



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:26 PM GMT

PDB ID : 1FM9
Title : THE 2.1 ANGSTROM RESOLUTION CRYSTAL STRUCTURE OF THE HETERODIMER OF THE HUMAN RXRALPHA AND PPARGAMMA LIGAND BINDING DOMAINS RESPECTIVELY BOUND WITH 9-CIS RETINOIC ACID AND GI262570 AND CO-ACTIVATOR PEPTIDES.
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Deposited on : 2000-08-16
Resolution : 2.10 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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

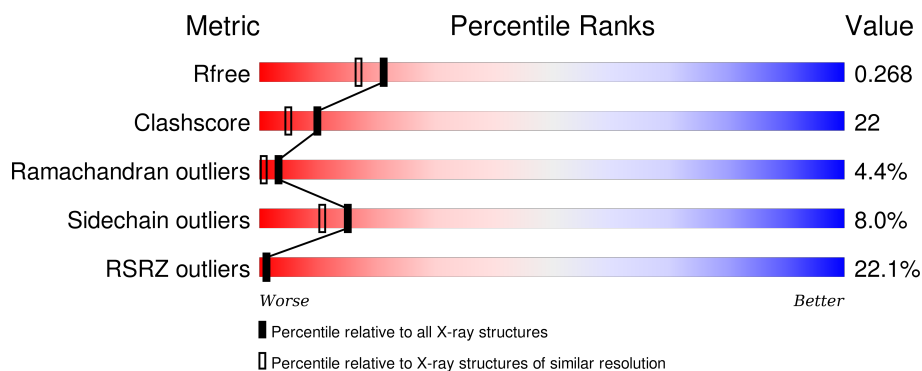
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	238	<div> <div>25%</div> <div>62%</div> <div>30%</div> <div>5%</div> <div>• •</div> </div>
2	D	272	<div> <div>16%</div> <div>67%</div> <div>28%</div> <div>5%</div> </div>
3	B	25	<div> <div>12%</div> <div>20%</div> <div>16%</div> <div>•</div> <div>60%</div> </div>
3	E	25	<div> <div>48%</div> <div>24%</div> <div>32%</div> <div>8%</div> <div>36%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4342 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 RETINOIC ACID RECEPTOR RXR-ALPHA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	232	Total	C	N	O	S	0	0	0
			1754	1117	306	321	10			

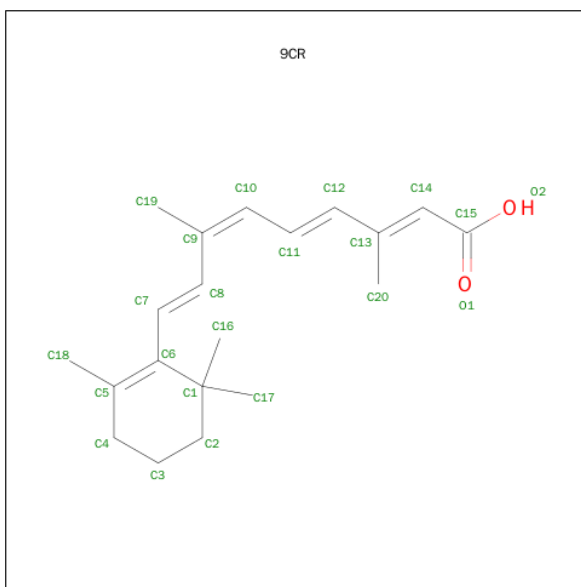
- Molecule 2 is a protein called PEROXISOME PROLIFERATOR ACTIVATED RECEPTOR GAMMA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	272	Total	C	N	O	S	0	0	0
			2179	1407	355	407	10			

- Molecule 3 is a protein called STEROID RECEPTOR COACTIVATOR.

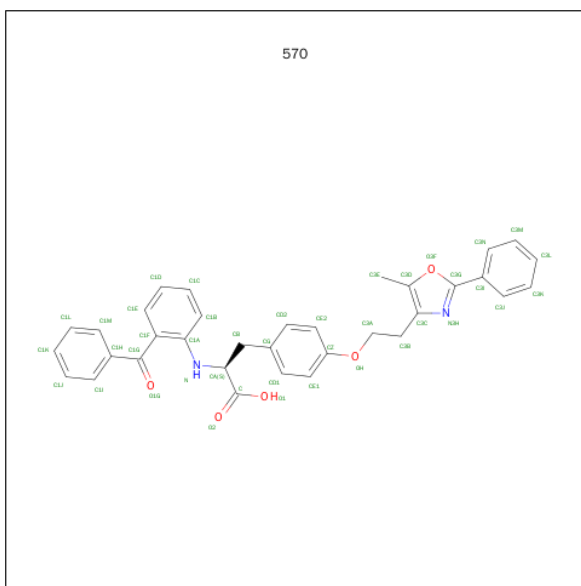
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	B	10	Total	C	N	O	0	0	0
			86	55	18	13			
3	E	16	Total	C	N	O	0	0	0
			129	79	27	23			

- Molecule 4 is (9CIS)-RETINOIC ACID (three-letter code: 9CR) (formula: C₂₀H₂₈O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			22	20	2		

- Molecule 5 is 2-(2-BENZOYL-PHENYLAMINO)-3-{4-[2-(5-METHYL-2-PHENYL-OXAZOL-4-YL)-ETHOXY]-PHENYL}-PROPIONIC ACID (three-letter code: 570) (formula: $C_{34}H_{30}N_2O_5$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	D	1	Total	C	N	O	0	0
			41	34	2	5		

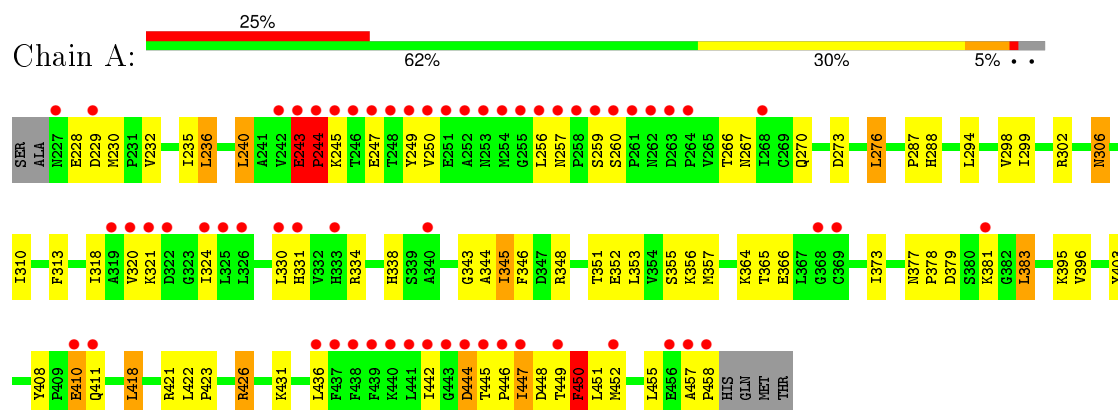
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	36	Total 36	O 36	0	0
6	D	89	Total 89	O 89	0	0
6	E	6	Total 6	O 6	0	0

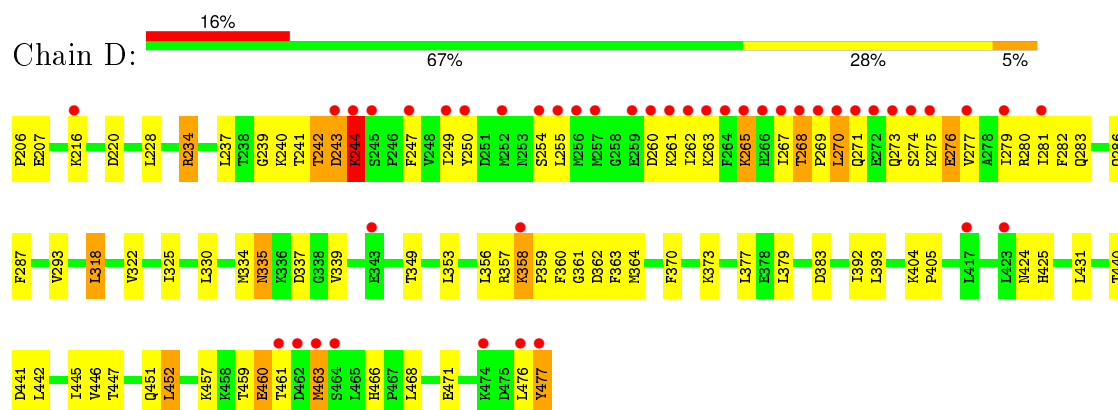
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

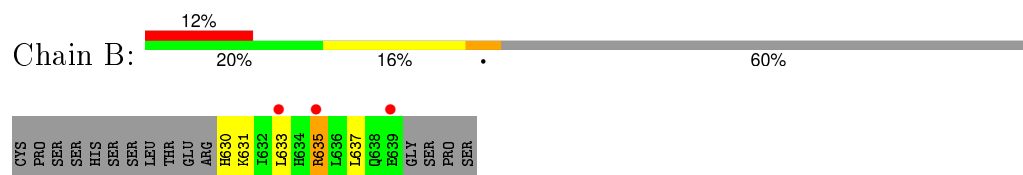
• Molecule 1: RETINOIC ACID RECEPTOR RXR-ALPHA



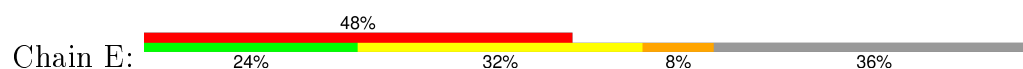
• Molecule 2: PEROXISOME PROLIFERATOR ACTIVATED RECEPTOR GAMMA



• Molecule 3: STEROID RECEPTOR COACTIVATOR



• Molecule 3: STEROID RECEPTOR COACTIVATOR



CYS	PRO	SER	SER	HIS	SER	SER	LEU	THR	E685	E686	H687	K688	I689	L690	H691	R692	L693	L694	Q695	E696	G697	S698	P699	S700
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	46.14Å 54.05Å 211.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.10 19.69 – 2.10	Depositor EDS
% Data completeness (in resolution range)	91.8 (20.00-2.10) 91.9 (19.69-2.10)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.14 (at 2.11Å)	Xtriage
Refinement program	CNS 0.5	Depositor
R, R_{free}	0.239 , 0.268 0.239 , 0.268	Depositor DCC
R_{free} test set	2895 reflections (9.95%)	DCC
Wilson B-factor (Å ²)	34.3	Xtriage
Anisotropy	0.254	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 57.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 29087 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4342	wwPDB-VP
Average B, all atoms (Å ²)	48.0	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, 0.375 respectively for untwinned datasets, and 0.333, 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: 9CR, 570

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	0.37	0/1789	0.64	2/2414 (0.1%)
2	D	0.42	0/2217	0.71	1/2987 (0.0%)
3	B	0.38	0/87	0.72	0/116
3	E	0.28	0/131	0.52	0/175
All	All	0.40	0/4224	0.68	3/5692 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	243	GLU	C-N-CD	-10.66	97.14	120.60
1	A	243	GLU	C-N-CA	6.33	148.58	122.00
2	D	206	PRO	N-CA-CB	5.39	109.77	103.30

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	1754	0	1733	80	0
2	D	2179	0	2237	99	0
3	B	86	0	86	7	0
3	E	129	0	125	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	22	0	27	4	0
5	D	41	0	29	5	0
6	A	36	0	0	3	0
6	D	89	0	0	2	0
6	E	6	0	0	1	0
All	All	4342	0	4237	187	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 22.

All (187) 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:268:THR:H	2:D:269:PRO:HD2	1.25	0.99
2:D:293:VAL:HG22	2:D:322:VAL:HG11	1.57	0.86
1:A:306:ASN:HD22	1:A:306:ASN:H	1.24	0.85
1:A:230:MET:CE	1:A:235:ILE:HD11	2.07	0.84
1:A:236:LEU:HD22	1:A:240:LEU:HD22	1.58	0.83
1:A:230:MET:HE3	1:A:235:ILE:HD11	1.60	0.83
1:A:338:HIS:HD2	1:A:343:GLY:HA3	1.46	0.81
2:D:242:THR:O	2:D:243:ASP:HB3	1.80	0.81
1:A:426:ARG:HH11	1:A:426:ARG:HB2	1.46	0.79
2:D:243:ASP:O	2:D:244:LYS:O	2.01	0.78
2:D:276:GLU:HG2	2:D:277:VAL:H	1.51	0.76
2:D:358:LYS:HB2	2:D:359:PRO:HD2	1.67	0.75
2:D:471:GLU:OE1	3:E:689:ILE:HG22	1.88	0.74
2:D:279:ILE:O	2:D:283:GLN:HG2	1.86	0.74
2:D:457:LYS:HA	2:D:463:MET:HG3	1.70	0.74
2:D:268:THR:N	2:D:269:PRO:HD2	2.00	0.73
2:D:268:THR:H	2:D:269:PRO:CD	2.00	0.73
2:D:357:ARG:HG3	2:D:357:ARG:HH11	1.54	0.72
2:D:357:ARG:NH1	2:D:358:LYS:CG	2.53	0.72
2:D:357:ARG:HH12	2:D:358:LYS:HB2	1.56	0.70
2:D:357:ARG:NH1	2:D:358:LYS:HB2	2.07	0.70
2:D:358:LYS:HB2	2:D:359:PRO:CD	2.22	0.69
2:D:466:HIS:HD2	2:D:468:LEU:H	1.40	0.69
2:D:335:ASN:HD22	2:D:335:ASN:C	1.96	0.68
1:A:306:ASN:H	1:A:306:ASN:ND2	1.92	0.66
1:A:426:ARG:NH1	1:A:426:ARG:HB2	2.09	0.66
2:D:383:ASP:OD2	2:D:425:HIS:HE1	1.77	0.66
2:D:358:LYS:CB	2:D:359:PRO:CD	2.73	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:325:ILE:HD11	2:D:392:ILE:HG13	1.79	0.65
1:A:338:HIS:CD2	1:A:343:GLY:HA3	2.29	0.65
2:D:243:ASP:OD1	2:D:243:ASP:O	2.15	0.65
2:D:286:GLN:OE1	5:D:200:570:H1C	1.98	0.64
2:D:273:GLN:HA	2:D:280:ARG:HD3	1.78	0.64
2:D:275:LYS:O	2:D:276:GLU:O	2.15	0.64
2:D:293:VAL:HG22	2:D:322:VAL:CG1	2.28	0.64
2:D:466:HIS:CD2	2:D:468:LEU:H	2.16	0.63
1:A:366:GLU:HG2	1:A:418:LEU:HD21	1.79	0.62
3:E:685:GLU:HB2	3:E:687:HIS:CE1	2.33	0.62
2:D:262:ILE:HA	2:D:265:LYS:HD2	1.80	0.62
1:A:320:VAL:HG21	1:A:331:HIS:CE1	2.34	0.62
1:A:343:GLY:O	1:A:344:ALA:HB3	2.00	0.60
2:D:377:LEU:HD22	2:D:431:LEU:HD12	1.82	0.60
2:D:357:ARG:HG3	2:D:357:ARG:NH1	2.16	0.60
2:D:357:ARG:NH1	2:D:358:LYS:HG2	2.16	0.59
2:D:446:VAL:HG11	2:D:477:TYR:HE1	1.68	0.59
1:A:447:ILE:HG12	1:A:452:MET:CE	2.33	0.59
2:D:262:ILE:HA	2:D:265:LYS:CD	2.32	0.59
2:D:247:PHE:HE2	2:D:254:SER:O	1.86	0.58
3:E:686:ARG:HH12	3:E:691:HIS:CD2	2.22	0.58
1:A:243:GLU:HB2	1:A:244:PRO:O	2.02	0.58
2:D:357:ARG:NH1	2:D:358:LYS:CB	2.67	0.57
1:A:320:VAL:HG21	1:A:331:HIS:NE2	2.20	0.57
2:D:276:GLU:HG2	2:D:277:VAL:N	2.20	0.57
1:A:423:PRO:HG3	2:D:440:THR:HG22	1.87	0.57
3:E:685:GLU:HG2	3:E:686:ARG:H	1.68	0.56
2:D:239:GLY:O	2:D:240:LYS:HD3	2.04	0.56
1:A:228:GLU:O	1:A:229:ASP:HB2	2.05	0.56
1:A:320:VAL:HG22	1:A:321:LYS:N	2.21	0.56
4:A:201:9CR:H19	4:A:201:9CR:H8	1.88	0.56
1:A:298:VAL:HG13	3:B:633:LEU:HD12	1.88	0.56
1:A:436:LEU:HD22	4:A:201:9CR:H16	1.88	0.55
2:D:270:LEU:HD22	2:D:270:LEU:O	2.06	0.55
1:A:306:ASN:ND2	6:A:10:HOH:O	2.40	0.54
2:D:281:ILE:HD12	2:D:356:LEU:HD11	1.88	0.54
2:D:377:LEU:HD22	2:D:431:LEU:CD1	2.38	0.54
2:D:325:ILE:HD11	2:D:392:ILE:CG1	2.38	0.54
1:A:298:VAL:HA	3:B:633:LEU:HD11	1.89	0.54
1:A:451:LEU:O	1:A:455:LEU:HD13	2.08	0.53
1:A:447:ILE:HD11	1:A:452:MET:SD	2.50	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408:TYR:HA	1:A:410:GLU:OE2	2.08	0.52
2:D:270:LEU:HD21	6:D:90:HOH:O	2.09	0.52
1:A:243:GLU:OE1	1:A:364:LYS:NZ	2.43	0.51
2:D:335:ASN:ND2	2:D:337:ASP:H	2.08	0.51
2:D:260:ASP:HA	2:D:263:LYS:HG3	1.93	0.51
1:A:345:ILE:HG12	1:A:431:LYS:CD	2.40	0.51
2:D:275:LYS:O	2:D:279:ILE:HG22	2.11	0.51
2:D:357:ARG:HH11	2:D:358:LYS:HG2	1.74	0.51
1:A:423:PRO:CG	2:D:440:THR:HG22	2.41	0.51
2:D:249:ILE:HD13	2:D:255:LEU:HD12	1.92	0.51
1:A:313:PHE:CE1	1:A:324:ILE:HD13	2.46	0.51
1:A:273:ASP:OD1	1:A:449:THR:O	2.29	0.51
1:A:318:ILE:HD11	1:A:357:MET:HB3	1.93	0.50
1:A:351:THR:HA	1:A:355:SER:OG	2.10	0.50
1:A:267:ASN:HB3	1:A:330:LEU:HD13	1.93	0.50
2:D:261:LYS:O	2:D:265:LYS:HG3	2.12	0.50
1:A:310:ILE:HA	1:A:313:PHE:CE2	2.46	0.50
1:A:423:PRO:CD	2:D:440:THR:HG22	2.42	0.49
1:A:266:THR:O	1:A:270:GLN:HG3	2.12	0.49
2:D:334:MET:HG2	2:D:339:VAL:HB	1.94	0.49
1:A:423:PRO:N	2:D:440:THR:HG22	2.27	0.49
3:E:685:GLU:OE1	3:E:685:GLU:N	2.45	0.49
1:A:298:VAL:HG22	3:B:633:LEU:CD1	2.43	0.49
1:A:267:ASN:HD22	1:A:330:LEU:HD11	1.78	0.49
2:D:477:TYR:C	2:D:477:TYR:CD1	2.83	0.49
2:D:356:LEU:HB2	2:D:361:GLY:HA2	1.93	0.49
1:A:365:THR:HG21	1:A:403:TYR:CE2	2.48	0.48
2:D:447:THR:HG22	2:D:477:TYR:OXT	2.12	0.48
1:A:228:GLU:O	1:A:228:GLU:HG3	2.12	0.48
2:D:370:PHE:HB2	2:D:445:ILE:HD11	1.95	0.48
2:D:377:LEU:HB2	2:D:379:LEU:HD12	1.95	0.48
1:A:243:GLU:HB3	1:A:244:PRO:C	2.34	0.48
1:A:449:THR:O	1:A:450:PHE:CB	2.59	0.48
1:A:426:ARG:CB	1:A:426:ARG:HH11	2.19	0.48
2:D:275:LYS:O	2:D:279:ILE:CG2	2.61	0.48
2:D:262:ILE:HA	2:D:265:LYS:CG	2.44	0.48
2:D:260:ASP:O	2:D:263:LYS:HB2	2.13	0.48
2:D:270:LEU:C	2:D:270:LEU:HD22	2.35	0.47
1:A:230:MET:CE	1:A:396:VAL:HG22	2.44	0.47
2:D:468:LEU:CD1	3:E:689:ILE:HG23	2.44	0.47
2:D:468:LEU:HD13	3:E:689:ILE:CG2	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:243:ASP:CG	2:D:243:ASP:O	2.52	0.47
2:D:273:GLN:HA	2:D:280:ARG:CD	2.45	0.47
3:E:695:GLN:CA	3:E:695:GLN:HE21	2.28	0.47
1:A:306:ASN:N	1:A:306:ASN:ND2	2.63	0.47
2:D:466:HIS:HD2	2:D:468:LEU:N	2.10	0.47
1:A:444:ASP:O	1:A:445:THR:OG1	2.33	0.47
2:D:404:LYS:HB3	2:D:405:PRO:HD3	1.97	0.47
2:D:255:LEU:HD23	2:D:277:VAL:HG23	1.95	0.47
3:E:686:ARG:NE	3:E:686:ARG:HA	2.29	0.47
1:A:273:ASP:CG	1:A:448:ASP:HB2	2.36	0.46
1:A:381:LYS:HG3	1:A:381:LYS:H	1.45	0.46
2:D:447:THR:O	2:D:451:GLN:HG3	2.16	0.46
2:D:234:ARG:NH2	2:D:334:MET:O	2.46	0.46
6:A:23:HOH:O	2:D:440:THR:HG23	2.16	0.46
1:A:299:ILE:HG21	1:A:383:LEU:HD13	1.98	0.46
2:D:282:PHE:CZ	5:D:200:570:H1D	2.51	0.46
1:A:330:LEU:HD23	1:A:331:HIS:N	2.31	0.45
1:A:436:LEU:CD2	4:A:201:9CR:H16	2.46	0.45
1:A:373:ILE:HG13	1:A:396:VAL:HG11	1.99	0.45
2:D:287:PHE:C	2:D:287:PHE:CD1	2.90	0.45
4:A:201:9CR:H8	4:A:201:9CR:H10	1.80	0.45
1:A:243:GLU:CB	1:A:244:PRO:C	2.85	0.45
3:B:633:LEU:C	3:B:633:LEU:HD13	2.37	0.45
1:A:421:ARG:NH2	6:A:25:HOH:O	2.50	0.45
1:A:230:MET:SD	1:A:287:PRO:HG2	2.56	0.44
1:A:228:GLU:O	1:A:229:ASP:CB	2.65	0.44
2:D:335:ASN:ND2	2:D:335:ASN:C	2.68	0.44
2:D:250:TYR:N	2:D:250:TYR:CD1	2.85	0.44
1:A:422:LEU:CB	2:D:440:THR:HG21	2.47	0.44
3:B:631:LYS:O	3:B:635:ARG:HB2	2.18	0.44
1:A:229:ASP:O	1:A:288:HIS:NE2	2.51	0.44
3:B:633:LEU:HD13	3:B:637:LEU:HD12	1.99	0.44
2:D:273:GLN:HG2	2:D:274:SER:H	1.83	0.44
2:D:216:LYS:NZ	2:D:220:ASP:OD1	2.50	0.44
1:A:353:LEU:O	1:A:357:MET:HG3	2.18	0.43
2:D:353:LEU:HD13	2:D:364:MET:HG3	2.01	0.43
1:A:230:MET:HE2	1:A:396:VAL:HG22	1.99	0.43
2:D:277:VAL:HG22	2:D:277:VAL:O	2.18	0.43
3:E:688:LYS:O	3:E:691:HIS:HB2	2.18	0.43
2:D:468:LEU:HD13	3:E:689:ILE:HG23	2.01	0.43
1:A:294:LEU:C	1:A:294:LEU:HD23	2.38	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:348:ARG:O	1:A:352:GLU:HB2	2.19	0.43
1:A:344:ALA:O	1:A:348:ARG:HG3	2.19	0.42
2:D:286:GLN:HB2	5:D:200:570:C1C	2.49	0.42
2:D:273:GLN:HG2	2:D:274:SER:N	2.33	0.42
1:A:356:LYS:HG3	1:A:421:ARG:NH1	2.34	0.42
2:D:249:ILE:HD13	2:D:255:LEU:CD1	2.49	0.42
2:D:360:PHE:HD1	5:D:200:570:H1L	1.84	0.42
2:D:459:THR:C	2:D:460:GLU:HG2	2.39	0.42
2:D:476:LEU:CG	2:D:477:TYR:H	2.32	0.42
3:E:698:SER:HA	3:E:699:PRO:HD3	1.92	0.42
1:A:320:VAL:CG2	1:A:321:LYS:N	2.83	0.42
1:A:276:LEU:HB3	1:A:450:PHE:CE1	2.54	0.42
1:A:267:ASN:HD22	1:A:330:LEU:CD1	2.32	0.42
2:D:468:LEU:CD1	3:E:689:ILE:CG2	2.98	0.42
1:A:276:LEU:HA	1:A:276:LEU:HD12	1.85	0.41
1:A:379:ASP:OD1	2:D:373:LYS:HE2	2.20	0.41
1:A:229:ASP:HB3	1:A:395:LYS:HD3	2.01	0.41
1:A:313:PHE:C	1:A:313:PHE:CD1	2.93	0.41
3:E:695:GLN:OE1	6:E:65:HOH:O	2.22	0.41
2:D:318:LEU:HD21	3:E:694:LEU:HD21	2.02	0.41
1:A:348:ARG:NH1	1:A:431:LYS:HE3	2.35	0.41
1:A:302:ARG:HD2	3:B:630:HIS:CD2	2.55	0.41
1:A:345:ILE:CG2	1:A:346:PHE:N	2.84	0.41
2:D:330:LEU:O	2:D:334:MET:HG3	2.20	0.41
2:D:392:ILE:HG22	2:D:393:LEU:HD22	2.03	0.41
2:D:377:LEU:HB2	2:D:379:LEU:CD1	2.51	0.41
2:D:476:LEU:HG	2:D:477:TYR:H	1.86	0.41
1:A:447:ILE:HG12	1:A:452:MET:HE1	2.01	0.41
1:A:247:GLU:C	1:A:249:TYR:H	2.23	0.41
2:D:441:ASP:OD2	6:D:66:HOH:O	2.21	0.41
1:A:366:GLU:CG	1:A:418:LEU:HD21	2.49	0.41
2:D:360:PHE:CD1	5:D:200:570:H1L	2.57	0.40
1:A:447:ILE:HD12	1:A:451:LEU:HG	2.02	0.40
1:A:377:ASN:HA	1:A:378:PRO:HD3	1.89	0.40
2:D:363:PHE:CZ	2:D:452:LEU:HB3	2.56	0.40
2:D:446:VAL:HG11	2:D:477:TYR:CE1	2.52	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	230/238 (97%)	196 (85%)	20 (9%)	14 (6%)	2	0
2	D	270/272 (99%)	255 (94%)	7 (3%)	8 (3%)	5	2
3	B	8/25 (32%)	7 (88%)	1 (12%)	0	100	100
3	E	14/25 (56%)	10 (71%)	3 (21%)	1 (7%)	1	0
All	All	522/560 (93%)	468 (90%)	31 (6%)	23 (4%)	3	1

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	243	GLU
1	A	244	PRO
1	A	257	ASN
1	A	259	SER
1	A	260	SER
1	A	446	PRO
2	D	243	ASP
2	D	244	LYS
2	D	276	GLU
2	D	358	LYS
1	A	245	LYS
1	A	256	LEU
1	A	442	ILE
1	A	450	PHE
2	D	267	ILE
1	A	250	VAL
2	D	265	LYS
2	D	271	GLN
1	A	457	ALA
2	D	268	THR
1	A	444	ASP
3	E	699	PRO
1	A	447	ILE

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	182/205 (89%)	167 (92%)	15 (8%)	14	10
2	D	243/245 (99%)	224 (92%)	19 (8%)	16	11
3	B	9/24 (38%)	8 (89%)	1 (11%)	8	4
3	E	14/24 (58%)	13 (93%)	1 (7%)	18	14
All	All	448/498 (90%)	412 (92%)	36 (8%)	15	11

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	232	VAL
1	A	236	LEU
1	A	240	LEU
1	A	244	PRO
1	A	276	LEU
1	A	306	ASN
1	A	334	ARG
1	A	345	ILE
1	A	383	LEU
1	A	410	GLU
1	A	411	GLN
1	A	418	LEU
1	A	426	ARG
1	A	450	PHE
1	A	458	PRO
2	D	207	GLU
2	D	228	LEU
2	D	234	ARG
2	D	237	LEU
2	D	241	THR
2	D	242	THR
2	D	244	LYS
2	D	270	LEU
2	D	318	LEU

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Mol	Chain	Res	Type
2	D	335	ASN
2	D	349	THR
2	D	362	ASP
2	D	424	ASN
2	D	442	LEU
2	D	452	LEU
2	D	460	GLU
2	D	461	THR
2	D	463	MET
2	D	477	TYR
3	B	635	ARG
3	E	695	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	267	ASN
1	A	270	GLN
1	A	306	ASN
1	A	338	HIS
2	D	273	GLN
2	D	308	ASN
2	D	335	ASN
2	D	375	ASN
2	D	424	ASN
2	D	425	HIS
2	D	454	GLN
2	D	466	HIS
3	B	634	HIS
3	B	638	GLN
3	E	691	HIS
3	E	695	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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
4	9CR	A	201	-	19,22,22	3.81	9 (47%)	26,30,30	2.55	11 (42%)
5	570	D	200	-	38,45,45	2.39	18 (47%)	48,61,61	1.77	4 (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	9CR	A	201	-	-	0/13/32/32	0/1/1/1
5	570	D	200	-	-	0/23/30/30	0/4/5/5

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	200	570	CB-CA	2.15	1.56	1.53
5	D	200	570	C3J-C3I	2.16	1.43	1.39
5	D	200	570	OH-CZ	2.33	1.43	1.37
5	D	200	570	CD1-CG	2.35	1.43	1.38
5	D	200	570	CD2-CG	2.37	1.43	1.38
5	D	200	570	CE1-CZ	2.45	1.43	1.38
5	D	200	570	C1M-C1H	2.45	1.43	1.39
5	D	200	570	C3G-N3H	2.49	1.41	1.34
5	D	200	570	C3N-C3I	2.57	1.44	1.39
4	A	201	9CR	C2-C3	2.64	1.59	1.52
4	A	201	9CR	C14-C13	2.74	1.39	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	200	570	C1B-C1A	2.81	1.44	1.39
5	D	200	570	C1E-C1F	2.94	1.44	1.39
4	A	201	9CR	C4-C5	2.95	1.57	1.51
4	A	201	9CR	C2-C1	3.11	1.61	1.54
4	A	201	9CR	C20-C13	3.28	1.57	1.50
5	D	200	570	C1I-C1H	3.37	1.45	1.39
5	D	200	570	CA-N	3.53	1.51	1.45
5	D	200	570	C3E-C3D	3.64	1.53	1.48
5	D	200	570	C1F-C1A	3.65	1.47	1.41
5	D	200	570	C1A-N	3.96	1.46	1.37
4	A	201	9CR	C10-C9	4.01	1.41	1.35
5	D	200	570	CE2-CZ	4.25	1.47	1.38
4	A	201	9CR	C16-C1	4.32	1.63	1.53
5	D	200	570	C1F-C1G	4.49	1.59	1.50
4	A	201	9CR	C1-C6	9.17	1.66	1.53
4	A	201	9CR	C5-C6	9.87	1.50	1.34

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	201	9CR	C16-C1-C6	-4.91	102.61	110.30
5	D	200	570	O1G-C1G-C1H	-4.61	112.75	120.12
4	A	201	9CR	C19-C9-C10	-4.14	116.79	122.90
4	A	201	9CR	C11-C10-C9	-3.51	122.12	127.20
4	A	201	9CR	C1-C6-C5	-2.83	118.51	122.66
4	A	201	9CR	C7-C6-C5	2.31	126.66	121.37
4	A	201	9CR	C8-C7-C6	2.60	135.13	127.32
4	A	201	9CR	C2-C1-C6	3.32	115.62	110.36
4	A	201	9CR	C20-C13-C12	3.38	123.72	118.10
4	A	201	9CR	C7-C8-C9	3.90	132.16	126.22
5	D	200	570	C1A-N-CA	3.94	133.24	124.75
4	A	201	9CR	C19-C9-C8	3.99	124.73	118.10
4	A	201	9CR	C17-C1-C6	5.22	118.49	110.30
5	D	200	570	C1A-C1F-C1G	6.37	126.47	121.92
5	D	200	570	C1F-C1G-C1H	6.41	130.39	119.47

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	201	9CR	4	0
5	D	200	570	5	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	232/238 (97%)	2.02	59 (25%) 1 1	23, 42, 109, 124	0
2	D	272/272 (100%)	1.42	43 (15%) 3 4	26, 39, 115, 131	0
3	B	10/25 (40%)	1.71	3 (30%) 1 1	50, 57, 69, 76	0
3	E	16/25 (64%)	3.84	12 (75%) 0 0	60, 68, 96, 98	0
All	All	530/560 (94%)	1.76	117 (22%) 1 1	23, 41, 110, 131	0

All (117) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	267	ILE	33.0
1	A	252	ALA	18.8
1	A	256	LEU	18.7
2	D	274	SER	18.6
2	D	266	HIS	17.2
1	A	249	TYR	14.4
2	D	264	PHE	14.4
2	D	270	LEU	14.2
1	A	441	LEU	13.6
2	D	265	LYS	13.3
1	A	259	SER	13.2
1	A	442	ILE	13.1
2	D	262	ILE	13.0
1	A	444	ASP	12.9
2	D	268	THR	12.8
1	A	445	THR	12.4
1	A	258	PRO	12.3
1	A	253	ASN	11.6
1	A	250	VAL	11.6
2	D	272	GLU	11.5
3	E	700	SER	11.5

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Mol	Chain	Res	Type	RSRZ
1	A	257	ASN	11.4
1	A	260	SER	11.3
2	D	269	PRO	11.2
1	A	446	PRO	11.0
1	A	227	ASN	10.8
1	A	262	ASN	10.8
1	A	248	THR	9.9
1	A	443	GLY	9.8
1	A	261	PRO	9.7
1	A	245	LYS	9.7
2	D	271	GLN	9.4
3	E	699	PRO	9.2
1	A	255	GLY	9.1
2	D	263	LYS	9.1
1	A	458	PRO	9.1
1	A	457	ALA	8.6
1	A	247	GLU	8.4
3	E	698	SER	8.3
1	A	254	MET	7.7
2	D	260	ASP	7.3
2	D	275	LYS	7.1
2	D	244	LYS	7.1
2	D	257	MET	7.0
3	E	685	GLU	6.2
1	A	251	GLU	6.1
2	D	273	GLN	5.9
1	A	244	PRO	5.8
2	D	461	THR	5.4
2	D	477	TYR	5.4
1	A	246	THR	5.1
2	D	261	LYS	4.9
1	A	456	GLU	4.6
1	A	447	ILE	4.3
2	D	259	GLU	4.0
3	E	696	GLU	3.9
2	D	463	MET	3.8
2	D	358	LYS	3.7
3	E	686	ARG	3.7
2	D	462	ASP	3.5
3	E	697	GLY	3.5
1	A	319	ALA	3.4
1	A	437	PHE	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	229	ASP	3.4
3	E	690	LEU	3.3
2	D	256	MET	3.3
2	D	252	MET	3.3
2	D	250	TYR	3.3
1	A	452	MET	3.2
3	B	633	LEU	3.2
1	A	340	ALA	3.2
1	A	264	PRO	3.1
1	A	438	PHE	3.1
1	A	322	ASP	3.1
2	D	243	ASP	3.0
2	D	343	GLU	3.0
2	D	281	ILE	3.0
2	D	277	VAL	3.0
2	D	255	LEU	2.9
2	D	245	SER	2.8
1	A	321	LYS	2.8
1	A	440	LYS	2.8
1	A	324	ILE	2.8
3	B	635	ARG	2.7
1	A	436	LEU	2.7
2	D	464	SER	2.7
1	A	439	PHE	2.7
2	D	247	PHE	2.7
3	B	639	GLU	2.6
1	A	381	LYS	2.6
1	A	242	VAL	2.5
1	A	325	LEU	2.5
1	A	410	GLU	2.5
3	E	695	GLN	2.5
2	D	279	ILE	2.5
1	A	331	HIS	2.4
3	E	694	LEU	2.4
1	A	263	ASP	2.3
1	A	333	HIS	2.3
3	E	692	ARG	2.2
2	D	249	ILE	2.2
3	E	693	LEU	2.2
2	D	423	LEU	2.2
1	A	411	GLN	2.2
2	D	417	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	330	LEU	2.1
1	A	369	CYS	2.1
1	A	268	ILE	2.1
2	D	216	LYS	2.1
2	D	474	LYS	2.1
2	D	254	SER	2.1
2	D	476	LEU	2.1
1	A	320	VAL	2.1
1	A	243	GLU	2.0
1	A	368	GLY	2.0
1	A	326	LEU	2.0
1	A	449	THR	2.0

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 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	9CR	A	201	22/22	0.88	0.25	1.80	43,48,50,50	0
5	570	D	200	41/41	0.93	0.17	0.30	33,38,44,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.