



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:34 AM UTC

PDB ID : 3HM1 / pdb_00003hm1
Title : Crystal structure of human Estrogen Receptor Alpha Ligand-Binding Domain in complex with a Glucocorticoid Receptor Interacting Protein 1 Nr Box II Peptide and estrone ((8R,9S,13S,14S)-3-hydroxy-13-methyl-7,8,9,11,12,14,15,16-octahydro-6H-cyclopenta[a]phenanthren-17-one)
Authors : Rajan, S.S.; Kim, Y.; Vanek, K.; Liwanag, M.; Joachimiak, A.; Greene, G.L.
Deposited on : 2009-05-28
Resolution : 2.33 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](#) ①) 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

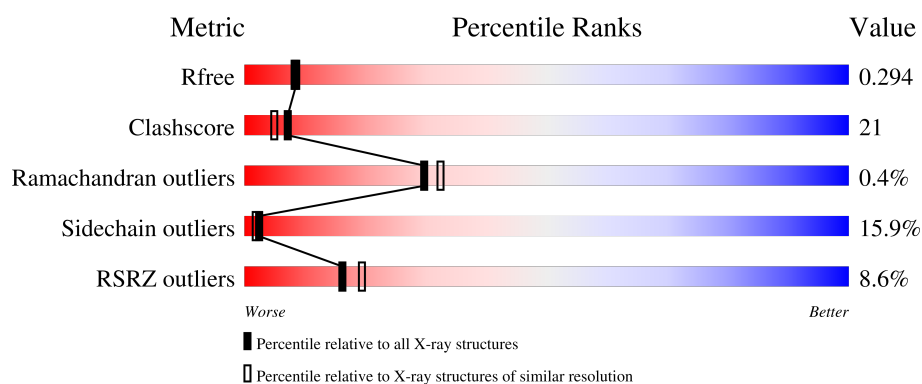
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.33 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	3031 (2.36-2.32)
Clashscore	190562	3127 (2.36-2.32)
Ramachandran outliers	187476	3095 (2.36-2.32)
Sidechain outliers	187428	3095 (2.36-2.32)
RSRZ outliers	180081	3033 (2.36-2.32)

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	253	9% 50% 32% 11% 7%
1	B	253	5% 59% 26% 8% 6%
2	C	13	15% 31% 31% 15% 23%
2	D	13	23% 31% 31% 8% 31%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4062 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 receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	235	Total	C	N	O	S	0	3	0
			1911	1223	327	339	22			
1	B	239	Total	C	N	O	S	0	2	0
			1933	1237	334	340	22			

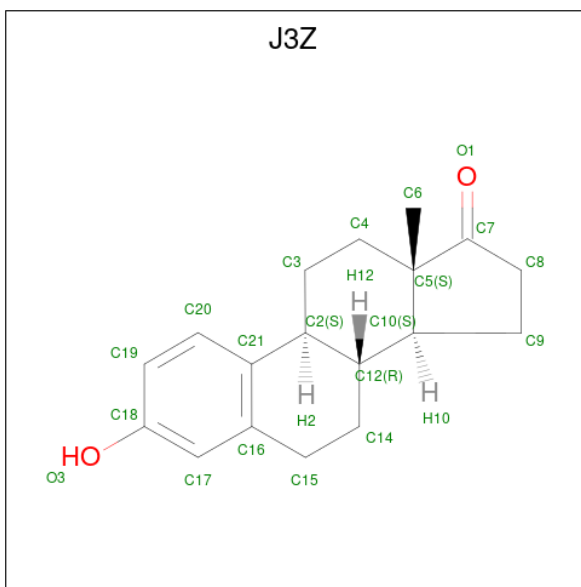
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	537	SER	TYR	engineered mutation	UNP P03372
B	537	SER	TYR	engineered mutation	UNP P03372

- Molecule 2 is a protein called Nuclear receptor coactivator 2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	10	Total	C	N	O	0	0	0
			89	57	19	13			
2	D	9	Total	C	N	O	0	0	0
			79	51	16	12			

- Molecule 3 is (9beta,13alpha)-3-hydroxyestra-1,3,5(10)-trien-17-one (CCD ID: J3Z) (formula: C₁₈H₂₂O₂).



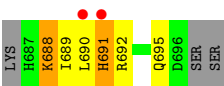
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			20	18	2		
3	B	1	Total	C	O	0	0
			20	18	2		

- Molecule 4 is water.

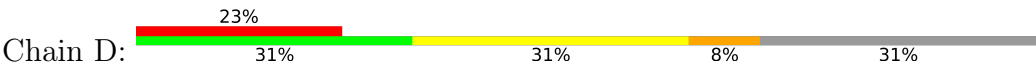
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	8	Total	O	0	0
			8	8		
4	B	2	Total	O	0	0
			2	2		

- Molecule 1: Estrogen receptor





● Molecule 2: Nuclear receptor coactivator 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.27Å 82.67Å 58.99Å 90.00° 109.18° 90.00°	Depositor
Resolution (Å)	33.37 – 2.33 33.37 – 2.33	Depositor EDS
% Data completeness (in resolution range)	86.0 (33.37-2.33) 86.1 (33.37-2.33)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.86 (at 2.34Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.231 , 0.289 0.241 , 0.294	Depositor DCC
R_{free} test set	966 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	39.3	Xtriage
Anisotropy	0.919	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 40.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.035 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4062	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.21% 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: CME, J3Z

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	0.39	0/1926	0.82	5/2595 (0.2%)
1	B	0.40	0/1946	0.75	2/2624 (0.1%)
2	C	0.29	0/90	0.53	0/119
2	D	0.32	0/79	0.67	0/104
All	All	0.39	0/4041	0.78	7/5442 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	335	ARG	CA-C-N	6.71	127.62	120.11
1	A	335	ARG	C-N-CA	6.71	127.62	120.11
1	B	335	ARG	CA-C-N	5.91	126.25	119.92
1	B	335	ARG	C-N-CA	5.91	126.25	119.92
1	A	323	GLU	CA-C-N	5.51	123.63	119.66
1	A	323	GLU	C-N-CA	5.51	123.63	119.66
1	A	315	MET	N-CA-C	-5.20	105.52	111.14

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1911	0	1952	96	0
1	B	1933	0	1982	77	0
2	C	89	0	95	4	0
2	D	79	0	88	6	0
3	A	20	0	22	1	0
3	B	20	0	22	3	0
4	A	8	0	0	1	0
4	B	2	0	0	0	0
All	All	4062	0	4161	171	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 21.

All (171) 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:D:688:LYS:HG2	2:D:691:HIS:HB2	1.46	0.95
1:B:335:ARG:HG3	1:B:336:PRO:HD2	1.59	0.83
1:B:522:MET:HA	1:B:522:MET:HE2	1.60	0.82
1:B:330:GLU:HG2	1:B:330:GLU:O	1.82	0.78
1:A:393:TRP:O	1:A:396:MET:HB2	1.84	0.77
1:A:331:TYR:HB2	1:A:345:LEU:HD21	1.67	0.76
1:A:481:LYS:HZ2	1:A:481:LYS:HB3	1.50	0.76
2:C:688:LYS:HG3	2:C:691:HIS:HB3	1.67	0.76
1:B:524:HIS:NE2	1:B:528:MET:HE3	2.01	0.76
1:B:524:HIS:CE1	1:B:528:MET:HE3	2.20	0.75
1:B:410:LEU:HD22	1:B:414:GLN:HG2	1.69	0.74
1:B:335:ARG:HG3	1:B:336:PRO:CD	2.17	0.74
2:D:688:LYS:HG3	2:D:691:HIS:H	1.54	0.73
1:A:543:MET:HE1	2:C:690:LEU:HD22	1.69	0.73
1:A:526:TYR:CE2	1:A:530:CME:HE3	2.25	0.71
1:B:304:ASN:H	1:B:304:ASN:HD22	1.35	0.71
1:A:377:HIS:NE2	1:A:460:THR:HG23	2.06	0.71
1:A:498:GLN:HA	1:A:501[A]:HIS:NE2	2.07	0.70
1:B:525:LEU:HD11	1:B:536:LEU:HD11	1.72	0.70
1:A:315:MET:HE3	1:A:481:LYS:HZ2	1.57	0.68
1:A:391:LEU:HA	1:A:394:ARG:NH1	2.08	0.68
1:B:474:HIS:O	1:B:478:VAL:HG12	1.93	0.68
1:A:508:LEU:HG	1:B:509:LEU:CD1	2.24	0.67
1:A:526:TYR:HE2	1:A:530:CME:HE3	1.57	0.67
1:A:385:GLU:HG2	1:A:514:ILE:HG22	1.76	0.66
1:A:315:MET:HE3	1:A:481:LYS:NZ	2.11	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:522:MET:HA	1:B:522:MET:CE	2.27	0.65
1:A:322:ALA:HA	1:A:363:ARG:NH2	2.12	0.64
1:A:481:LYS:HB3	1:A:481:LYS:NZ	2.13	0.64
1:A:354:LEU:O	1:A:358:ILE:HG12	1.97	0.64
1:B:383:TRP:NE1	1:B:543:MET:HB3	2.15	0.62
1:A:349:LEU:HD11	1:A:405:ALA:HB2	1.81	0.62
1:B:304:ASN:HD22	1:B:304:ASN:N	1.98	0.62
1:A:459[A]:TYR:CE2	1:B:513[A]:HIS:HB3	2.35	0.62
1:A:525:LEU:HD13	3:A:2:J3Z:H36	1.82	0.61
1:A:308:LEU:HD21	1:A:477:ARG:HB3	1.83	0.61
1:B:438:MET:HG3	1:B:438:MET:O	2.00	0.61
1:B:443:GLU:HB3	1:B:489:LEU:HD21	1.83	0.60
1:A:487:ILE:CD1	1:A:504:LEU:HD22	2.32	0.59
1:B:347:THR:HG21	1:B:536:LEU:HD21	1.85	0.59
1:A:429:LEU:HD12	1:A:430:ALA:N	2.17	0.58
2:D:688:LYS:CG	2:D:691:HIS:H	2.16	0.58
1:B:343:MET:HE2	1:B:528:MET:HG3	1.85	0.57
1:B:497:LEU:HD22	1:B:501[A]:HIS:HE2	1.68	0.57
1:B:421:MET:HE1	3:B:1:J3Z:H10	1.87	0.56
1:B:496:THR:O	1:B:500:GLN:HG3	2.04	0.56
1:A:360:TRP:CZ2	1:A:449:LYS:HG2	2.40	0.56
1:A:371:THR:HG22	1:A:373:HIS:H	1.71	0.56
1:B:346:LEU:HB3	3:B:1:J3Z:H23	1.87	0.56
1:B:388:MET:HG2	1:B:428:LEU:HD21	1.87	0.56
1:A:391:LEU:HD11	1:A:402:LEU:HB3	1.88	0.56
1:B:437:MET:O	1:B:437:MET:HG2	2.06	0.55
1:B:549:LEU:HD22	1:B:549:LEU:N	2.22	0.55
1:A:528:MET:HE3	1:A:536:LEU:HD21	1.88	0.55
1:A:329:SER:H	1:A:407:ASN:HD21	1.52	0.55
1:B:526:TYR:CD2	1:B:549:LEU:HD23	2.42	0.55
1:A:459[A]:TYR:CE1	1:B:513[A]:HIS:CD2	2.95	0.54
1:B:339:GLU:HA	1:B:418:VAL:HG22	1.90	0.53
1:B:311:THR:HG23	1:B:314:GLN:OE1	2.08	0.53
1:A:527:SER:O	1:A:531:LYS:HG2	2.08	0.53
1:B:377:HIS:CD2	1:B:461:PHE:CE2	2.97	0.53
1:B:497:LEU:HD22	1:B:501[A]:HIS:NE2	2.23	0.53
1:A:314:GLN:HA	1:A:317:SER:HB2	1.90	0.53
1:A:456:SER:HA	1:A:515:ARG:NH2	2.25	0.52
1:A:344:GLY:HA2	1:A:534:VAL:HG21	1.90	0.52
1:B:427:MET:HE3	1:B:517:MET:HG2	1.91	0.52
1:A:351:ASP:OD2	1:A:540:LEU:HB2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:329:SER:H	1:A:407:ASN:ND2	2.07	0.52
1:A:498:GLN:HA	1:A:501[A]:HIS:CD2	2.45	0.52
1:A:329:SER:HB3	1:A:407:ASN:HD21	1.75	0.52
1:B:522:MET:HE1	1:B:544:LEU:HD22	1.91	0.51
1:A:344:GLY:CA	1:A:534:VAL:HG21	2.42	0.50
1:B:329:SER:C	1:B:331:TYR:H	2.19	0.50
1:A:487:ILE:HD13	1:A:504:LEU:HD22	1.92	0.50
1:B:486:LEU:HD23	1:B:504:LEU:HD12	1.92	0.50
1:A:523:GLU:CD	1:A:548[B]:ARG:HH12	2.20	0.50
1:A:352:ARG:NH2	4:A:5:HOH:O	2.43	0.50
1:A:474:HIS:O	1:A:478:VAL:HG12	2.11	0.50
1:B:305:SER:C	1:B:307:ALA:H	2.20	0.50
1:A:459[B]:TYR:HE2	1:B:434:ARG:HG2	1.76	0.49
1:A:342:MET:HE2	1:A:418:VAL:HG23	1.93	0.49
1:B:424:ILE:HA	1:B:427:MET:HE2	1.94	0.49
1:A:498:GLN:HA	1:A:501[A]:HIS:CE1	2.48	0.49
1:B:382:ALA:HB2	1:B:456:SER:OG	2.12	0.49
1:A:342:MET:HE2	1:A:418:VAL:CG2	2.43	0.48
1:B:506:GLN:O	1:B:510:ILE:HG13	2.13	0.48
1:A:416:LYS:O	1:A:417:CME:C	2.61	0.48
1:A:497:LEU:O	1:A:500:GLN:HB2	2.14	0.48
1:A:421:MET:CE	1:A:421:MET:HA	2.43	0.48
1:B:498:GLN:HA	1:B:501[B]:HIS:ND1	2.28	0.48
1:A:393:TRP:HB2	1:A:445:PHE:CE1	2.48	0.48
1:B:514:ILE:HA	1:B:517:MET:HE2	1.96	0.48
1:A:332:ASP:C	1:A:334:THR:H	2.21	0.48
1:A:391:LEU:HA	1:A:394:ARG:HH12	1.77	0.48
1:A:376:VAL:O	1:A:380:GLU:HG3	2.14	0.48
1:A:514:ILE:HA	1:A:517:MET:HE2	1.95	0.48
1:B:525:LEU:HG	1:B:544:LEU:CD2	2.44	0.48
1:A:525:LEU:HD12	1:A:528:MET:CE	2.44	0.48
1:A:528:MET:HG3	1:A:533:VAL:HG23	1.95	0.48
1:B:487:ILE:HG23	1:B:500:GLN:NE2	2.29	0.47
1:A:331:TYR:CB	1:A:345:LEU:HD21	2.43	0.47
1:B:498:GLN:HA	1:B:501[B]:HIS:CE1	2.50	0.47
1:A:349:LEU:O	1:A:353:GLU:HG3	2.15	0.47
1:B:385:GLU:HG3	1:B:518:SER:HB2	1.98	0.46
1:A:343:MET:SD	1:A:418:VAL:HG21	2.55	0.46
1:B:410:LEU:HD23	1:B:414:GLN:HE21	1.80	0.46
1:B:372:LEU:HD21	2:D:695:GLN:CD	2.40	0.46
1:B:522:MET:HE1	1:B:525:LEU:HD23	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:538:ASP:O	1:A:538:ASP:OD2	2.34	0.46
1:A:379:LEU:O	1:A:383:TRP:HB3	2.16	0.46
1:B:305:SER:C	1:B:307:ALA:N	2.74	0.46
1:A:352:ARG:O	1:A:355:VAL:HG13	2.16	0.45
1:A:385:GLU:CG	1:A:514:ILE:HG22	2.44	0.45
1:A:367:PHE:CE1	1:A:453:LEU:HD11	2.51	0.45
1:A:421:MET:HA	1:A:421:MET:HE3	1.97	0.45
1:A:308:LEU:HD21	1:A:477:ARG:HE	1.82	0.45
1:A:351:ASP:OD1	1:A:537:SER:HB3	2.16	0.45
1:A:458:VAL:HG23	1:A:475:ILE:HD12	1.99	0.44
1:B:307:ALA:HA	1:B:310:LEU:CD1	2.47	0.44
1:B:403:LEU:HD12	1:B:409:LEU:HD11	1.98	0.44
2:D:688:LYS:HD2	2:D:690:LEU:HB3	2.00	0.44
1:A:409:LEU:H	1:A:409:LEU:CD1	2.31	0.44
2:C:691:HIS:O	2:C:695:GLN:HG3	2.18	0.44
1:A:459[B]:TYR:CE2	1:B:434:ARG:HG2	2.51	0.44
1:A:396:MET:O	1:A:436:ARG:HD3	2.18	0.44
1:A:458:VAL:HG12	1:A:459[B]:TYR:CD1	2.53	0.44
1:B:385:GLU:HB3	1:B:452:ILE:HG21	2.00	0.43
1:A:339:GLU:HG3	1:A:340:ALA:N	2.34	0.43
1:A:421:MET:CE	1:A:424:ILE:HD12	2.48	0.43
1:A:456:SER:HA	1:A:515:ARG:HH22	1.83	0.43
2:C:691:HIS:CD2	2:C:692:ARG:HG3	2.54	0.43
1:B:439:ASN:O	1:B:440:LEU:C	2.62	0.43
1:A:343:MET:HG2	1:A:418:VAL:HG21	2.01	0.43
1:B:510:ILE:O	1:B:514:ILE:HG13	2.18	0.43
1:B:472:LYS:N	1:B:472:LYS:HE2	2.34	0.43
1:A:459[B]:TYR:HE2	1:B:434:ARG:CG	2.32	0.43
1:A:497:LEU:HD13	1:B:497:LEU:HD21	2.00	0.43
1:A:387:LEU:HD12	1:A:387:LEU:HA	1.90	0.43
1:A:520:LYS:HA	1:A:520:LYS:HD3	1.83	0.42
1:A:339:GLU:HG3	1:A:340:ALA:H	1.84	0.42
1:B:306:LEU:O	1:B:306:LEU:HG	2.19	0.42
1:B:541:LEU:HD12	1:B:541:LEU:HA	1.87	0.42
2:D:688:LYS:CG	2:D:691:HIS:HB2	2.33	0.42
1:A:315:MET:O	1:A:319:LEU:HG	2.20	0.42
1:A:380:GLU:O	1:A:547:HIS:HE1	2.03	0.42
1:A:409:LEU:H	1:A:409:LEU:HD12	1.84	0.42
1:B:525:LEU:HD22	3:B:1:J3Z:H36	2.02	0.42
1:B:396:MET:HE3	1:B:396:MET:HB3	1.75	0.42
1:A:379:LEU:HD12	1:A:379:LEU:HA	1.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:354:LEU:HD12	1:B:354:LEU:HA	1.83	0.42
1:B:544:LEU:HD13	1:B:544:LEU:HA	1.78	0.42
1:B:548:ARG:C	1:B:549:LEU:HD13	2.45	0.42
1:A:392:VAL:HG13	1:A:432:SER:HA	2.02	0.41
1:A:409:LEU:HD13	1:A:409:LEU:O	2.20	0.41
1:B:388:MET:HE1	1:B:517:MET:O	2.20	0.41
1:A:350:ALA:O	1:A:354:LEU:HB2	2.20	0.41
1:A:434:ARG:HE	1:A:437:MET:HE3	1.84	0.41
1:B:310:LEU:O	1:B:481:LYS:HE2	2.20	0.41
1:A:508:LEU:HD12	1:A:508:LEU:HA	1.87	0.41
1:A:515:ARG:HD2	1:B:516:HIS:HB2	2.02	0.41
1:A:391:LEU:HD12	1:A:391:LEU:O	2.21	0.41
1:B:303:LYS:HB2	1:B:303:LYS:HE3	1.94	0.41
1:B:456:SER:HA	1:B:515:ARG:NH2	2.36	0.41
1:A:501[B]:HIS:CD2	1:B:501[B]:HIS:HB3	2.56	0.41
1:A:407:ASN:H	1:A:407:ASN:HD22	1.69	0.41
1:B:342:MET:HE3	1:B:342:MET:HB2	1.86	0.40
1:B:403:LEU:CD1	1:B:409:LEU:HD11	2.51	0.40
1:A:323:GLU:HA	1:A:324:PRO:HD3	1.84	0.40
1:A:459[A]:TYR:CE1	1:B:513[A]:HIS:HD2	2.39	0.40
1:B:372:LEU:HD22	1:B:375:GLN:NE2	2.37	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	230/253 (91%)	224 (97%)	5 (2%)	1 (0%)	30	32
1	B	234/253 (92%)	225 (96%)	8 (3%)	1 (0%)	30	32
2	C	8/13 (62%)	8 (100%)	0	0	100	100
2	D	7/13 (54%)	7 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	479/532 (90%)	464 (97%)	13 (3%)	2 (0%)	30	32

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	396	MET
1	A	333	PRO

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	211/226 (93%)	176 (83%)	35 (17%)	2	1
1	B	214/226 (95%)	183 (86%)	31 (14%)	3	2
2	C	10/13 (77%)	7 (70%)	3 (30%)	0	0
2	D	9/13 (69%)	7 (78%)	2 (22%)	1	0
All	All	444/478 (93%)	373 (84%)	71 (16%)	2	2

All (71) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	310	LEU
1	A	311	THR
1	A	317	SER
1	A	321	ASP
1	A	327	LEU
1	A	330	GLU
1	A	335	ARG
1	A	339	GLU
1	A	345	LEU
1	A	349	LEU
1	A	354	LEU
1	A	355	VAL
1	A	363	ARG

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Mol	Chain	Res	Type
1	A	373	HIS
1	A	379	LEU
1	A	385	GLU
1	A	407	ASN
1	A	409	LEU
1	A	416	LYS
1	A	419	GLU
1	A	421	MET
1	A	434	ARG
1	A	458	VAL
1	A	459[A]	TYR
1	A	459[B]	TYR
1	A	470	GLU
1	A	478	VAL
1	A	481	LYS
1	A	487	ILE
1	A	492	LYS
1	A	498	GLN
1	A	508	LEU
1	A	539	LEU
1	A	542	GLU
1	A	543	MET
1	B	303	LYS
1	B	304	ASN
1	B	306	LEU
1	B	308	LEU
1	B	330	GLU
1	B	347	THR
1	B	354	LEU
1	B	372	LEU
1	B	373	HIS
1	B	378	LEU
1	B	396	MET
1	B	397	GLU
1	B	403	LEU
1	B	414	GLN
1	B	424	ILE
1	B	437	MET
1	B	438	MET
1	B	460	THR
1	B	462	LEU
1	B	472	LYS

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Mol	Chain	Res	Type
1	B	478	VAL
1	B	489	LEU
1	B	497	LEU
1	B	508	LEU
1	B	509	LEU
1	B	522	MET
1	B	523	GLU
1	B	533	VAL
1	B	542	GLU
1	B	544	LEU
1	B	549	LEU
2	C	688	LYS
2	C	689	ILE
2	C	691	HIS
2	D	688	LYS
2	D	693	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	356	HIS
1	A	375	GLN
1	A	398	HIS
1	A	407	ASN
1	A	439	ASN
1	A	441	GLN
1	A	455	ASN
1	A	488	HIS
1	A	499	GLN
1	A	502	GLN
1	A	513	HIS
1	A	516	HIS
1	A	547	HIS
1	B	304	ASN
1	B	375	GLN
1	B	377	HIS
1	B	414	GLN
1	B	455	ASN
1	B	500	GLN
1	B	502	GLN
1	B	506	GLN
1	B	532	ASN

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Mol	Chain	Res	Type
2	C	695	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	CME	A	381	1	8,9,10	0.52	0	6,9,11	0.57	0
1	CME	A	417	1	8,9,10	0.61	0	6,9,11	1.61	1 (16%)
1	CME	B	381	1	8,9,10	0.55	0	6,9,11	2.18	1 (16%)
1	CME	B	530	1	8,9,10	0.40	0	6,9,11	0.68	0
1	CME	B	417	1	8,9,10	0.51	0	6,9,11	0.67	0
1	CME	A	530	1	8,9,10	0.50	0	6,9,11	0.63	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
1	CME	A	381	1	-	1/5/8/10	-
1	CME	A	417	1	-	2/5/8/10	-
1	CME	B	381	1	-	1/5/8/10	-
1	CME	B	530	1	-	0/5/8/10	-
1	CME	B	417	1	-	0/5/8/10	-
1	CME	A	530	1	-	2/5/8/10	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	381	CME	CB-SG-SD	-4.79	91.46	103.86
1	A	417	CME	CB-SG-SD	3.38	112.61	103.86

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	417	CME	CA-CB-SG-SD
1	B	381	CME	N-CA-CB-SG
1	A	530	CME	CA-CB-SG-SD
1	A	381	CME	CZ-CE-SD-SG
1	A	417	CME	CZ-CE-SD-SG
1	A	530	CME	CZ-CE-SD-SG

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	417	CME	1	0
1	A	530	CME	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	J3Z	A	2	-	23,23,23	0.91	0	36,36,36	1.54	8 (22%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	J3Z	B	1	-	23,23,23	0.94	1 (4%)	36,36,36	1.53	9 (25%)

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	J3Z	A	2	-	-	-	0/4/4/4
3	J3Z	B	1	-	-	-	0/4/4/4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1	J3Z	C21-C2	-2.04	1.49	1.52

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1	J3Z	C2-C12-C10	-3.77	103.44	108.68
3	A	2	J3Z	C8-C7-C5	3.70	112.24	108.56
3	A	2	J3Z	C2-C12-C10	-3.34	104.03	108.68
3	A	2	J3Z	C9-C10-C5	3.27	106.89	104.08
3	B	1	J3Z	C5-C10-C12	-2.94	110.04	113.13
3	B	1	J3Z	C8-C7-C5	2.84	111.38	108.56
3	B	1	J3Z	C9-C8-C7	-2.79	103.01	105.70
3	B	1	J3Z	C15-C14-C12	2.70	115.11	110.61
3	B	1	J3Z	C6-C5-C7	2.62	110.01	105.18
3	A	2	J3Z	C14-C12-C2	2.60	111.76	109.27
3	A	2	J3Z	C5-C10-C12	-2.56	110.44	113.13
3	A	2	J3Z	C9-C8-C7	-2.31	103.47	105.70
3	B	1	J3Z	C14-C12-C2	2.20	111.38	109.27
3	B	1	J3Z	C9-C10-C5	2.14	105.92	104.08
3	B	1	J3Z	O1-C7-C8	-2.13	123.16	125.75
3	A	2	J3Z	C15-C14-C12	2.06	114.05	110.61
3	A	2	J3Z	O1-C7-C8	-2.03	123.29	125.75

There are no chirality outliers.

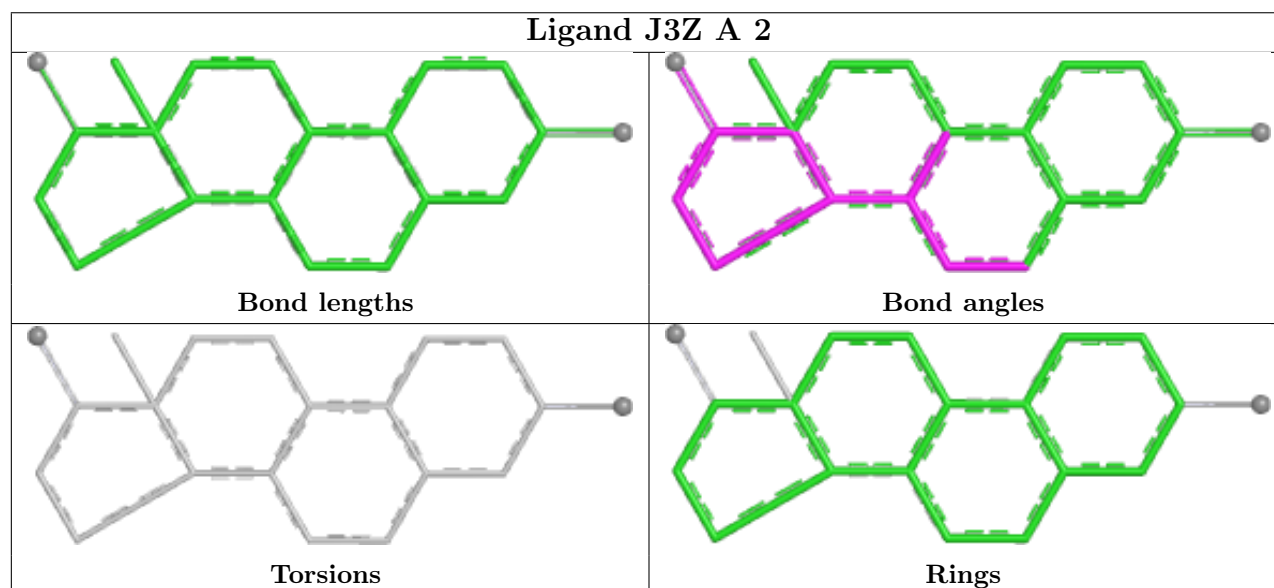
There are no torsion outliers.

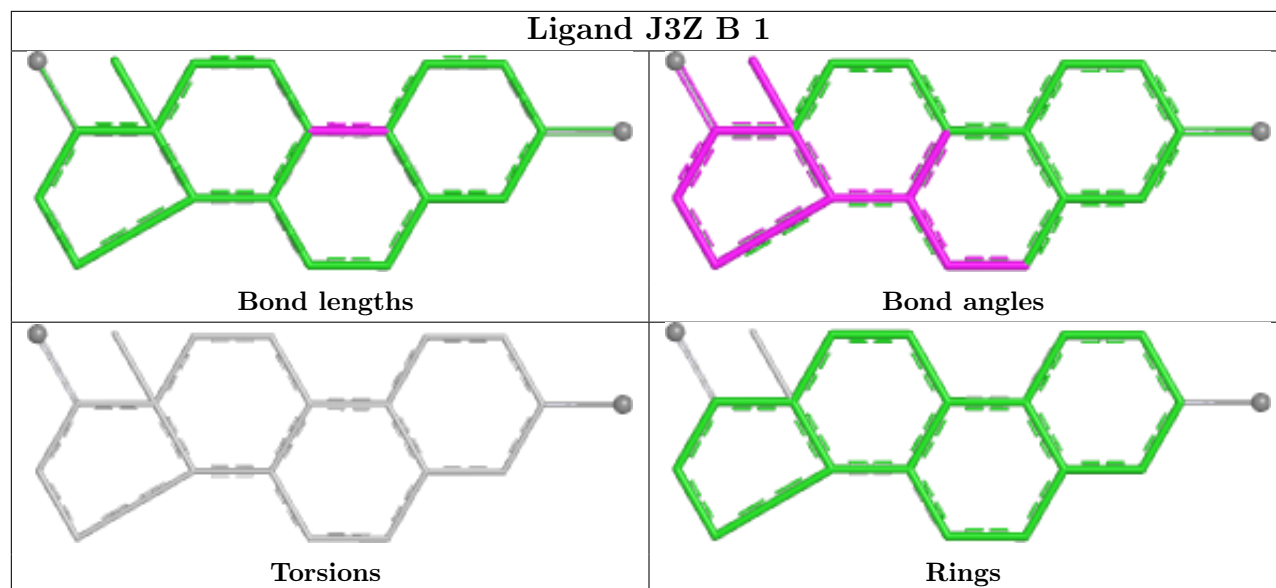
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2	J3Z	1	0
3	B	1	J3Z	3	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	232/253 (91%)	0.76	24 (10%)	12 14	36, 65, 115, 139	3 (1%)
1	B	236/253 (93%)	0.63	13 (5%)	30 36	31, 64, 103, 137	2 (0%)
2	C	10/13 (76%)	1.67	2 (20%)	3 3	92, 112, 126, 128	0
2	D	9/13 (69%)	1.46	3 (33%)	1 1	77, 94, 106, 108	0
All	All	487/532 (91%)	0.73	42 (8%)	16 19	31, 65, 112, 139	5 (1%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	336	PRO	5.3
1	B	460	THR	4.5
1	A	306	LEU	4.1
1	A	351	ASP	4.1
1	A	470	GLU	3.7
1	B	462	LEU	3.5
1	A	422	VAL	3.3
1	A	309	SER	3.2
1	B	334	THR	3.2
1	A	418	VAL	3.1
2	C	690	LEU	3.0
1	B	461	PHE	2.9
1	A	371	THR	2.9
1	B	309	SER	2.8
1	A	333	PRO	2.8
1	A	310	LEU	2.7
1	A	525	LEU	2.6
1	A	538	ASP	2.6
2	C	691	HIS	2.6
1	A	548[A]	ARG	2.5
1	A	461	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	458	VAL	2.4
1	B	459	TYR	2.4
1	A	473	ASP	2.4
1	B	351	ASP	2.4
1	B	548	ARG	2.4
1	A	533	VAL	2.3
2	D	695	GLN	2.3
1	A	368	VAL	2.3
1	B	338	SER	2.2
1	B	549	LEU	2.2
1	A	334	THR	2.2
1	B	371	THR	2.2
2	D	696	ASP	2.2
1	B	550	HIS	2.2
1	A	459[A]	TYR	2.1
1	A	307	ALA	2.1
1	A	420	GLY	2.1
1	A	410	LEU	2.1
2	D	689	ILE	2.1
1	B	333	PRO	2.1
1	A	337	PHE	2.1

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
1	CME	A	381	10/11	0.75	0.15	42,53,77,129	4
1	CME	A	530	10/11	0.77	0.16	88,90,126,196	3
1	CME	B	530	10/11	0.78	0.13	61,67,109,109	0
1	CME	B	417	10/11	0.79	0.12	69,78,132,155	4
1	CME	B	381	10/11	0.80	0.16	42,75,107,192	3
1	CME	A	417	10/11	0.85	0.14	87,115,118,119	4

6.3 Carbohydrates [i](#)

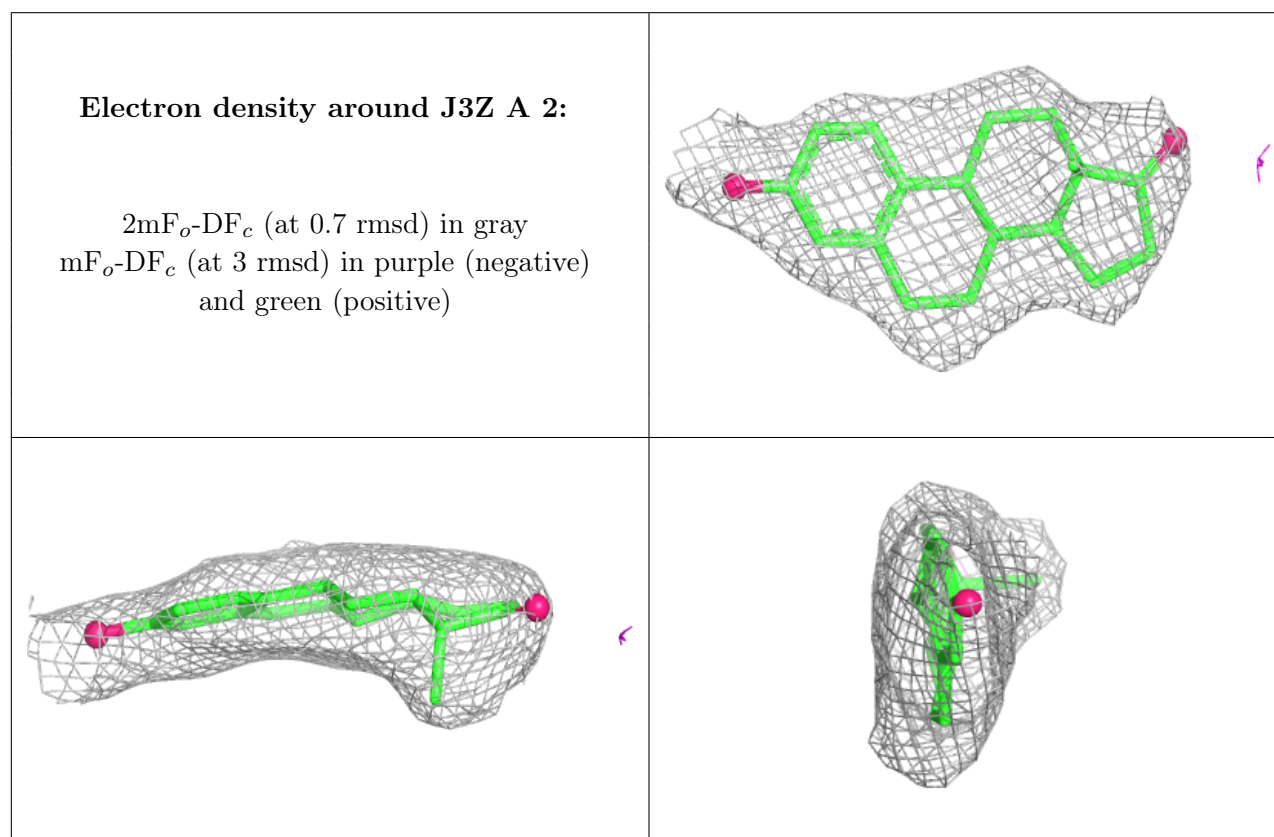
There are no oligosaccharides in this entry.

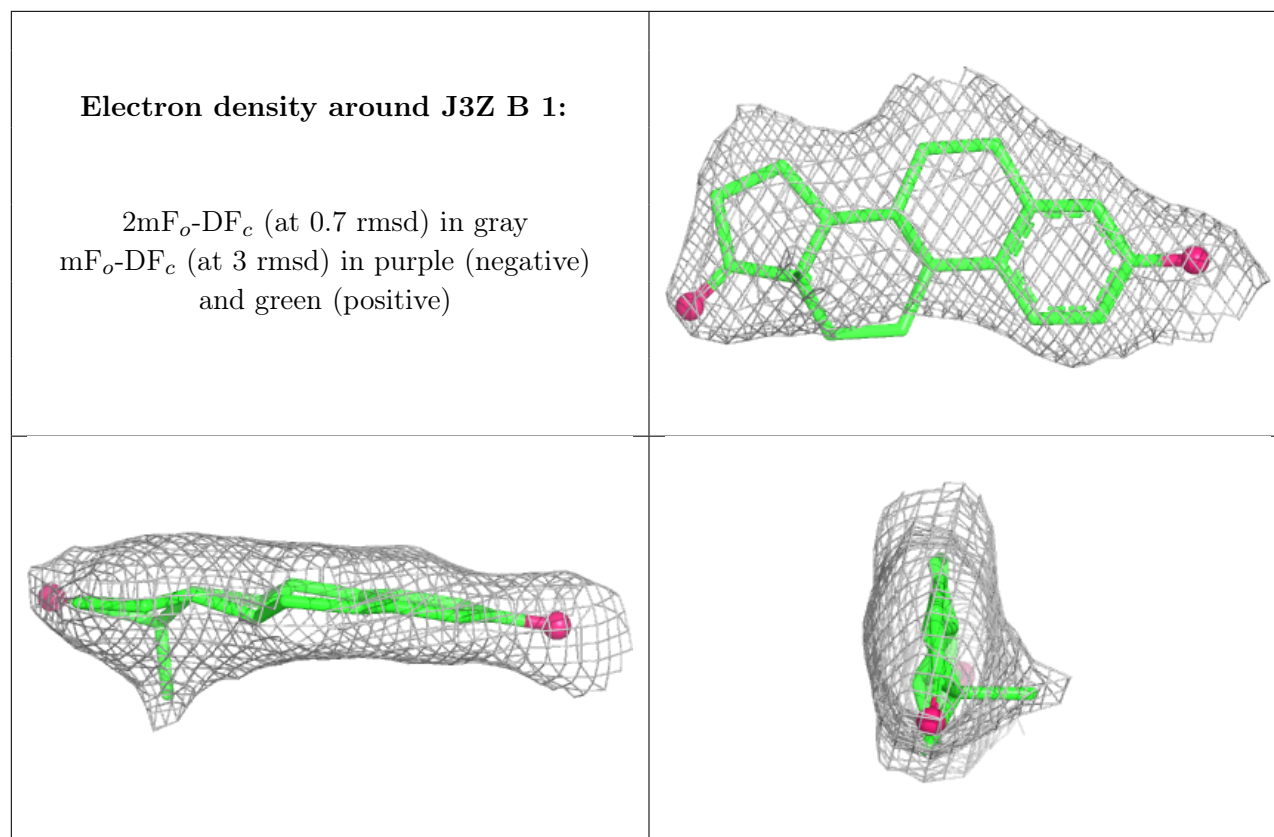
6.4 Ligands ⓘ

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
3	J3Z	A	2	20/20	0.93	0.08	38,51,64,65	0
3	J3Z	B	1	20/20	0.95	0.08	41,51,61,63	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.