

a)CYP81AA2 rigid receptor, all 25 docking results

Global docking result analysis

25 VINA docking runs of the ligand object 2 to the receptor object 1 yielded the following results, sorted by binding energy [more positive energies indicate stronger binding, and negative energies mean no binding]

Run	Bind.energy[kcal/mol]	Dissoc. constant [pM]	Contacting receptor residues
001	000007.9230	00000001557270.2500	PHE 116 ILE 218 VAL 219 MET 306 ILE 309 ALA 310 THR 314 SER
375 LEU 378 LEU 379 CYS 445 ALA 484 SER 485 HEM A 600			
002	000007.9230	00000001557270.2500	PHE 116 ILE 218 VAL 219 MET 306 ILE 309 ALA 310 THR 314 SER
375 LEU 378 LEU 379 CYS 445 ALA 484 SER 485 HEM A 600			
003	000007.7130	00000002219709.0000	VAL 115 PHE 116 ILE 218 VAL 219 MET 306 ALA 310 THR 314 SER
375 LEU 378 LEU 379 CYS 445 ALA 484 SER 485 HEM A 600			
004	000007.7130	00000002219709.0000	VAL 115 PHE 116 ILE 218 VAL 219 MET 306 ALA 310 THR 314 SER
375 LEU 378 LEU 379 CYS 445 ALA 484 SER 485 HEM A 600			
005	000007.5570	00000002888321.7500	PHE 116 PHE 215 ILE 218 MET 306 ILE 309 ALA 310 THR 314 PRO
374 SER 375 LEU 379 CYS 445 ALA 484 SER 485 MET 486 HEM A 600			
006	000007.5570	00000002888321.7500	PHE 116 PHE 215 ILE 218 MET 306 ILE 309 ALA 310 THR 314 PRO
374 SER 375 LEU 379 CYS 445 ALA 484 SER 485 MET 486 HEM A 600			
007	000007.5200	00000003074446.2500	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ILE 309 GLU 313 THR
314 ALA 317 PRO 374 SER 375 LEU 379 SER 485 MET 486			
008	000007.5200	00000003074446.2500	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ILE 309 GLU 313 THR
314 ALA 317 PRO 374 SER 375 LEU 379 SER 485 MET 486			
009	000007.3530	00000004075487.2500	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ALA 310 THR 314 SER
375 LEU 379 CYS 445 ALA 484 SER 485 HEM A 600			
010	000007.3530	00000004075487.2500	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ALA 310 THR 314 SER
375 LEU 379 CYS 445 ALA 484 SER 485 HEM A 600			
011	000007.3310	00000004229663.0000	LEU 107 PHE 113 VAL 115 PHE 116 LEU 214 PHE 215 ILE 218 ARG
249 THR 305 MET 306 ILE 309 THR 314 SER 375 LEU 379 SER 485			
012	000007.1980	00000005294147.5000	VAL 115 PHE 116 ILE 218 MET 306 ILE 309 ALA 310 THR 314 SER
375 LEU 379 CYS 445 HEM A 600			
013	000007.1790	00000005466674.5000	LEU 107 PHE 113 VAL 115 PHE 116 LEU 214 PHE 215 ILE 218 ARG
249 MET 306 ILE 309 GLU 313 THR 314 SER 375 LEU 379 SER 485			
014	000007.0950	00000006299354.5000	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ILE 309 GLU 313 THR
314 ALA 317 PRO 374 SER 375 LEU 379 SER 485 MET 486			
015	000007.0760	00000006504639.5000	LEU 51 TYR 52 SER 375 PRO 376 LEU 377 LEU 378 LEU 379 PRO
380 MET 399 LEU 401 GLY 482 ALA 483 ALA 484 MET 486			
016	000007.0360	00000006958948.5000	LEU 107 THR 108 PHE 113 VAL 115 PHE 116 LEU 214 PHE 215 ILE
218 ARG 249 THR 305 MET 306 ILE 309 ALA 310 GLU 313 THR 314 LEU 379			
017	000007.0260	00000007077399.5000	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ALA 310 THR 314 PRO
374 SER 375 LEU 379 CYS 445 ALA 484 SER 485 HEM A 600			
018	000006.8570	00000009413527.0000	PHE 116 PHE 215 ILE 218 ILE 309 GLU 313 THR 314 SER 375 LEU
378 LEU 379 PRO 380 SER 485			

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019 |          000006.8020 | 00000010329229.0000 | VAL  115 PHE  116 ILE  218 VAL  219 MET  306 ILE  309 ALA  310 THR
314 SER  375 LEU  378 LEU  379 PRO  380 SER  485
020 |          000006.6950 | 00000012373710.0000 | VAL  115 PHE  116 PHE  215 ILE  218 MET  306 ILE  309 ALA  310 GLU
313 THR  314 LEU  379 CYS  445 HEM A 600
021 |          000006.6820 | 00000012648210.0000 | ILE  46 LEU  51 VAL  219 SER  375 PRO  376 LEU  377 LEU  378 PRO
380 ALA  483 ALA  484 SER  485 MET  486
022 |          000006.6490 | 00000013372678.0000 | VAL  115 PHE  116 PHE  215 ILE  218 MET  306 ILE  309 THR  314 PRO
374 SER  375 LEU  379 SER  485 HEM A 600
023 |          000006.6480 | 00000013395267.0000 | PHE  116 PHE  215 ILE  218 ILE  309 ALA  310 GLU  313 THR  314 SER
375 LEU  378 LEU  379 PRO  380 SER  485
024 |          000006.5540 | 00000015698361.0000 | VAL  115 PHE  116 PHE  215 ILE  218 VAL  219 MET  306 ILE  309 ALA
310 GLU  313 THR  314 LEU  379 ALA  484 SER  485 HEM A 600
025 |          000006.4300 | 00000019352964.0000 | PHE  116 PHE  215 ILE  218 ILE  309 ALA  310 GLU  313 THR  314 PRO
374 SER  375 LEU  377 LEU  378 LEU  379 PRO  380 ALA  484 SER  485 MET  486

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After clustering the 25 runs, the following 7 distinct complex conformations were found:
[They all differ by at least 5.0 A heavy atom RMSD]

Clu |Bind.energy[kcal/mol]|Dissoc. constant [pM]| Contacting receptor residues

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-----+-----+-----+-----
001 |          000007.9230 | 00000001557270.2500 | PHE  116 ILE  218 VAL  219 MET  306 ILE  309 ALA  310 THR  314 SER
375 LEU  378 LEU  379 CYS  445 ALA  484 SER  485 HEM A 600
002 |          000007.7130 | 00000002219709.0000 | VAL  115 PHE  116 ILE  218 VAL  219 MET  306 ALA  310 THR  314 SER
375 LEU  378 LEU  379 CYS  445 ALA  484 SER  485 HEM A 600
003 |          000007.5200 | 00000003074446.2500 | VAL  115 PHE  116 PHE  215 ILE  218 MET  306 ILE  309 GLU  313 THR
314 ALA  317 PRO  374 SER  375 LEU  379 SER  485 MET  486
004 |          000007.3310 | 00000004229663.0000 | LEU  107 PHE  113 VAL  115 PHE  116 LEU  214 PHE  215 ILE  218 ARG
249 THR  305 MET  306 ILE  309 THR  314 SER  375 LEU  379 SER  485
005 |          000007.0760 | 00000006504639.5000 | LEU  51 TYR  52 SER  375 PRO  376 LEU  377 LEU  378 LEU  379 PRO
380 MET  399 LEU  401 GLY  482 ALA  483 ALA  484 MET  486
006 |          000007.0360 | 00000006958948.5000 | LEU  107 THR  108 PHE  113 VAL  115 PHE  116 LEU  214 PHE  215 ILE
218 ARG  249 THR  305 MET  306 ILE  309 ALA  310 GLU  313 THR  314 LEU  379
007 |          000006.6820 | 00000012648210.0000 | ILE  46 LEU  51 VAL  219 SER  375 PRO  376 LEU  377 LEU  378 PRO
380 ALA  483 ALA  484 SER  485 MET  486

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b) CYP81AA2 flexible receptor sidechains, all 100 docking results

Ensemble docking result analysis

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The ligand was docked 20 times ["Run"] with VINA against each of the 5 receptors ["Rec"] in the ensemble, yielding the following 100 results, sorted by binding energy:

[More positive energies indicate stronger binding, and negative energies mean no binding]

Rec	Run	Bind.energy[kcal/mol]	Dissoc. constant [pM]	Contacting receptor residues
002	001	000008.2560	00000000887711.3125	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ILE 309 ALA 310
THR	314 PRO	374 SER 375 LEU	379 CYS 445 ALA	484 SER 485 MET 486 HEM A 600
005	001	000008.2050	00000000967509.5625	VAL 115 PHE 116 PHE 215 ILE 218 ILE 309 ALA 310 THR 314
PRO	374 SER	375 LEU 379 CYS	445 ALA 484 SER	485 MET 486 HEM A 600
003	001	000008.0680	00000001219207.3750	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ILE 309 ALA 310
THR	314 PRO	374 SER 375 LEU	379 CYS 445 ALA	484 SER 485 MET 486 HEM A 600
004	001	000007.9640	00000001453150.5000	PHE 215 ILE 218 ILE 309 ALA 310 THR 314 PRO 374 SER 375
CYS	445 ALA	484 SER 485 MET	486 HEM A 600	
005	002	000007.9390	00000001515778.8750	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ILE 309 GLU 313
THR	314 ALA	317 PRO 374 SER	375 LEU 379 SER	485
001	001	000007.9190	00000001567819.3750	VAL 115 PHE 116 ILE 218 VAL 219 MET 306 ILE 309 ALA 310
THR	314 SER	375 LEU 378 LEU	379 CYS 445 ALA	484 SER 485 HEM A 600
005	003	000007.8210	00000001849826.5000	VAL 115 PHE 116 PHE 215 ILE 218 VAL 219 MET 306 ILE 309
ALA	310 THR	314 SER 375 LEU	378 LEU 379 CYS	445 ALA 484 SER 485 HEM A 600
004	002	000007.8170	00000001862357.3750	VAL 115 PHE 116 ILE 218 MET 306 ALA 310 THR 314 PRO 374
SER	375 LEU	379 CYS 445 HEM A 600		
005	004	000007.7460	00000002099455.7500	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ALA 310 THR 314
PRO	374 SER	375 LEU 379 ALA	484 SER 485 HEM A 600	
001	002	000007.7200	00000002193638.0000	VAL 115 PHE 116 ILE 218 VAL 219 MET 306 ALA 310 THR 314
PRO	374 SER	375 LEU 378 LEU	379 CYS 445 ALA	484 SER 485 HEM A 600
003	002	000007.7140	00000002215965.7500	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ALA 310 THR 314
PRO	374 SER	375 LEU 379 CYS	445 ALA 484 SER	485 HEM A 600
002	002	000007.6980	00000002276623.2500	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ALA 310 THR 314
PRO	374 SER	375 LEU 379 CYS	445 ALA 484 SER	485 HEM A 600
004	003	000007.6310	00000002549193.7500	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ALA 310 THR 314
PRO	374 SER	375 LEU 379 CYS	445 ALA 484 SER	485 HEM A 600
005	005	000007.6150	00000002618972.7500	VAL 115 PHE 116 ILE 218 VAL 219 MET 306 ALA 310 SER 375
LEU	378 LEU	379 CYS 445 ALA	484 SER 485 HEM A 600	
001	003	000007.5550	00000002898088.0000	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ALA 310 THR 314
PRO	374 SER	375 LEU 379 CYS	445 ALA 484 SER	485 MET 486 HEM A 600
004	004	000007.5550	00000002898088.0000	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ALA 310 THR 314
PRO	374 SER	375 LEU 379 SER	485 HEM A 600	
001	004	000007.5170	00000003090053.0000	VAL 115 PHE 116 PHE 215 ILE 218 MET 306 ILE 309 GLU 313
THR	314 ALA	317 PRO 374 SER	375 LEU 379 SER	485 MET 486


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001 | 013 |      000006.3100 | 00000023697836.0000 | PHE  215 ILE  218 VAL  219 ILE  309 ALA  310 GLU  313 THR  314
SER  375 LEU  378 LEU  379 PRO  380 ALA  484 SER  485
002 | 015 |      000006.2560 | 00000025959194.0000 | LEU  51 TYR  52 PHE  116 VAL  219 SER  375 PRO  376 LEU  377
LEU  378 LEU  379 PRO  380 ALA  483 ALA  484
002 | 016 |      000006.2550 | 00000026003044.0000 | LEU  51 PHE  116 ILE  218 VAL  219 SER  375 PRO  376 LEU  377
LEU  378 LEU  379 ALA  484
001 | 014 |      000006.2230 | 00000027446092.0000 | VAL  115 PHE  116 PHE  215 ILE  218 MET  306 ILE  309 ALA  310
THR  314 PRO  374 SER  375 LEU  377 LEU  378 LEU  379 PRO  380 ALA  484 SER  485 MET  486
002 | 017 |      000006.2170 | 00000027725448.0000 | ILE  46 LEU  51 TYR  52 VAL  219 SER  375 PRO  376 LEU  377 LEU
378 ALA  483 ALA  484 SER  485 MET  486
003 | 018 |      000006.1790 | 00000029561940.0000 | LEU  51 PHE  116 PHE  215 ILE  218 VAL  219 SER  375 PRO  376
LEU  377 LEU  378 LEU  379 PRO  380 ALA  484 SER  485
001 | 015 |      000006.1740 | 00000029812472.0000 | LEU  51 VAL  115 PHE  116 ILE  218 VAL  219 SER  375 PRO  376
LEU  377 LEU  378 LEU  379 ALA  484
001 | 016 |      000006.1690 | 00000030065128.0000 | LEU  51 VAL  115 PHE  116 MET  306 PRO  374 SER  375 PRO  376
LEU  377 LEU  378 LEU  379 PRO  380 ALA  484 SER  485 MET  486
001 | 017 |      000006.1680 | 00000030115914.0000 | VAL  115 PHE  116 PHE  215 MET  306 ILE  309 ALA  310 GLU  313
THR  314 ALA  317 PRO  374 SER  375 LEU  379 SER  485 MET  486 HEM A 600
002 | 018 |      000006.1590 | 00000030576878.0000 | ILE  46 LYS  47 PRO  48 GLY  49 PRO  50 LEU  51 VAL  219 PHE
226 ALA  483 ALA  484
003 | 019 |      000006.1550 | 00000030784010.0000 | ILE  46 LYS  47 PRO  48 GLY  49 PRO  50 LEU  51 VAL  219 ALA
483 ALA  484 SER  485
001 | 018 |      000006.1100 | 00000033213192.0000 | LEU  51 THR  101 PHE  116 ILE  218 VAL  219 PRO  374 SER  375
PRO  376 LEU  377 LEU  378 LEU  379 PRO  380 ALA  484 SER  485 MET  486
002 | 019 |      000006.1100 | 00000033213192.0000 | ILE  46 LEU  51 TYR  52 VAL  219 PRO  374 SER  375 PRO  376 LEU
377 LEU  378 GLY  482 ALA  483 ALA  484 SER  485 MET  486
002 | 020 |      000006.1010 | 00000033721564.0000 | LEU  51 PHE  116 ILE  218 VAL  219 PRO  374 SER  375 PRO  376
LEU  377 LEU  378 LEU  379 PRO  380 ALA  484 SER  485 MET  486
003 | 020 |      000006.0880 | 00000034469648.0000 | PHE  116 PHE  215 ILE  218 MET  306 ILE  309 ALA  310 GLU  313
THR  314 PRO  374 SER  375 PRO  376 LEU  377 LEU  378 LEU  379 PRO  380 ALA  484 SER  485 MET  486
001 | 019 |      000006.0620 | 00000036015968.0000 | VAL  115 PHE  116 LEU  214 PHE  215 ILE  218 MET  306 ILE  309
GLU  313 THR  314 PRO  374 SER  375 LEU  379 SER  485 MET  486
001 | 020 |      000006.0360 | 00000037631656.0000 | VAL  115 PHE  116 ILE  218 VAL  219 MET  306 ILE  309 ALA  310
SER  375 LEU  378 LEU  379 ALA  484

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After clustering the 100 runs, the following 9 distinct complex conformations were found:

[They all differ by at least 5.0 A heavy ligand atom RMSD, and have been saved as YOb files with terminal number "Num"]

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Num | Rec | Run | Bind.energy[kcal/mol] | Dissoc. constant [pM] | Contacting receptor residues
-----+-----+-----+-----+-----+-----
001 | 002 | 001 |      000008.2560 | 00000000887711.3125 | VAL  115 PHE  116 PHE  215 ILE  218 MET  306 ILE  309 ALA
310 THR  314 PRO  374 SER  375 LEU  379 CYS  445 ALA  484 SER  485 MET  486 HEM A 600
002 | 001 | 002 |      000007.7200 | 00000002193638.0000 | VAL  115 PHE  116 ILE  218 VAL  219 MET  306 ALA  310 THR
314 PRO  374 SER  375 LEU  378 LEU  379 CYS  445 ALA  484 SER  485 HEM A 600

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c)CYP81AA2(mut6) rigid receptor, all 25 docking results

Global docking result analysis

25 VINA docking runs of the ligand object 2 to the receptor object 1 yielded the following results, sorted by binding energy [more positive energies indicate stronger binding, and negative energies mean no binding]

Run	Bind.energy[kcal/mol]	Dissoc. constant [pM]	Contacting receptor residues
001 314 PRO	000008.4470 374 ALA 375 LEU	00000000643082.8125 379 SER 485 HEM A 600	VAL 115 PHE 116 PHE 215 THR 305 MET 306 ILE 309 ALA 310 THR
002 314 PRO	000008.4470 374 ALA 375 LEU	00000000643082.8125 379 SER 485 HEM A 600	VAL 115 PHE 116 PHE 215 THR 305 MET 306 ILE 309 ALA 310 THR
003 306 ILE	000008.2310 309 ALA 310 LEU	00000000925970.1875 379 HEM A 600	VAL 115 PHE 116 LEU 214 PHE 215 ILE 218 ARG 249 THR 305 MET
004 306 ILE	000008.2310 309 ALA 310 LEU	00000000925970.1875 379 HEM A 600	VAL 115 PHE 116 LEU 214 PHE 215 ILE 218 ARG 249 THR 305 MET
005 305 MET	000008.0280 306 ILE 309 ALA	00000001304361.5000 310 THR 314 ALA 484	LEU 107 VAL 115 LEU 214 PHE 215 ILE 218 ILE 243 ARG 249 THR
006 305 MET	000008.0280 306 ILE 309 ALA	00000001304361.5000 310 THR 314 ALA 484	LEU 107 VAL 115 LEU 214 PHE 215 ILE 218 ILE 243 ARG 249 THR
007 305 MET	000008.0100 306 ILE 309 ALA	00000001344597.0000 310 LEU 379 HEM A 600	LEU 107 VAL 115 LEU 214 PHE 215 ILE 218 ILE 243 ARG 249 THR
008 305 MET	000008.0100 306 ILE 309 ALA	00000001344597.0000 310 LEU 379 HEM A 600	LEU 107 VAL 115 LEU 214 PHE 215 ILE 218 ILE 243 ARG 249 THR
009 218 ILE	000008.0030 243 ARG 249 THR	00000001360577.2500 305 MET 306 ILE 309 ALA 310 THR 314	LEU 107 THR 108 PHE 113 VAL 115 PHE 116 LEU 214 PHE 215 ILE
010 218 ILE	000008.0030 243 ARG 249 THR	00000001360577.2500 305 MET 306 ILE 309 ALA 310 THR 314	LEU 107 THR 108 PHE 113 VAL 115 PHE 116 LEU 214 PHE 215 ILE
011 305 MET	000007.9940 306 ILE 309 ALA	00000001381402.7500 310 LEU 379 ALA 484 HEM A 600	VAL 115 PHE 116 LEU 214 PHE 215 ILE 218 ILE 243 ARG 249 THR
012 218 ARG	000007.7560 249 GLY 302 THR	00000002064318.1250 305 MET 306 ILE 309 ALA 310 LEU 379	LEU 107 THR 108 PHE 113 VAL 115 PHE 116 LEU 214 PHE 215 ILE
013 306 ILE	000007.6370 309 ALA 310 LEU	00000002523508.5000 379 ALA 484 HEM A 600	LEU 107 PHE 113 VAL 115 PHE 116 PHE 215 ILE 218 THR 305 MET
014 LEU 379 PRO	000007.6230 380 LEU 401 ALA	00000002583847.7500 481 GLY 482 THR 483 ALA 484	GLY 49 PRO 50 LEU 51 TYR 52 VAL 219 PRO 376 LEU 377 SER 378
015 306 ILE	000007.5370 309 ALA 310 LEU	00000002987485.0000 379 ALA 484 HEM A 600	VAL 115 LEU 214 PHE 215 ILE 218 ILE 243 ARG 249 THR 305 MET
016 309 ALA	000007.4010 310 ALA 375 LEU	00000003758331.2500 379 ALA 484 SER 485 HEM A 600	VAL 115 PHE 116 LEU 214 PHE 215 ILE 218 THR 305 MET 306 ILE
017 310 THR	000007.3400 314 ALA 375 LEU	00000004165898.2500 379 ALA 484 HEM A 600	PHE 113 VAL 115 PHE 116 ILE 218 THR 305 MET 306 ILE 309 ALA

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018 |          000007.2920 | 00000004517447.5000 | GLY  49 PRO  50 LEU  51 TYR  52 VAL  219 ALA  375 PRO  376 LEU  377
SER  378 LEU  379 ALA  481 GLY  482 THR  483 ALA  484
019 |          000007.2890 | 00000004540379.5000 | LEU  107 PHE  113 VAL  115 PHE  116 PHE  215 ILE  218 THR  305 MET
306 ILE  309 ALA  310 THR  314 ALA  484 SER  485
020 |          000007.2020 | 00000005258525.5000 | LEU  107 PHE  113 VAL  115 PHE  116 LEU  214 PHE  215 ILE  218 ILE
243 GLY  246 ARG  249 THR  305 MET  306 ILE  309 ALA  310 LEU  379 ALA  484
021 |          000007.1530 | 00000005711911.5000 | PHE  116 PHE  215 ILE  218 ILE  309 ALA  310 GLU  313 THR  314 ALA
375 SER  378 LEU  379 PRO  380 ALA  484 SER  485
022 |          000007.0800 | 00000006460873.0000 | PHE  116 PHE  215 ILE  218 ILE  309 ALA  310 THR  314 ALA  375 LEU
379 ALA  484 HEM A 600
023 |          000007.0090 | 00000007283412.0000 | PHE  116 ILE  218 ILE  309 ALA  310 THR  314 ALA  375 SER  378 LEU
379 ALA  484 SER  485
024 |          000006.9850 | 00000007584502.5000 | VAL  115 PHE  116 PHE  215 ILE  218 THR  305 MET  306 ILE  309 ALA
310 THR  314 ALA  375 ALA  484 SER  485
025 |          000006.8740 | 00000009147264.0000 | VAL  115 PHE  116 ILE  218 MET  306 ILE  309 ALA  310 THR  314 ALA
375 SER  378 LEU  379 PRO  380 ALA  484

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After clustering the 25 runs, the following 7 distinct complex conformations were found:
[They all differ by at least 5.0 A heavy atom RMSD]

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Clu |Bind.energy[kcal/mol]|Dissoc. constant [pM]| Contacting receptor residues
-----+-----+-----+-----
001 |          000008.4470 | 00000000643082.8125 | VAL  115 PHE  116 PHE  215 THR  305 MET  306 ILE  309 ALA  310 THR
314 PRO  374 ALA  375 LEU  379 SER  485 HEM A 600
002 |          000008.2310 | 00000000925970.1875 | VAL  115 PHE  116 LEU  214 PHE  215 ILE  218 ARG  249 THR  305 MET
306 ILE  309 ALA  310 LEU  379 HEM A 600
003 |          000007.7560 | 00000002064318.1250 | LEU  107 THR  108 PHE  113 VAL  115 PHE  116 LEU  214 PHE  215 ILE
218 ARG  249 GLY  302 THR  305 MET  306 ILE  309 ALA  310 LEU  379
004 |          000007.6230 | 00000002583847.7500 | GLY  49 PRO  50 LEU  51 TYR  52 VAL  219 PRO  376 LEU  377 SER  378
LEU  379 PRO  380 LEU  401 ALA  481 GLY  482 THR  483 ALA  484
005 |          000007.2920 | 00000004517447.5000 | GLY  49 PRO  50 LEU  51 TYR  52 VAL  219 ALA  375 PRO  376 LEU  377
SER  378 LEU  379 ALA  481 GLY  482 THR  483 ALA  484
006 |          000007.1530 | 00000005711911.5000 | PHE  116 PHE  215 ILE  218 ILE  309 ALA  310 GLU  313 THR  314 ALA
375 SER  378 LEU  379 PRO  380 ALA  484 SER  485
007 |          000006.8740 | 00000009147264.0000 | VAL  115 PHE  116 ILE  218 MET  306 ILE  309 ALA  310 THR  314 ALA
375 SER  378 LEU  379 PRO  380 ALA  484

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d) CYP81AA2(mut6) flexible receptor sidechains, all 100 docking results

Ensemble docking result analysis

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The ligand was docked 20 times ["Run"] with VINA against each of the 5 receptors ["Rec"] in the ensemble, yielding the following 100 results, sorted by binding energy:

[More positive energies indicate stronger binding, and negative energies mean no binding]

Rec	Run	Bind.energy[kcal/mol]	Dissoc. constant [pM]	Contacting receptor residues
001	001	000008.4390	00000000651824.9375	VAL 115 PHE 116 PHE 215 THR 305 MET 306 ILE 309 ALA 310 THR 314 PRO 374 ALA 375 LEU 379 SER 485 HEM A 600
001	002	000008.2390	00000000913551.2500	VAL 115 PHE 116 LEU 214 PHE 215 ILE 218 ARG 249 THR 305 MET 306 ILE 309 ALA 310 LEU 379 HEM A 600
001	003	000008.0320	00000001295585.1250	LEU 107 VAL 115 LEU 214 PHE 215 ILE 218 ILE 243 ARG 249 THR 305 MET 306 ILE 309 ALA 310 THR 314 ALA 484
001	004	000008.0020	00000001362875.6250	LEU 107 VAL 115 LEU 214 PHE 215 ILE 218 ILE 243 ARG 249 THR 305 MET 306 ILE 309 ALA 310 LEU 379 HEM A 600
001	005	000008.0020	00000001362875.6250	LEU 107 VAL 115 PHE 116 LEU 214 PHE 215 ILE 218 ILE 243 ARG 249 THR 305 MET 306 ILE 309 ALA 310 LEU 379 ALA 484 HEM A 600
001	006	000007.9960	00000001376747.5000	LEU 107 THR 108 PHE 113 VAL 115 PHE 116 LEU 214 PHE 215 ILE 218 ILE 243 ARG 249 THR 305 MET 306 ILE 309 ALA 310 THR 314
001	007	000007.9070	00000001599897.5000	VAL 115 PHE 116 PHE 215 ILE 218 THR 305 MET 306 ILE 309 ALA 310 ALA 379 ALA 379 ALA 484 HEM A 600
001	008	000007.8600	00000001731982.8750	LEU 107 PHE 113 VAL 115 PHE 116 LEU 214 PHE 215 ILE 218 ILE 243 ARG 249 THR 305 MET 306 ILE 309 ALA 310
004	001	000007.8340	00000001809680.2500	ARG 98 PRO 99 LYS 100 THR 101 SER 103 GLY 104 PHE 116 ILE 218 VAL 219 LEU 379 PRO 380
001	009	000007.7610	00000002046970.3750	PHE 113 VAL 115 PHE 116 PHE 215 ILE 218 THR 305 MET 306 ILE 309 ALA 310 THR 314 SER 485 HEM A 600
002	001	000007.7520	00000002078302.0000	THR 101 SER 103 GLY 104 LEU 107 PHE 113 PHE 116 LEU 214 PRO 217 ILE 218 VAL 219 PRO 223 ILE 243 ARG 249 THR 305 MET 306 ILE 309
003	001	000007.7520	00000002078302.0000	THR 101 SER 103 GLY 104 LEU 107 PHE 113 PHE 116 LEU 214 PRO 217 ILE 218 VAL 219 PRO 223 ILE 243 ARG 249 THR 305 MET 306 ILE 309
005	001	000007.7280	00000002164217.5000	VAL 115 PHE 215 ILE 218 MET 306 ALA 310 THR 314 ALA 375 LEU 379 CYS 445 PRO 446 ALA 484 SER 485 HEM A 600
001	010	000007.7140	00000002215965.7500	VAL 115 PHE 116 PHE 215 MET 306 ALA 310 THR 314 PRO 374 ALA 375 LEU 379 ALA 484 SER 485 HEM A 600
001	011	000007.7010	00000002265125.0000	LEU 107 THR 108 PHE 113 VAL 115 PHE 116 LEU 214 PHE 215 ILE 218 ARG 249 GLY 302 THR 305 MET 306 ILE 309 ALA 310 LEU 379
001	012	000007.6870	00000002319286.0000	LEU 107 VAL 115 PHE 116 PHE 215 ILE 218 THR 305 MET 306 ILE 309 ALA 310 LEU 379 ALA 484 HEM A 600
001	013	000007.6450	00000002489663.7500	LEU 107 PHE 113 VAL 115 LEU 214 PHE 215 ILE 218 ILE 243 ARG 249 THR 305 MET 306 ILE 309 ALA 310 LEU 379 ALA 484

003		016			000006.6750		00000012798531.0000		VAL	115	PHE	116	PHE	215	ILE	218	MET	306	THR	314	PRO	374
ALA		375	LEU	379	ALA		484	SER	485	LEU		486	HEM	A	600							
002		017			000006.6610		00000013104554.0000		LEU	107	PHE	116	LEU	214	PHE	215	ILE	218	ILE	243	ARG	249
MET		306	ILE	309	ALA		310	LEU	379	ALA		484										
003		017			000006.6610		00000013104554.0000		LEU	107	PHE	116	LEU	214	PHE	215	ILE	218	ILE	243	ARG	249
MET		306	ILE	309	ALA		310	LEU	379	ALA		484										
004		016			000006.6540		00000013260299.0000		VAL	115	PHE	116	PHE	215	ILE	218	MET	306	THR	314	PRO	374
ALA		375	LEU	379	ALA		484	SER	485	LEU		486	HEM	A	600							
002		018			000006.6450		00000013463266.0000		LYS	173	LYS	212	PHE	215	PHE	216	GLN	480	SER	487	ARG	488
LYS		489	LYS	490	PRO		491	LEU	492													
003		018			000006.6450		00000013463266.0000		LYS	173	LYS	212	PHE	215	PHE	216	GLN	480	SER	487	ARG	488
LYS		489	LYS	490	PRO		491	LEU	492													
005		018			000006.6350		00000013692430.0000		ARG	98	PRO	99	LEU	107	PHE	113	VAL	115	PHE	116	PRO	217
ILE		218	ILE	243	LEU		379	PRO	380													
004		017			000006.6250		00000013925494.0000		LYS	173	LYS	212	PHE	215	PHE	216	GLN	480	THR	483	SER	487
ARG		488	LYS	489	LYS		490	PRO	491	LEU		492										
005		019			000006.6140		00000014186450.0000		THR	101	SER	103	GLY	104	LEU	107	PHE	113	VAL	115	PHE	116
LEU		214	ILE	243	ARG		249	THR	305	MET		306	ILE	309								
004		018			000006.5990		00000014550197.0000		THR	101	SER	103	LEU	107	PHE	116	PRO	217	ILE	218	PRO	223
ILE		242	ILE	243	LEU		379	PRO	380													
005		020			000006.5790		00000015049742.0000		THR	101	SER	103	GLY	104	LEU	107	PHE	113	LEU	214	PRO	217
ILE		218	PRO	223	ILE		243	GLY	246	ARG		249	MET	306	ILE	309						
004		019			000006.5100		00000016908564.0000		LEU	107	PHE	116	PHE	215	ILE	218	THR	305	MET	306	ILE	309
ALA		310	THR	314	LEU		379	ALA	484													
002		019			000006.4890		00000017518622.0000		PHE	116	PHE	215	ILE	218	MET	306	ILE	309	ALA	310	THR	314
ALA		375	LEU	379	ALA		484	HEM	A	600												
003		019			000006.4890		00000017518622.0000		PHE	116	PHE	215	ILE	218	MET	306	ILE	309	ALA	310	THR	314
ALA		375	LEU	379	ALA		484	HEM	A	600												
004		020			000006.4790		00000017816814.0000		VAL	115	PHE	116	PHE	215	ILE	218	MET	306	ILE	309	ALA	310
GLU		313	THR	314	PRO		374	ALA	375	SER		378	LEU	379	SER		485	HEM	A	600		
002		020			000006.4690		00000018120080.0000		PHE	116	PHE	215	ILE	218	THR	305	MET	306	ILE	309	ALA	310
THR		314	PRO	374	ALA		375	LEU	379	ALA		484	SER	485								
003		020			000006.4690		00000018120080.0000		PHE	116	PHE	215	ILE	218	THR	305	MET	306	ILE	309	ALA	310
THR		314	PRO	374	ALA		375	LEU	379	ALA		484	SER	485								

After clustering the 100 runs, the following 12 distinct complex conformations were found:

[They all differ by at least 5.0 A heavy ligand atom RMSD, and have been saved as YOb files with terminal number "Num"]

Num		Rec		Run		Bind.energy[kcal/mol]		Dissoc. constant [pM]		Contacting receptor residues
001		001		001		000008.4390		00000000651824.9375		VAL 115 PHE 116 PHE 215 THR 305 MET 306 ILE 309 ALA
310		THR		314		PRO		374		ALA 375 LEU 379 SER 485 HEM A 600
002		001		002		000008.2390		00000000913551.2500		VAL 115 PHE 116 LEU 214 PHE 215 ILE 218 ARG 249 THR
305		MET		306		ILE		309		ALA 310 LEU 379 HEM A 600

003 001 003	000007.8600	00000001731982.8750	LEU	107 PHE	113 VAL	115 PHE	116 LEU	214 PHE	215 ILE
218 ILE 243 ARG	249 THR 305 MET	306 ILE 309 ALA	310						
004 004 001	000007.8340	00000001809680.2500	ARG	98 PRO	99 LYS	100 THR	101 SER	103 GLY	104 PHE
116 ILE 218 VAL	219 LEU 379 PRO	380							
005 002 001	000007.7520	00000002078302.0000	THR	101 SER	103 GLY	104 LEU	107 PHE	113 PHE	116 LEU
214 PRO 217 ILE	218 VAL 219 PRO	223 ILE 243 ARG	249 THR	305 MET	306 ILE	309			
006 001 004	000007.7140	00000002215965.7500	VAL	115 PHE	116 PHE	215 MET	306 ALA	310 THR	314 PRO
374 ALA 375 LEU	379 ALA 484 SER	485 HEM A 600							
007 001 005	000007.6280	00000002562134.2500	GLY	49 PRO	50 LEU	51 TYR	52 VAL	219 PRO	376 LEU
377 SER 378 LEU	379 PRO 380 LEU	401 ALA 481 GLY	482 THR	483 ALA	484				
008 005 002	000007.5590	00000002878588.2500	THR	101 SER	103 GLY	104 LEU	107 PHE	113 VAL	115 PHE
116 LEU 214 PRO	217 ILE 218 PRO	223 ILE 242 ILE	243 MET	306 ILE	309				
009 004 003	000007.4920	00000003223229.2500	ARG	98 PRO	99 LYS	100 THR	101 GLY	104 THR	108 PHE
116 ILE 218 VAL	219 PRO 380								
010 002 003	000007.4880	00000003245063.7500	LYS	173 LYS	212 PHE	215 PHE	216 VAL	219 GLN	480 GLY
482 THR 483 SER	487								
011 004 005	000007.3240	00000004279931.5000	GLY	49 PRO	50 LEU	51 TYR	52 VAL	219 ALA	375 PRO
376 LEU 377 SER	378 LEU 379 ALA	481 GLY 482 THR	483 ALA	484					
012 005 007	000007.2040	00000005240805.0000	LYS	173 LYS	212 PHE	215 PHE	216 ILE	218 VAL	219 GLN
480 GLY 482 THR	483 ALA 484 SER	487 PRO 491							