



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 12, 2016 – 03:54 PM EDT

PDB ID : 5DX3
Title : Estrogen Receptor Alpha Ligand Binding Domain Y537S Mutant in Complex with Stapled Peptide SRC2-P3 and Estradiol
Authors : Speltz, T.E.; Fanning, S.W.; Mayne, C.G.; Tajkhorshid, E.; Greene, G.L.; Moore, T.W.
Deposited on : 2015-09-23
Resolution : 2.09 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20027790
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20027790

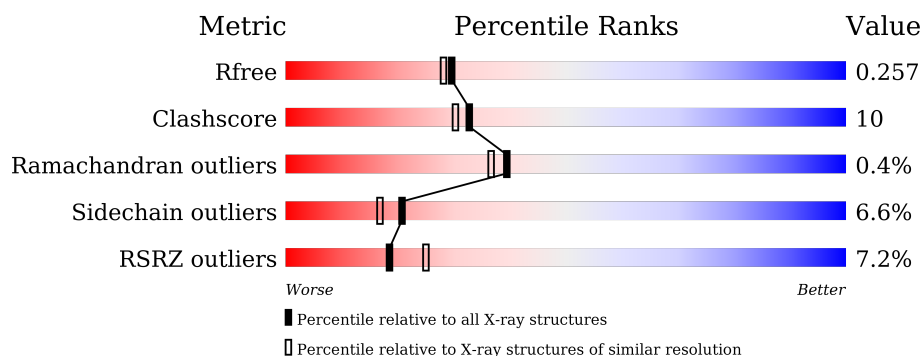
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.09 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}	91344	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	261	<div> <div>3%</div> <div>75%</div> <div>11%</div> <div>11%</div> </div>
2	B	261	<div> <div>9%</div> <div>66%</div> <div>17%</div> <div>14%</div> </div>
3	C	13	<div> <div>8%</div> <div>69%</div> <div>15%</div> <div>15%</div> </div>
3	D	13	<div> <div>23%</div> <div>23%</div> <div>46%</div> <div>31%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4185 atoms, of which 216 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	231	Total	C	N	O	S	2	5	0
			1841	1183	317	322	19			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	294	MET	-	initiating methionine	UNP P03372
A	295	ASP	-	expression tag	UNP P03372
A	296	PRO	-	expression tag	UNP P03372
A	537	SER	TYR	engineered mutation	UNP P03372

- Molecule 2 is a protein called Estrogen receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	224	Total	C	N	O	S	0	0	0
			1743	1113	303	309	18			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	294	MET	-	initiating methionine	UNP P03372
B	295	ASP	-	expression tag	UNP P03372
B	296	PRO	-	expression tag	UNP P03372
B	417	MET	CYS	conflict	UNP P03372
B	537	SER	TYR	engineered mutation	UNP P03372

- Molecule 3 is a protein called Stapled Peptide SRC2-P3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	13	Total	C	H	N	O	0	0	1
			210	65	108	21	16			

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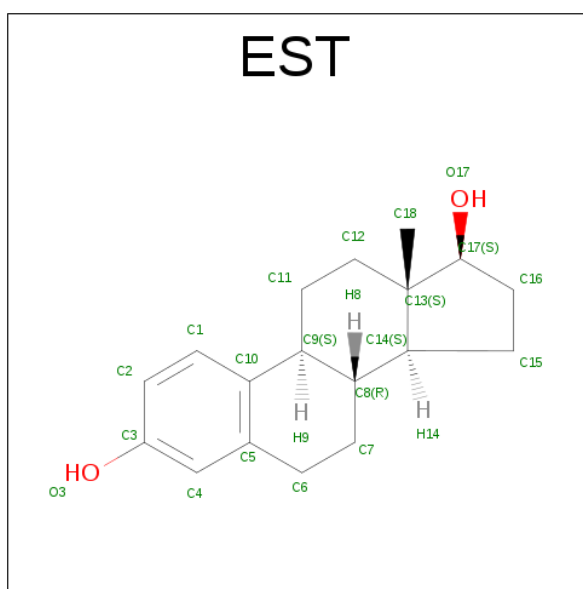
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	13	Total	C	H	N	O	0	0	1
			210	65	108	21	16			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1	ACE	-	expression tag	UNP Q15596
C	4	5GM	ILE	conflict	UNP Q15596
C	13	NH2	-	expression tag	UNP Q15596
D	1	ACE	-	expression tag	UNP Q15596
D	4	5GM	ILE	conflict	UNP Q15596
D	13	NH2	-	expression tag	UNP Q15596

- Molecule 4 is ESTRADIOL (three-letter code: EST) (formula: $C_{18}H_{24}O_2$).



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

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Cl	0	0
			1	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	89	Total 89	O 89	0	0
6	B	49	Total 49	O 49	0	0
6	C	1	Total 1	O 1	0	0
6	D	1	Total 1	O 1	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.16 Å 85.85 Å 58.76 Å 90.00° 109.09° 90.00°	Depositor
Resolution (Å)	26.54 – 2.09 26.54 – 2.09	Depositor EDS
% Data completeness (in resolution range)	95.7 (26.54-2.09) 95.7 (26.54-2.09)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 2.10 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.210 , 0.257 0.212 , 0.257	Depositor DCC
R_{free} test set	1445 reflections (4.84%)	DCC
Wilson B-factor (Å ²)	27.8	Xtriage
Anisotropy	0.131	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.028 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4185	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.49% 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: EST, ACE, 5GM, CL, NH2, MK8

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	0.35	1/1883 (0.1%)	0.57	2/2548 (0.1%)
2	B	0.37	1/1770 (0.1%)	0.57	2/2390 (0.1%)
3	C	1.74	0/79	1.65	2/101 (2.0%)
3	D	1.73	3/79 (3.8%)	1.48	1/101 (1.0%)
All	All	0.50	5/3811 (0.1%)	0.64	7/5140 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	362	LYS	CD-CE	8.08	1.71	1.51
3	D	5	LEU	CA-C	-7.49	1.33	1.52
3	D	7	ARG	CA-C	-6.13	1.37	1.52
1	A	353	GLU	CG-CD	6.09	1.61	1.51
3	D	6	HIS	CA-C	-5.76	1.38	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	362	LYS	CD-CE-NZ	-8.45	92.26	111.70
2	B	310	LEU	CA-CB-CG	6.88	131.12	115.30
3	D	7	ARG	NE-CZ-NH1	-6.73	116.93	120.30
3	C	11	ASP	CB-CG-OD2	-6.58	112.38	118.30
1	A	511[A]	LEU	CA-CB-CG	5.57	128.12	115.30
1	A	511[B]	LEU	CA-CB-CG	5.57	128.12	115.30
3	C	2	HIS	N-CA-CB	5.08	119.74	110.60

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	1841	0	1884	30	1
2	B	1743	0	1770	48	1
3	C	102	108	92	1	1
3	D	102	108	92	16	0
4	A	20	0	23	1	0
4	B	20	0	23	3	0
5	A	1	0	0	0	0
6	A	89	0	0	5	2
6	B	49	0	0	1	1
6	C	1	0	0	0	0
6	D	1	0	0	0	0
All	All	3969	216	3884	80	3

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 (80) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:362:LYS:CE	3:D:9:LEU:O	1.80	1.27
2:B:362:LYS:HE2	3:D:9:LEU:O	1.48	1.12
1:A:353:GLU:OE1	4:A:601:EST:O3	1.80	0.99
2:B:362:LYS:HE3	3:D:9:LEU:O	1.73	0.88
3:D:11:ASP:O	3:D:13:NH2:N	2.10	0.85
1:A:363:ARG:HH21	1:A:363:ARG:HG2	1.49	0.78
1:A:437:MET:HG2	6:A:788:HOH:O	1.85	0.77
2:B:315:MET:CE	2:B:481:LYS:HD2	2.15	0.76
2:B:421:MET:CE	4:B:1000:EST:H161	2.17	0.75
2:B:362:LYS:HD3	3:D:9:LEU:HD12	1.72	0.70
1:A:385:GLU:HG2	1:A:514:ILE:HG22	1.76	0.67
2:B:526:TYR:HA	2:B:544:LEU:HD21	1.76	0.66
2:B:421:MET:HE1	4:B:1000:EST:H161	1.75	0.66
1:A:413:ASN:HA	1:A:416:LYS:HE2	1.77	0.65
2:B:456:SER:HA	2:B:515:ARG:NH2	2.12	0.65
2:B:362:LYS:HE2	3:D:9:LEU:C	2.18	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:315:MET:CE	1:A:481:LYS:HG2	2.28	0.64
1:A:477:ARG:NH2	6:A:702:HOH:O	2.29	0.63
2:B:543:MET:CE	3:D:5:LEU:HD11	2.30	0.61
1:A:508:LEU:O	1:A:511[A]:LEU:HD22	2.01	0.61
2:B:315:MET:HE3	2:B:481:LYS:HD2	1.82	0.61
2:B:496:THR:O	2:B:500:GLN:HG3	2.01	0.60
1:A:368:VAL:HG22	6:A:706:HOH:O	2.02	0.59
2:B:362:LYS:CE	3:D:9:LEU:HD12	2.32	0.59
2:B:543:MET:HE1	3:D:5:LEU:HD11	1.85	0.59
2:B:362:LYS:CD	3:D:9:LEU:HD12	2.33	0.58
2:B:343:MET:CB	2:B:534:VAL:HG11	2.34	0.58
2:B:421:MET:HE3	4:B:1000:EST:H161	1.85	0.58
1:A:392:VAL:HG13	1:A:432:SER:HA	1.86	0.57
1:A:363:ARG:HG2	1:A:363:ARG:NH2	2.14	0.57
2:B:533:VAL:HG23	2:B:534:VAL:N	2.19	0.57
2:B:529:LYS:O	2:B:532:ASN:ND2	2.26	0.56
2:B:492:LYS:NZ	2:B:492:LYS:HB3	2.20	0.56
1:A:315:MET:HE1	1:A:481:LYS:HG2	1.87	0.55
2:B:308:LEU:HG	2:B:309:SER:H	1.71	0.55
1:A:342:MET:HG2	1:A:418:VAL:CG2	2.36	0.55
2:B:383:TRP:NE1	2:B:543:MET:HB3	2.23	0.54
2:B:311:THR:OG1	2:B:314:GLN:HG3	2.06	0.54
2:B:492:LYS:HB3	2:B:492:LYS:HZ3	1.72	0.52
2:B:423:GLU:O	2:B:427:MET:HE2	2.10	0.52
1:A:477:ARG:NH1	6:A:702:HOH:O	2.43	0.52
1:A:315:MET:HE3	1:A:481:LYS:HG2	1.92	0.51
1:A:342:MET:HG2	1:A:418:VAL:HG23	1.93	0.51
2:B:362:LYS:HE2	3:D:9:LEU:CA	2.41	0.51
2:B:343:MET:HB3	2:B:534:VAL:HG11	1.93	0.50
2:B:362:LYS:NZ	3:D:9:LEU:O	2.44	0.49
1:A:385:GLU:HG3	1:A:518:SER:HB2	1.94	0.49
1:A:497:LEU:HD21	2:B:497:LEU:CD1	2.42	0.49
2:B:492:LYS:CB	2:B:492:LYS:NZ	2.76	0.48
2:B:372:LEU:O	2:B:376:VAL:HG23	2.14	0.48
2:B:509:LEU:HD12	6:B:1145:HOH:O	2.14	0.48
1:A:316:VAL:HG21	1:A:489:LEU:HD21	1.96	0.47
2:B:308:LEU:HD23	2:B:310:LEU:H	1.80	0.47
1:A:412:ARG:HA	1:A:425:PHE:CD2	2.50	0.47
2:B:478:VAL:O	2:B:482:ILE:HG13	2.15	0.46
1:A:496:THR:O	1:A:500:GLN:HG3	2.15	0.46
3:D:8:MK8:HB1A	3:D:11:ASP:HB2	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:380:GLU:OE1	3:C:3:LYS:HE2	2.16	0.46
1:A:497:LEU:HD21	2:B:497:LEU:HD11	1.99	0.45
1:A:527:SER:O	1:A:531:LYS:HG3	2.17	0.45
3:D:1:ACE:O	3:D:2:HIS:CG	2.70	0.45
2:B:343:MET:HB2	2:B:534:VAL:HG11	2.00	0.44
2:B:440:LEU:HD12	2:B:444:GLU:HB2	2.00	0.43
1:A:422:VAL:HA	1:A:425:PHE:HB2	2.00	0.43
2:B:362:LYS:HE2	3:D:9:LEU:HA	2.00	0.43
2:B:497:LEU:HA	2:B:497:LEU:HD23	1.87	0.43
2:B:435:PHE:HE1	2:B:510:ILE:HG21	1.82	0.43
3:D:5:LEU:HD23	3:D:5:LEU:HA	1.81	0.43
2:B:533:VAL:CG2	2:B:534:VAL:N	2.81	0.43
2:B:440:LEU:CD1	2:B:444:GLU:HB2	2.48	0.43
2:B:327:LEU:HD12	2:B:353:GLU:HG3	2.01	0.42
2:B:392:VAL:HG13	2:B:432:SER:HA	2.02	0.42
1:A:509:LEU:HD23	1:A:509:LEU:HA	1.86	0.41
2:B:427:MET:HB3	2:B:517:MET:SD	2.60	0.41
1:A:477:ARG:CZ	6:A:702:HOH:O	2.68	0.41
1:A:392:VAL:HG13	1:A:432:SER:CA	2.50	0.41
1:A:456:SER:HA	1:A:515:ARG:NH2	2.36	0.41
2:B:308:LEU:CG	2:B:309:SER:H	2.34	0.41
1:A:511[A]:LEU:HA	1:A:514:ILE:HB	2.03	0.40
2:B:396:MET:HE2	2:B:436:ARG:HA	2.03	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:786:HOH:O	6:B:1121:HOH:O[1_554]	2.05	0.15
3:C:10:GLN:HE21	6:A:717:HOH:O[2_857]	1.52	0.08
1:A:359:ASN:ND2	2:B:328:TYR:OH[1_554]	2.14	0.06

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	232/261 (89%)	226 (97%)	6 (3%)	0	100	100
2	B	216/261 (83%)	212 (98%)	3 (1%)	1 (0%)	34	30
3	C	9/13 (69%)	9 (100%)	0	0	100	100
3	D	9/13 (69%)	6 (67%)	2 (22%)	1 (11%)	0	0
All	All	466/548 (85%)	453 (97%)	11 (2%)	2 (0%)	39	37

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	D	12	SER
2	B	535	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	205/236 (87%)	198 (97%)	7 (3%)	44	45
2	B	191/236 (81%)	175 (92%)	16 (8%)	14	9
3	C	9/9 (100%)	7 (78%)	2 (22%)	1	0
3	D	9/9 (100%)	6 (67%)	3 (33%)	0	0
All	All	414/490 (84%)	386 (93%)	28 (7%)	21	16

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	308	LEU
1	A	342	MET
1	A	425	PHE
1	A	460	THR
1	A	481	LYS
1	A	511[A]	LEU

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Mol	Chain	Res	Type
1	A	511[B]	LEU
2	B	308	LEU
2	B	310	LEU
2	B	321	ASP
2	B	339	GLU
2	B	371	THR
2	B	401	LYS
2	B	421	MET
2	B	425	PHE
2	B	437	MET
2	B	472	LYS
2	B	473	ASP
2	B	481	LYS
2	B	532	ASN
2	B	534	VAL
2	B	535	PRO
2	B	542	GLU
3	C	2	HIS
3	C	11	ASP
3	D	7	ARG
3	D	9	LEU
3	D	12	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	519	ASN
2	B	519	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 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	$\# Z > 2$
3	5GM	C	4	3	6,9,10	1.17	1 (16%)	6,12,14	0.98	0
3	MK8	C	8	3	6,8,9	1.19	0	3,10,12	0.83	0
3	5GM	D	4	3	6,9,10	1.37	2 (33%)	6,12,14	0.69	0
3	MK8	D	8	3	6,8,9	2.20	1 (16%)	3,10,12	0.74	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	5GM	C	4	3	-	0/8/10/13	0/0/0/0
3	MK8	C	8	3	-	0/6/8/11	0/0/0/0
3	5GM	D	4	3	-	0/8/10/13	0/0/0/0
3	MK8	D	8	3	-	0/6/8/11	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	8	MK8	CB-CA	-4.98	1.49	1.55
3	D	4	5GM	CB-CC	-2.42	1.46	1.54
3	C	4	5GM	CE-CD	-2.34	1.35	1.50
3	D	4	5GM	CE-CD	-2.09	1.37	1.50

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	8	MK8	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates 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$
4	EST	A	601	-	23,23,23	2.09	9 (39%)	36,36,36	1.20	4 (11%)
4	EST	B	1000	-	23,23,23	2.05	11 (47%)	36,36,36	1.15	3 (8%)

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
4	EST	A	601	-	-	0/0/40/40	0/4/4/4
4	EST	B	1000	-	-	0/0/40/40	0/4/4/4

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	601	EST	C10-C9	-4.45	1.45	1.52
4	A	601	EST	C1-C10	-2.90	1.35	1.39
4	B	1000	EST	C10-C9	-2.71	1.48	1.52
4	B	1000	EST	C1-C10	-2.64	1.36	1.39
4	B	1000	EST	C13-C17	-2.63	1.49	1.54
4	A	601	EST	O3-C3	-2.57	1.31	1.37
4	B	1000	EST	O17-C17	-2.49	1.39	1.43
4	B	1000	EST	C12-C13	-2.35	1.49	1.54
4	B	1000	EST	C11-C9	-2.33	1.50	1.53
4	B	1000	EST	C9-C8	-2.30	1.51	1.54
4	A	601	EST	C18-C13	-2.28	1.50	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	601	EST	C13-C14	-2.24	1.50	1.55
4	A	601	EST	C4-C5	-2.22	1.35	1.39
4	B	1000	EST	C13-C14	-2.21	1.50	1.55
4	A	601	EST	C11-C9	-2.17	1.51	1.53
4	B	1000	EST	C4-C5	-2.17	1.35	1.39
4	B	1000	EST	O3-C3	-2.10	1.32	1.37
4	A	601	EST	C9-C8	-2.05	1.52	1.54
4	A	601	EST	C5-C10	3.80	1.46	1.39
4	B	1000	EST	C5-C10	4.35	1.47	1.39

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	601	EST	C16-C17-C13	-2.60	102.53	104.58
4	B	1000	EST	C9-C8-C14	-2.37	105.23	108.91
4	A	601	EST	C12-C11-C9	-2.34	109.44	112.17
4	B	1000	EST	C12-C11-C9	-2.29	109.49	112.17
4	A	601	EST	C9-C8-C14	-2.13	105.60	108.91
4	B	1000	EST	C10-C9-C8	2.02	113.99	111.53
4	A	601	EST	C10-C9-C8	3.17	115.40	111.53

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
4	A	601	EST	1	0
4	B	1000	EST	3	0

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	231/261 (88%)	0.08	7 (3%) 54 62	16, 29, 49, 72	2 (0%)
2	B	224/261 (85%)	0.64	23 (10%) 9 12	18, 39, 69, 84	2 (0%)
3	C	9/13 (69%)	0.62	1 (11%) 7 10	24, 31, 66, 68	0
3	D	9/13 (69%)	2.35	3 (33%) 0 1	50, 63, 85, 108	0
All	All	473/548 (86%)	0.40	34 (7%) 18 25	16, 33, 66, 108	4 (0%)

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	12	SER	8.1
2	B	530	CYS	5.6
2	B	461	PHE	5.4
3	D	2	HIS	5.3
1	A	308	LEU	4.4
3	C	12	SER	3.8
2	B	448	LEU	3.8
2	B	460	THR	3.8
2	B	533	VAL	3.6
3	D	11	ASP	3.5
2	B	452	ILE	3.3
2	B	511	LEU	3.2
2	B	386	ILE	3.2
1	A	461	PHE	3.2
2	B	389	ILE	3.2
1	A	477	ARG	3.1
1	A	459	TYR	3.1
2	B	451	ILE	3.0
2	B	420	GLY	3.0
2	B	321	ASP	3.0
2	B	450	SER	3.0

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Mol	Chain	Res	Type	RSRZ
2	B	419	GLU	2.8
2	B	310	LEU	2.8
2	B	308	LEU	2.7
2	B	473	ASP	2.7
2	B	369	ASP	2.6
2	B	309	SER	2.6
2	B	526	TYR	2.5
1	A	473	ASP	2.5
1	A	386	ILE	2.4
2	B	459	TYR	2.4
2	B	387	LEU	2.3
2	B	370	LEU	2.2
1	A	387	LEU	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	5GM	C	4	10/11	0.97	0.09	-	25,37,57,57	0
3	MK8	C	8	9/10	0.96	0.08	-	21,25,31,34	0
3	MK8	D	8	9/10	0.93	0.13	-	46,61,66,68	0
3	5GM	D	4	10/11	0.91	0.13	-	49,65,70,70	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
4	EST	A	601	20/20	0.94	0.14	0.06	15,23,31,39	0
4	EST	B	1000	20/20	0.94	0.14	-0.08	23,35,40,41	0
5	CL	A	602	1/1	0.99	0.14	-	32,32,32,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.