C1-C2-O2
4	A	330	EDO	O1-C1-C2-O2
2	A	327	NAP	PA-O3-PN-O5D
2	B	327	NAP	PA-O3-PN-O5D
2	A	327	NAP	C4D-C5D-O5D-PN
2	A	327	NAP	O4B-C4B-C5B-O5B
2	A	327	NAP	C5B-O5B-PA-O3
2	B	327	NAP	C4D-C5D-O5D-PN
2	A	327	NAP	PN-O3-PA-O2A
4	A	331	EDO	O1-C1-C2-O2
4	B	328	EDO	O1-C1-C2-O2
2	B	327	NAP	PN-O3-PA-O1A
2	B	327	NAP	PN-O3-PA-O2A

There are no ring outliers.

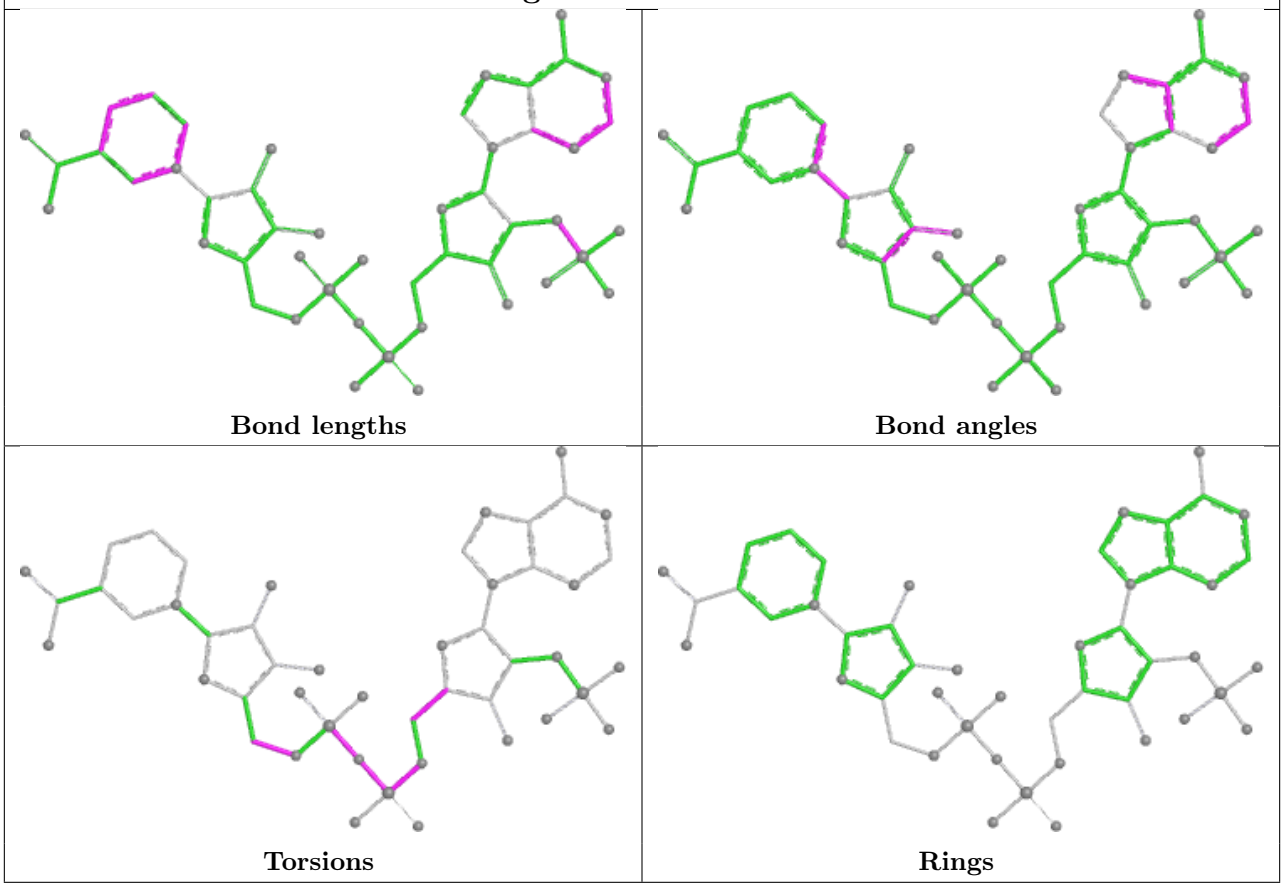
8 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	328	BME	1	0
4	A	329	EDO	3	0
2	B	327	NAP	3	0
5	A	332	ASD	3	0
2	A	327	NAP	5	0
4	B	330	EDO	2	0
5	B	331	ASD	3	0
4	A	330	EDO	1	0

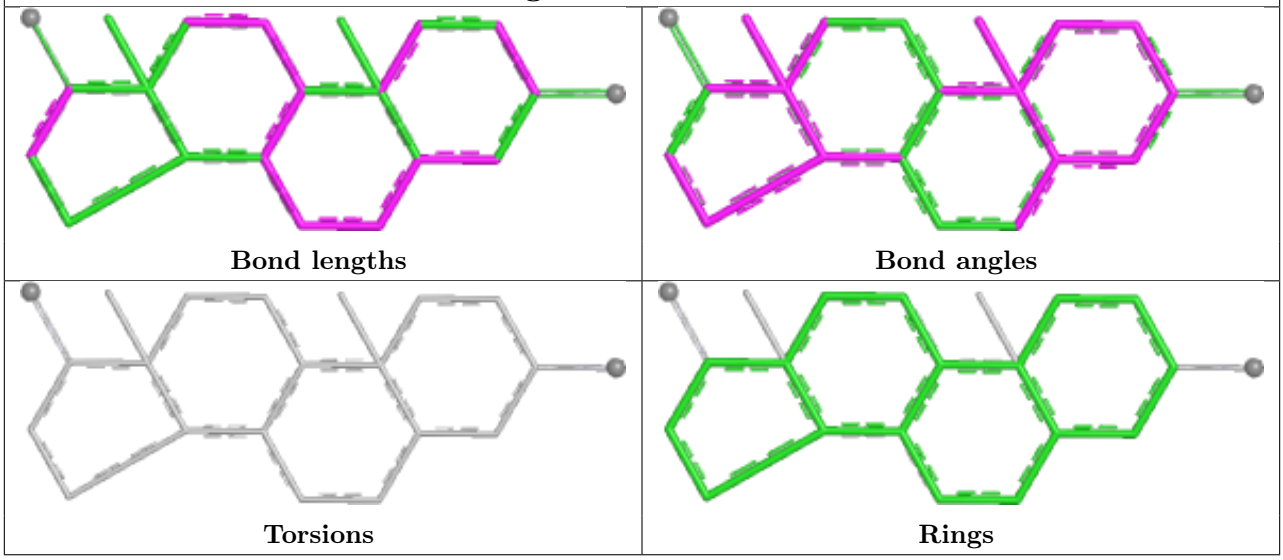
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

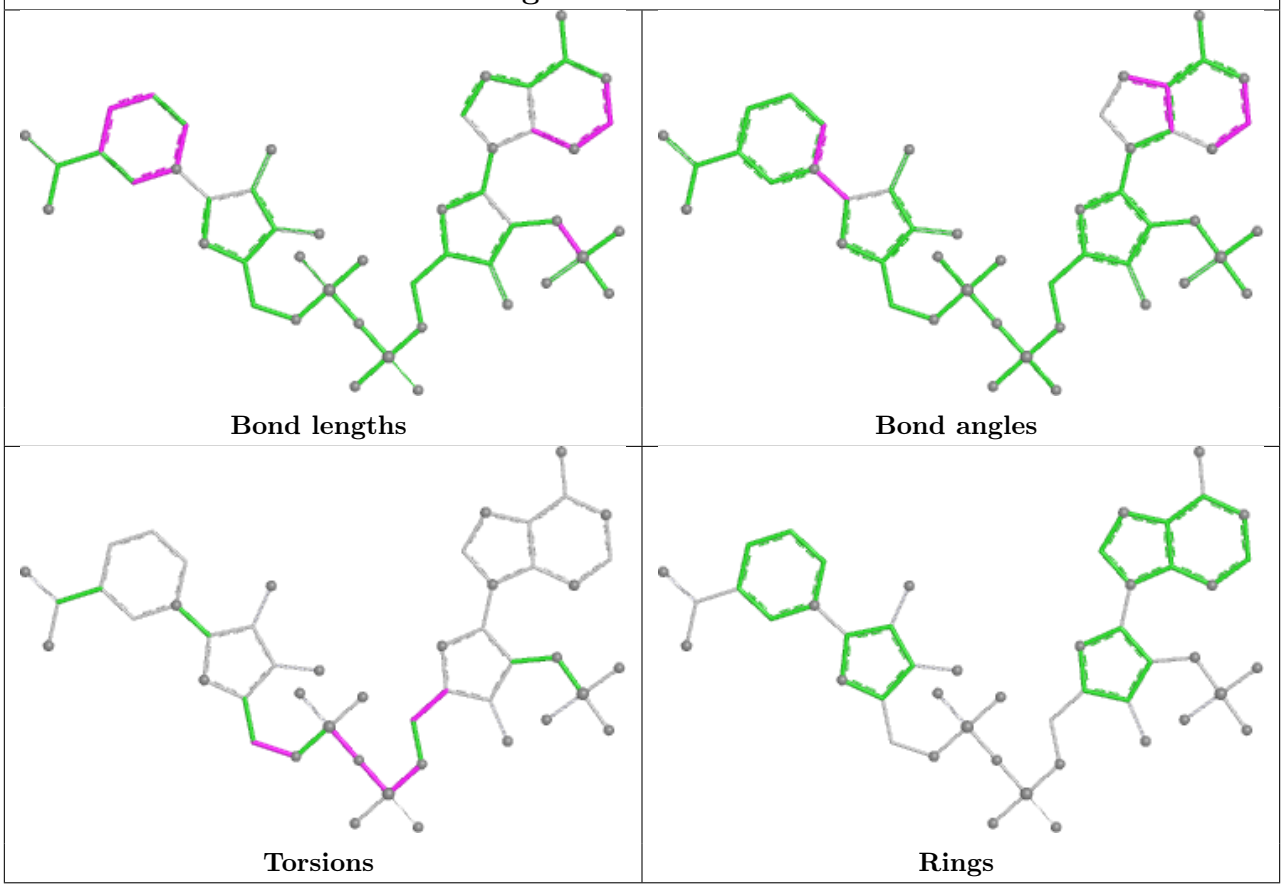
Ligand NAP B 327



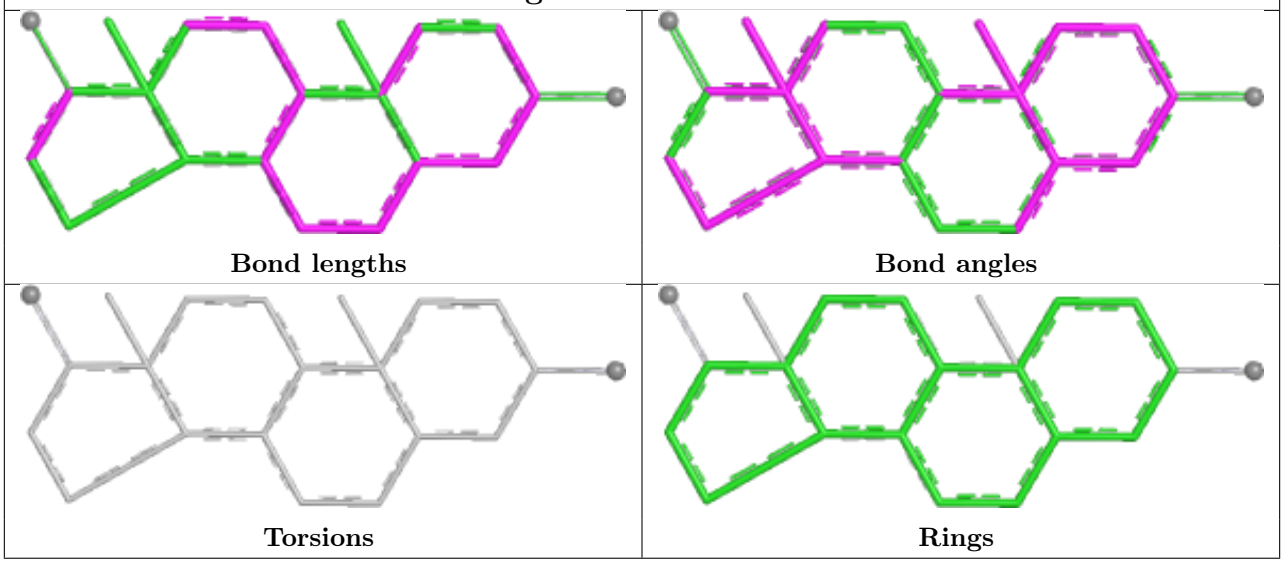
Ligand ASD A 332



Ligand NAP A 327



Ligand ASD B 331



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	325/326 (99%)	-0.15	1 (0%)	94	93	3, 11, 24, 31	5 (1%)
1	B	325/326 (99%)	0.15	16 (4%)	29	28	4, 13, 27, 34	3 (0%)
All	All	650/652 (99%)	0.00	17 (2%)	56	54	3, 12, 26, 34	8 (1%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	137	ASN	4.6
1	B	2	ASP	4.2
1	B	138	GLY	3.7
1	B	30	LYS	3.7
1	B	129	ASP	3.3
1	A	232	ASN	3.0
1	B	135	ASP	3.0
1	B	136	GLU	2.8
1	B	128	GLY	2.6
1	B	184	GLY	2.5
1	B	143	HIS	2.5
1	B	316	ASP	2.5
1	B	134	ARG	2.4
1	B	178	LEU	2.4
1	B	229	ILE	2.3
1	B	234	SER	2.2
1	B	315	ARG	2.2

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 monosaccharides in this entry.

6.4 Ligands (i)

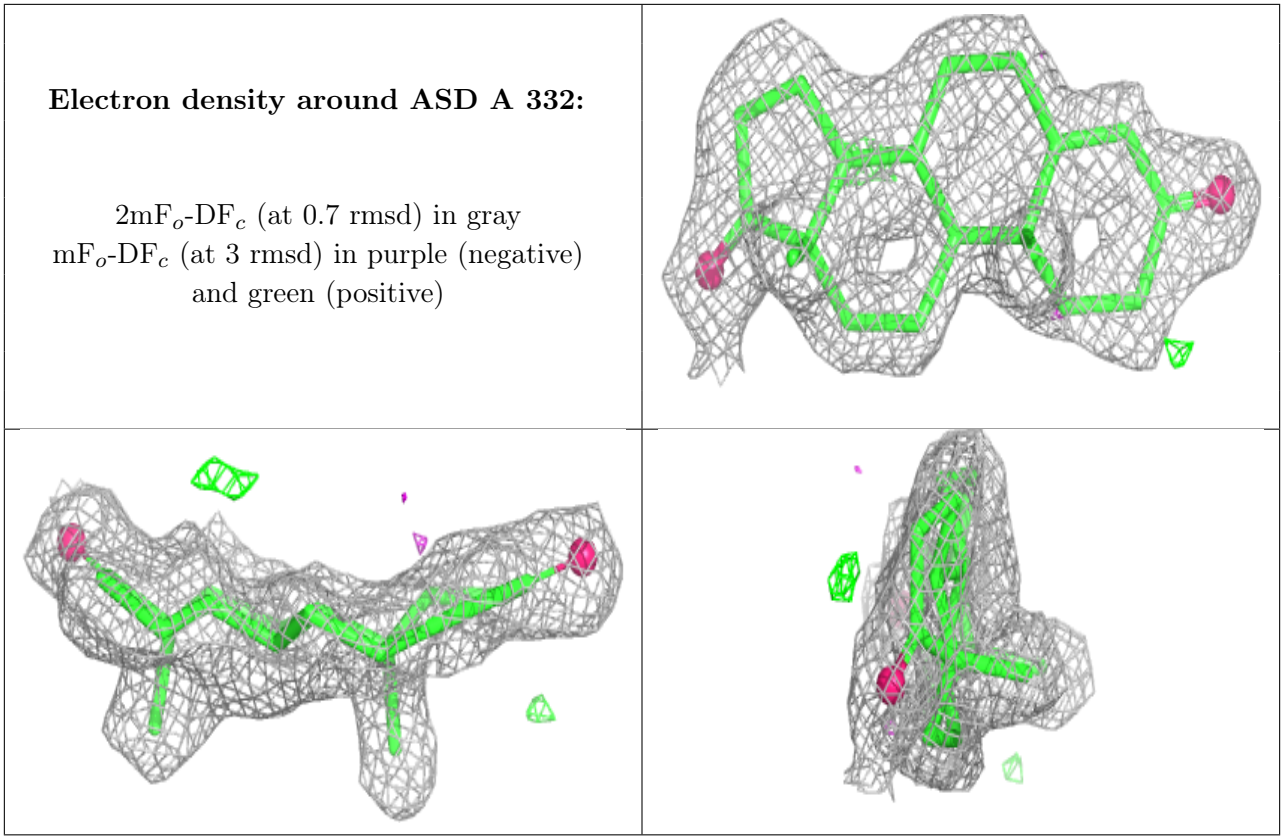
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	EDO	B	330	4/4	0.64	0.31	39,40,41,41	0
4	EDO	B	329	4/4	0.81	0.19	36,37,37,38	0
4	EDO	B	328	4/4	0.82	0.20	34,34,36,38	0
4	EDO	A	331	4/4	0.83	0.18	28,28,29,30	0
4	EDO	A	329	4/4	0.86	0.18	21,21,22,22	0
5	ASD	A	332	21/21	0.86	0.17	21,22,22,23	0
4	EDO	A	330	4/4	0.88	0.14	21,24,25,25	0
3	BME	A	328	4/4	0.89	0.22	38,38,39,39	0
5	ASD	B	331	21/21	0.91	0.15	20,23,28,30	0
2	NAP	B	327	48/48	0.96	0.11	6,11,16,25	0
2	NAP	A	327	48/48	0.96	0.10	4,8,11,18	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.

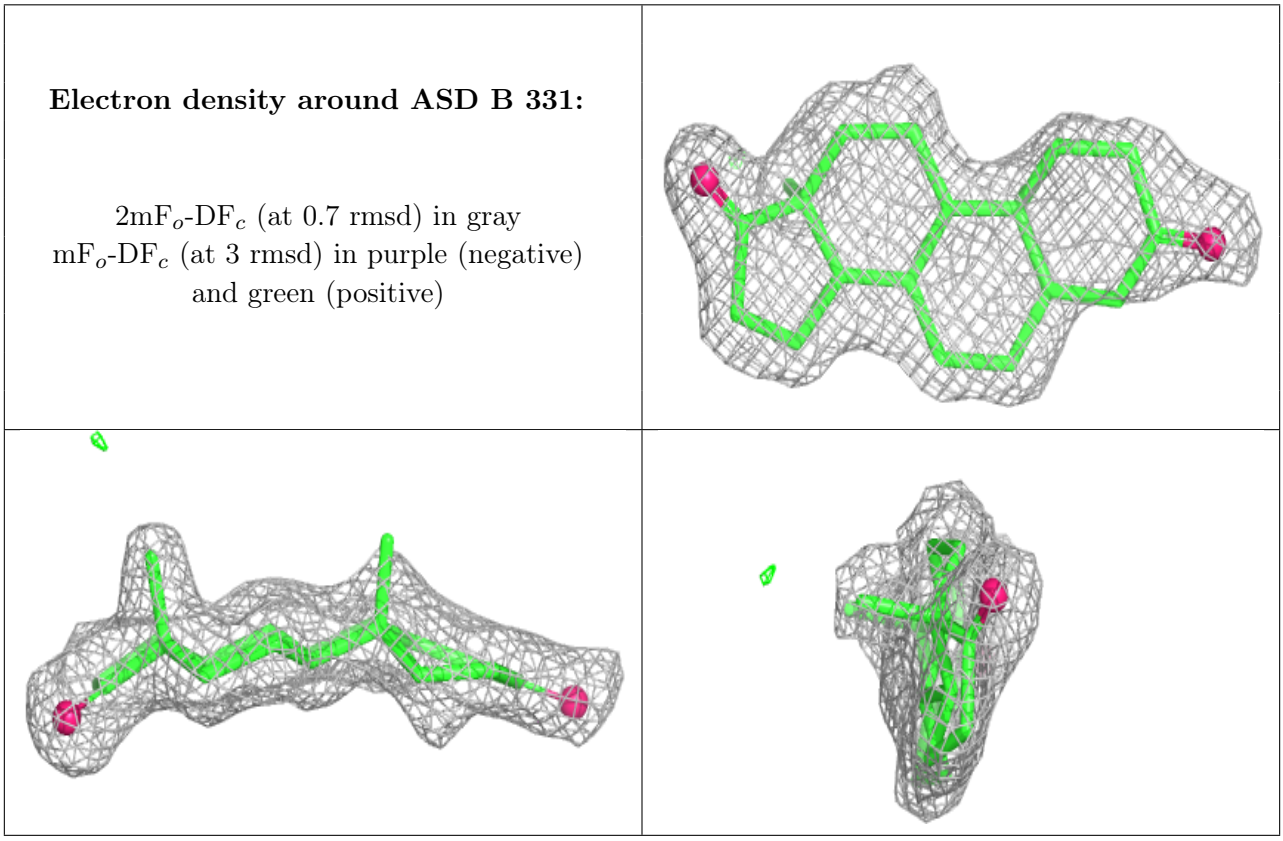
Electron density around ASD A 332:

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



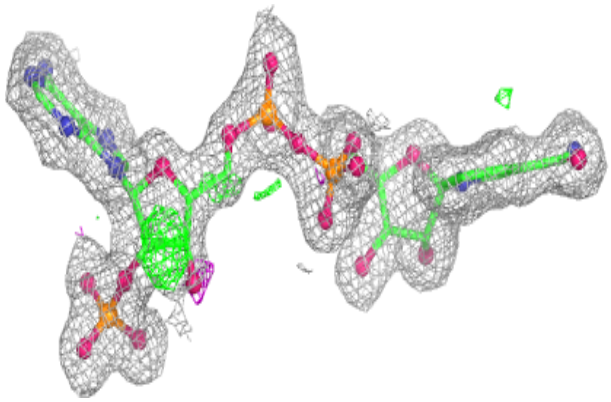
Electron density around ASD B 331:

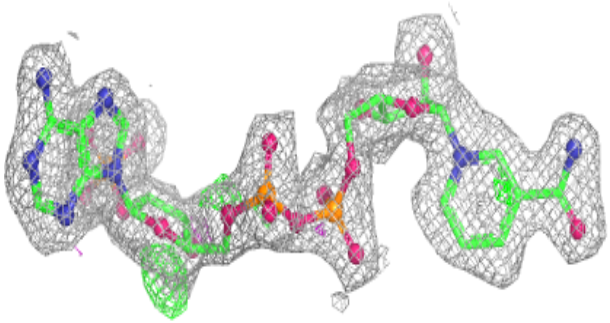
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

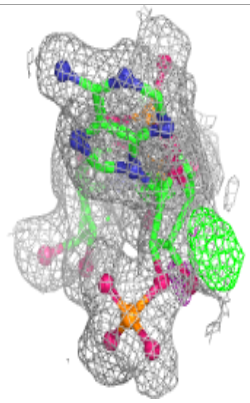


Electron density around NAP B 327:

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

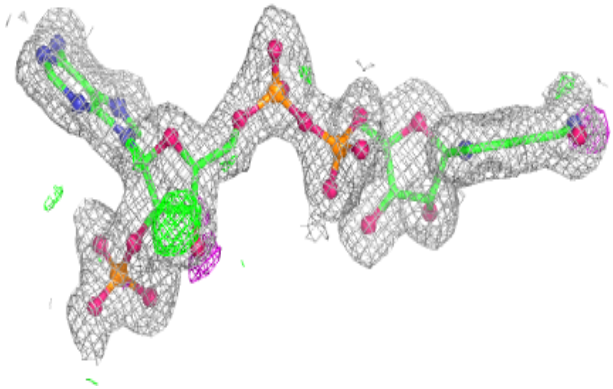


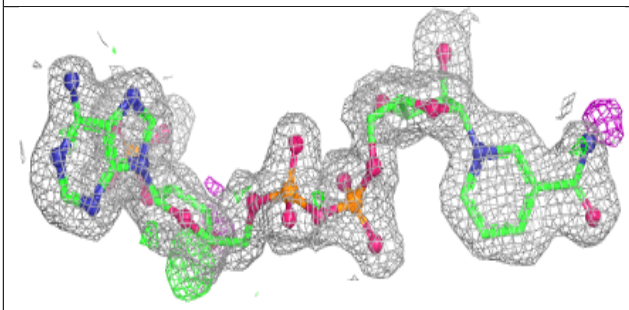


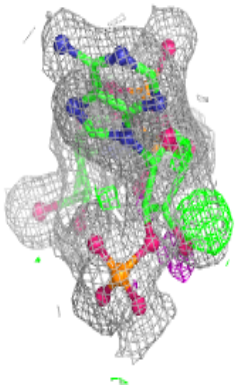


Electron density around NAP A 327:

$2mF_o$-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 (i)

There are no such residues in this entry.