



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2018 – 06:59 am GMT

PDB ID : 1E7E
Title : HUMAN SERUM ALBUMIN COMPLEXED WITH DECANOIC ACID
(CAPRIC ACID)
Authors : Bhattacharya, A.A.; Gruene, T.; Curry, S.
Deposited on : 2000-08-29
Resolution : 2.50 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : trunk30967
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk30967

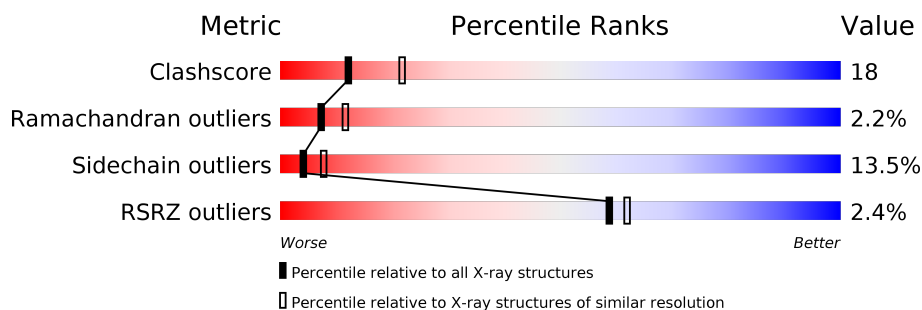
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	122126	4827 (2.50-2.50)
Ramachandran outliers	120053	4735 (2.50-2.50)
Sidechain outliers	120020	4737 (2.50-2.50)
RSRZ outliers	108989	4058 (2.50-2.50)

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	585	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	DKA	A	1007	-	-	-	X

2 Entry composition [i](#)

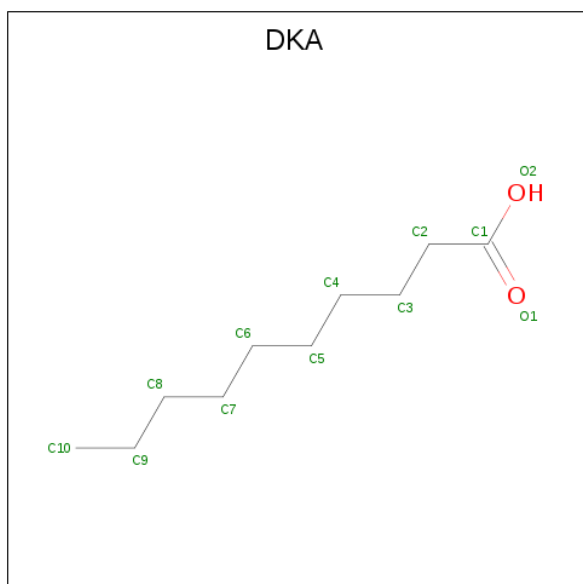
There are 3 unique types of molecules in this entry. The entry contains 4591 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SERUM ALBUMIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	582	Total	C	N	O	S	0	0	0
			4455	2821	747	846	41			

- Molecule 2 is DECANOIC ACID (three-letter code: DKA) (formula: $C_{10}H_{20}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			11	9	2		
2	A	1	Total	C	O	0	0
			10	8	2		
2	A	1	Total	C	O	0	0
			9	7	2		
2	A	1	Total	C	O	0	0
			12	10	2		
2	A	1	Total	C	O	0	0
			11	9	2		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 11 9 2	0	0
2	A	1	Total C 10 10	0	0
2	A	1	Total C O 12 10 2	0	0
2	A	1	Total C O 12 10 2	0	0
2	A	1	Total C 8 8	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	30	Total O 30 30	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	186.78Å 39.19Å 95.34Å 90.00° 105.15° 90.00°	Depositor
Resolution (Å)	40.00 – 2.50 37.46 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.1 (40.00-2.50) 97.1 (37.46-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.61 (at 2.51Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.220 , 0.271 0.212 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	51.2	Xtriage
Anisotropy	0.323	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 74.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4591	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.38% 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: DKA

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	0/4542	0.56	1/6157 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	94	GLN	N-CA-C	-5.64	95.76	111.00

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	4455	0	4228	164	0
2	A	106	0	152	10	0
3	A	30	0	0	2	0
All	All	4591	0	4380	164	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 18.

All (164) 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:548:MET:HG2	2:A:1005:DKA:H41	1.50	0.92
1:A:479:GLU:HG3	1:A:480:SER:N	1.89	0.87
1:A:479:GLU:OE2	1:A:487:CYS:HB2	1.75	0.86
1:A:59:ALA:HB3	1:A:62:CYS:SG	2.18	0.82
1:A:267:ASN:HD22	1:A:267:ASN:N	1.79	0.81
1:A:19:PHE:O	1:A:23:VAL:HG23	1.83	0.78
1:A:151:ALA:HB3	1:A:152:PRO:HD3	1.69	0.74
1:A:114:ARG:NH2	1:A:116:VAL:HG12	2.02	0.73
1:A:118:PRO:HG2	1:A:123:MET:HG3	1.70	0.73
1:A:49:PHE:O	1:A:52:THR:HB	1.91	0.70
1:A:29:GLN:HG2	1:A:147:PRO:HA	1.72	0.69
1:A:521:ARG:O	1:A:525:LYS:HG3	1.92	0.69
1:A:138:TYR:O	1:A:142:ILE:HG12	1.93	0.68
1:A:81:ARG:O	1:A:85:GLY:HA2	1.92	0.68
1:A:138:TYR:HB3	2:A:1001:DKA:H72	1.74	0.67
1:A:485:ARG:HB3	1:A:486:PRO:HD3	1.77	0.66
1:A:424:VAL:O	1:A:428:ARG:HG3	1.95	0.65
1:A:475:LYS:O	1:A:479:GLU:HB3	1.97	0.65
1:A:366:PRO:C	1:A:368:GLU:H	2.01	0.64
1:A:9:HIS:O	1:A:13:ASP:HB2	1.99	0.63
1:A:525:LYS:HD3	2:A:1005:DKA:H22	1.81	0.63
1:A:95:GLU:O	1:A:96:PRO:C	2.32	0.63
1:A:385:GLN:HG3	1:A:446:MET:HE1	1.82	0.62
1:A:510:HIS:O	1:A:513:ILE:HG23	1.99	0.61
1:A:267:ASN:ND2	1:A:267:ASN:N	2.49	0.61
1:A:208:GLU:HG2	1:A:212:LYS:HD2	1.83	0.61
1:A:77:VAL:O	1:A:80:LEU:HB2	2.00	0.61
1:A:499:PRO:HB3	1:A:535:HIS:O	2.00	0.60
1:A:265:CYS:O	1:A:268:GLN:HG3	2.00	0.60
1:A:16:GLU:O	1:A:20:LYS:HG3	2.01	0.59
1:A:107:ASP:O	1:A:147:PRO:HG3	2.02	0.59
1:A:133:THR:O	1:A:137:LYS:HB2	2.02	0.59
1:A:161:TYR:CD1	1:A:185:LEU:HD23	2.37	0.59
1:A:567:CYS:O	1:A:571:GLU:HB2	2.04	0.58
1:A:344:VAL:HG22	1:A:450:GLU:OE2	2.04	0.58
1:A:323:LYS:HE3	3:A:2026:HOH:O	2.04	0.58
1:A:446:MET:HB3	1:A:447:PRO:HD3	1.85	0.58
1:A:177:CYS:O	1:A:180:PRO:HD2	2.04	0.58
1:A:233:LYS:HE2	1:A:237:ASP:OD2	2.03	0.57
1:A:420:THR:N	1:A:421:PRO:HD2	2.18	0.57
1:A:86:GLU:O	1:A:89:ASP:HB2	2.04	0.57
1:A:513:ILE:HA	1:A:516:LEU:HD12	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:317:LYS:HG3	1:A:318:ASN:N	2.19	0.56
1:A:430:LEU:HD12	1:A:456:VAL:HG11	1.87	0.56
1:A:420:THR:O	1:A:424:VAL:HG23	2.04	0.56
1:A:420:THR:HG23	1:A:530:VAL:HG11	1.87	0.56
1:A:284:LEU:HD12	2:A:1002:DKA:H42	1.87	0.56
1:A:245:CYS:HA	1:A:253:CYS:HB2	1.88	0.56
1:A:540:THR:O	1:A:544:LEU:HB2	2.06	0.55
1:A:95:GLU:OE2	1:A:99:ASN:HB2	2.07	0.55
1:A:241:VAL:HG22	1:A:256:ASP:HB3	1.89	0.55
1:A:306:ALA:HA	1:A:310:VAL:HG23	1.89	0.55
1:A:114:ARG:HH21	1:A:116:VAL:HA	1.70	0.55
1:A:378:LYS:O	1:A:382:GLU:HG3	2.07	0.55
1:A:79:THR:HB	1:A:83:THR:OG1	2.07	0.55
1:A:488:PHE:HB3	2:A:1004:DKA:H52	1.89	0.54
1:A:511:ALA:O	1:A:514:CYS:SG	2.65	0.54
1:A:549:ASP:O	1:A:552:ALA:HB3	2.08	0.54
1:A:299:PRO:HB2	1:A:302:LEU:CD2	2.38	0.54
1:A:360:CYS:O	1:A:366:PRO:HG3	2.07	0.53
1:A:5:SER:HA	1:A:62:CYS:O	2.08	0.53
1:A:115:LEU:HD13	1:A:145:ARG:CZ	2.39	0.52
1:A:73:LYS:O	1:A:76:THR:HB	2.09	0.52
1:A:480:SER:HB2	2:A:1006:DKA:O2	2.10	0.52
1:A:510:HIS:O	1:A:512:ASP:N	2.43	0.52
1:A:36:PHE:CE1	1:A:137:LYS:HA	2.45	0.52
1:A:81:ARG:HG3	1:A:81:ARG:O	2.09	0.52
1:A:21:ALA:O	1:A:25:ILE:HG13	2.10	0.52
1:A:366:PRO:O	1:A:368:GLU:N	2.43	0.51
1:A:430:LEU:CD1	1:A:456:VAL:HG11	2.40	0.51
1:A:392:CYS:O	1:A:396:GLU:HG2	2.10	0.51
1:A:220:SER:HB2	1:A:335:ALA:HB3	1.92	0.51
1:A:513:ILE:HD11	1:A:555:VAL:CG1	2.41	0.51
1:A:302:LEU:O	1:A:303:PRO:O	2.28	0.51
1:A:400:GLU:O	1:A:404:GLN:HG3	2.11	0.50
1:A:465:GLU:O	1:A:468:PRO:HD3	2.10	0.50
1:A:227:GLU:O	1:A:230:GLU:N	2.44	0.50
1:A:306:ALA:HA	1:A:310:VAL:CG2	2.41	0.50
1:A:291:ALA:HA	2:A:1007:DKA:H82	1.93	0.50
1:A:487:CYS:O	1:A:490:ALA:HB3	2.12	0.50
1:A:214:TRP:CD1	1:A:343:VAL:HG11	2.47	0.49
1:A:33:GLN:HB2	1:A:84:TYR:CZ	2.47	0.49
1:A:240:LYS:HE3	1:A:244:GLU:OE2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:GLN:HE22	1:A:107:ASP:H	1.61	0.49
1:A:224:PRO:HD2	1:A:296:ASP:HB3	1.95	0.49
1:A:513:ILE:HD11	1:A:555:VAL:HG13	1.94	0.49
1:A:77:VAL:HG12	1:A:80:LEU:H	1.78	0.48
1:A:415:VAL:HG23	1:A:415:VAL:O	2.13	0.48
1:A:214:TRP:CZ2	1:A:218:ARG:HD3	2.48	0.48
1:A:366:PRO:C	1:A:368:GLU:N	2.64	0.48
1:A:495:GLU:OE1	1:A:495:GLU:HA	2.13	0.48
1:A:480:SER:OG	1:A:483:ASN:HB2	2.14	0.48
1:A:510:HIS:C	1:A:512:ASP:H	2.17	0.48
1:A:11:PHE:CZ	1:A:51:LYS:HD2	2.49	0.48
1:A:532:LEU:HD12	1:A:532:LEU:O	2.13	0.48
1:A:296:ASP:OD1	1:A:297:GLU:N	2.46	0.47
1:A:225:LYS:HE3	1:A:297:GLU:O	2.14	0.47
1:A:10:ARG:O	1:A:14:LEU:HB2	2.15	0.47
1:A:267:ASN:O	1:A:271:ILE:HD12	2.15	0.47
1:A:433:VAL:HG22	1:A:452:TYR:HB3	1.96	0.47
1:A:5:SER:OG	1:A:62:CYS:HB3	2.14	0.47
1:A:560:LYS:HB3	1:A:560:LYS:HE2	1.60	0.47
1:A:160:ARG:NH2	1:A:188:GLU:OE1	2.39	0.47
1:A:480:SER:OG	1:A:483:ASN:N	2.47	0.47
1:A:6:GLU:O	1:A:10:ARG:HG2	2.14	0.47
1:A:161:TYR:CE1	1:A:185:LEU:HD23	2.49	0.46
1:A:165:PHE:CE1	1:A:178:LEU:HD21	2.50	0.46
1:A:194:ALA:HB2	2:A:1008:DKA:H91	1.97	0.46
1:A:459:GLN:HG3	2:A:1008:DKA:H72	1.97	0.46
1:A:16:GLU:OE1	1:A:51:LYS:HE2	2.15	0.45
1:A:383:GLU:HB3	1:A:384:PRO:HD3	1.98	0.45
1:A:516:LEU:HD22	1:A:520:GLU:CB	2.47	0.45
1:A:80:LEU:HA	1:A:84:TYR:HD2	1.81	0.45
1:A:315:VAL:HG12	1:A:316:CYS:N	2.32	0.45
1:A:39:HIS:HD2	1:A:140:TYR:CE2	2.34	0.45
1:A:33:GLN:HB2	1:A:84:TYR:CE1	2.52	0.45
1:A:518:GLU:O	1:A:522:GLN:HG3	2.17	0.45
1:A:383:GLU:HB3	1:A:384:PRO:CD	2.47	0.45
1:A:76:THR:HG22	1:A:76:THR:O	2.17	0.44
1:A:214:TRP:CH2	1:A:218:ARG:HD3	2.52	0.44
1:A:479:GLU:HG3	1:A:480:SER:H	1.80	0.44
1:A:5:SER:HB3	1:A:8:ALA:HB3	1.98	0.44
1:A:420:THR:HG23	1:A:530:VAL:CG1	2.47	0.44
1:A:420:THR:N	1:A:421:PRO:CD	2.80	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:409:VAL:O	1:A:413:LYS:HG3	2.17	0.44
1:A:464:HIS:CE1	1:A:469:VAL:H	2.36	0.44
1:A:412:THR:OG1	1:A:423:LEU:HD13	2.18	0.43
1:A:483:ASN:C	1:A:486:PRO:HD2	2.39	0.43
1:A:77:VAL:HG12	1:A:80:LEU:HB2	2.00	0.43
1:A:366:PRO:O	1:A:369:CYS:N	2.50	0.43
1:A:530:VAL:HG12	1:A:534:LYS:HE3	1.99	0.43
1:A:257:ARG:HH21	1:A:287:SER:CB	2.32	0.43
1:A:459:GLN:O	1:A:463:LEU:HG	2.18	0.43
1:A:426:VAL:O	1:A:430:LEU:HD13	2.19	0.43
1:A:87:MET:HE2	1:A:87:MET:HB2	1.92	0.43
1:A:583:LEU:O	1:A:584:GLY:O	2.37	0.43
1:A:118:PRO:HG2	1:A:123:MET:CG	2.45	0.42
1:A:441:PRO:O	1:A:443:ALA:N	2.51	0.42
1:A:481:LEU:HD12	2:A:1009:DKA:H81	2.01	0.42
1:A:408:LEU:HD13	1:A:427:SER:CB	2.49	0.42
1:A:14:LEU:CD1	1:A:283:LEU:HD21	2.50	0.42
1:A:564:LYS:C	1:A:566:THR:H	2.23	0.42
1:A:192:SER:O	1:A:196:GLN:HG3	2.19	0.42
1:A:34:CYS:HA	1:A:35:PRO:HD3	1.88	0.42
1:A:558:CYS:C	1:A:560:LYS:H	2.22	0.42
1:A:234:LEU:O	1:A:238:LEU:HB2	2.20	0.42
1:A:201:ALA:O	1:A:205:LYS:HB2	2.19	0.42
1:A:378:LYS:CB	1:A:379:PRO:HD3	2.51	0.41
1:A:251:LEU:HD23	1:A:251:LEU:HA	1.75	0.41
1:A:408:LEU:CD2	1:A:526:GLN:HG2	2.50	0.41
1:A:257:ARG:HE	1:A:257:ARG:HB3	1.68	0.41
1:A:357:LEU:HA	1:A:357:LEU:HD23	1.90	0.41
1:A:66:LEU:HD23	1:A:66:LEU:HA	1.83	0.41
1:A:59:ALA:O	1:A:62:CYS:HB2	2.20	0.41
1:A:299:PRO:HB2	1:A:302:LEU:HD23	2.03	0.41
1:A:442:GLU:HA	1:A:445:ARG:HD2	2.02	0.41
1:A:136:LYS:HD2	1:A:136:LYS:HA	1.66	0.41
1:A:575:LEU:C	1:A:577:ALA:N	2.73	0.41
1:A:311:GLU:O	1:A:312:SER:C	2.59	0.41
1:A:240:LYS:O	1:A:244:GLU:HG3	2.20	0.40
1:A:564:LYS:O	1:A:566:THR:N	2.54	0.40
1:A:95:GLU:HG3	3:A:2006:HOH:O	2.22	0.40
1:A:299:PRO:O	1:A:302:LEU:HD21	2.21	0.40
1:A:61:ASN:C	1:A:63:ASP:N	2.74	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	580/585 (99%)	520 (90%)	47 (8%)	13 (2%)	7 11

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	303	PRO
1	A	364	ALA
1	A	511	ALA
1	A	293	VAL
1	A	367	HIS
1	A	312	SER
1	A	314	ASP
1	A	442	GLU
1	A	497	TYR
1	A	565	GLU
1	A	499	PRO
1	A	120	VAL
1	A	85	GLY

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	460/511 (90%)	398 (86%)	62 (14%)	4 8

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	SER
1	A	7	VAL
1	A	14	LEU
1	A	31	LEU
1	A	51	LYS
1	A	52	THR
1	A	81	ARG
1	A	93	LYS
1	A	100	GLU
1	A	103	LEU
1	A	107	ASP
1	A	114	ARG
1	A	116	VAL
1	A	131	GLU
1	A	137	LYS
1	A	138	TYR
1	A	140	TYR
1	A	153	GLU
1	A	156	PHE
1	A	179	LEU
1	A	185	LEU
1	A	203	LEU
1	A	205	LYS
1	A	218	ARG
1	A	230	GLU
1	A	233	LYS
1	A	238	LEU
1	A	262	LYS
1	A	267	ASN
1	A	274	LYS
1	A	283	LEU
1	A	293	VAL
1	A	302	LEU
1	A	315	VAL
1	A	317	LYS
1	A	329	MET
1	A	334	TYR
1	A	337	ARG
1	A	359	LYS
1	A	373	VAL
1	A	376	GLU
1	A	380	LEU
1	A	387	LEU

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Mol	Chain	Res	Type
1	A	398	LEU
1	A	408	LEU
1	A	419	SER
1	A	425	GLU
1	A	435	SER
1	A	436	LYS
1	A	450	GLU
1	A	453	LEU
1	A	457	LEU
1	A	479	GLU
1	A	480	SER
1	A	481	LEU
1	A	484	ARG
1	A	498	VAL
1	A	512	ASP
1	A	513	ILE
1	A	525	LYS
1	A	540	THR
1	A	544	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	39	HIS
1	A	105	HIS
1	A	242	HIS
1	A	247	HIS
1	A	267	ASN
1	A	535	HIS

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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

10 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	$\# Z > 2$
2	DKA	A	1001	-	7,10,11	0.32	0	6,10,11	1.01	1 (16%)
2	DKA	A	1002	-	6,9,11	0.22	0	5,9,11	1.15	1 (20%)
2	DKA	A	1003	-	5,8,11	0.30	0	4,8,11	1.14	1 (25%)
2	DKA	A	1004	-	8,11,11	0.19	0	7,11,11	0.99	1 (14%)
2	DKA	A	1005	-	7,10,11	0.18	0	6,10,11	1.08	1 (16%)
2	DKA	A	1006	-	7,10,11	0.22	0	6,10,11	1.03	1 (16%)
2	DKA	A	1007	-	9,9,11	0.34	0	8,8,11	0.50	0
2	DKA	A	1008	-	8,11,11	0.19	0	7,11,11	0.98	1 (14%)
2	DKA	A	1009	-	8,11,11	0.22	0	7,11,11	0.97	1 (14%)
2	DKA	A	1010	-	7,7,11	0.26	0	6,6,11	0.45	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
2	DKA	A	1001	-	-	0/6/8/9	0/0/0/0
2	DKA	A	1002	-	-	0/5/7/9	0/0/0/0
2	DKA	A	1003	-	-	0/4/6/9	0/0/0/0
2	DKA	A	1004	-	-	0/7/9/9	0/0/0/0
2	DKA	A	1005	-	-	0/6/8/9	0/0/0/0
2	DKA	A	1006	-	-	0/6/8/9	0/0/0/0
2	DKA	A	1007	-	-	0/7/7/9	0/0/0/0
2	DKA	A	1008	-	-	0/7/9/9	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	DKA	A	1009	-	-	0/7/9/9	0/0/0/0
2	DKA	A	1010	-	-	0/5/5/9	0/0/0/0

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1002	DKA	C4-C3-C2	-2.40	104.22	113.75
2	A	1005	DKA	C4-C3-C2	-2.36	104.40	113.75
2	A	1001	DKA	C4-C3-C2	-2.35	104.45	113.75
2	A	1004	DKA	C4-C3-C2	-2.31	104.60	113.75
2	A	1009	DKA	C4-C3-C2	-2.30	104.62	113.75
2	A	1008	DKA	C4-C3-C2	-2.27	104.74	113.75
2	A	1006	DKA	C4-C3-C2	-2.25	104.83	113.75
2	A	1003	DKA	C4-C3-C2	-2.18	105.11	113.75

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1001	DKA	1	0
2	A	1002	DKA	1	0
2	A	1004	DKA	1	0
2	A	1005	DKA	2	0
2	A	1006	DKA	1	0
2	A	1007	DKA	1	0
2	A	1008	DKA	2	0
2	A	1009	DKA	1	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 [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	582/585 (99%)	0.03	14 (2%) 59 62	26, 54, 93, 100	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	555	VAL	3.9
1	A	561	ALA	3.7
1	A	513	ILE	3.4
1	A	502	PHE	3.1
1	A	562	ASP	2.8
1	A	558	CYS	2.4
1	A	363	ALA	2.4
1	A	443	ALA	2.4
1	A	553	ALA	2.3
1	A	559	CYS	2.3
1	A	507	PHE	2.3
1	A	563	ASP	2.1
1	A	302	LEU	2.0
1	A	301	ASP	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. 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(\AA^2)	Q<0.9
2	DKA	A	1007	10/12	0.61	0.41	70,74,78,78	0
2	DKA	A	1009	12/12	0.81	0.30	73,76,87,88	0
2	DKA	A	1008	12/12	0.81	0.31	73,76,86,86	0
2	DKA	A	1006	11/12	0.81	0.35	64,77,91,92	0
2	DKA	A	1005	11/12	0.88	0.24	54,59,74,75	0
2	DKA	A	1001	11/12	0.88	0.31	75,79,80,80	0
2	DKA	A	1004	12/12	0.89	0.31	53,58,62,63	0
2	DKA	A	1010	8/12	0.90	0.28	47,52,60,60	0
2	DKA	A	1003	9/12	0.92	0.28	62,64,69,71	0
2	DKA	A	1002	10/12	0.95	0.15	45,47,58,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.