

COEFICIENTE DE CORRELACIÓN CUERPO ROSARIO - TAJO 238

	Au	Ag	Al	As	B	Be	Be	Bi	Ca	Cd	Co	Cr	Cu	Fe	Ga	Hg	K	La	Mg	Mn	Mo	Na	Ni	P	Pb	S	Sb	Sc	Sr	Ti	Ti	U	V	W	Zn	
Au	0.357750085																																			
Ag	-0.220807652	1																																		
Al	0.358109237	0.9987944	1																																	
As	0.31187247	0.8913957	-0.07587774	1																																
B	-0.180936716	0.46424245	-0.12429577	0.44371541	1																															
Be	0.418330013	0.1496576	0.24632125	0.14758037	0.5925064	1																														
Bi	0.239045722	0.69606672	0.05473806	0.66322066	0.94587571	0.60553007	1																													
Ca	0.418330013	0.1496576	-0.20526771	0.14758037	0.13046561	0.30276504	-0.1	1																												
Cd	0.35099928	0.97842681	-0.30372291	0.98646078	0.7785449	0.37409062	0.14683353	0.53273583	0.14683353	1																										
Co	0.3479445	0.27536346	0.8493457	0.87011839	0.4367701	0.08813831	0.14555563	0.49488913	0.14555563	0.18193611	1																									
Cr	0.198237982	0.85483885	0.04539368	0.83240789	0.98072641	0.58920032	0.1194176	0.95081793	0.07078599	0.73418019	0.46351743	1																								
Cu	0.327019146	0.88730803	-0.35666884	0.90921191	0.5819615	0.22806237	0.13615357	0.28657085	0.13745027	0.96341004	0.05624514	0.530042	1																							
Fe	0.356768361	0.83238025	-0.06953631	0.91702853	0.99084082	0.54908004	0.14155599	0.90204297	0.15102174	0.83942601	0.47388924	0.97263818	0.66542029	1																						
Ga	0.239045722	0.64476732	-0.3968509	0.67807049	0.22831483	0	0.1	-0.1	0.1	0.78876597	-0.14855563	0.17249209	0.9245476	0.33173152	1																					
Hg	0.218586777	0.90843882	-0.10504133	0.89488178	0.93423346	0.68628254	0.19061026	0.82941222	0.24727818	0.83226837	0.34942942	0.92529906	0.87896745	0.94136636	0.37606889	1																				
K	-0.338421033	-0.51428011	0.92106657	-0.50746525	-0.39579076	-0.34290363	0.0808981	-0.25887392	-0.27505354	-0.53149017	0.45216867	-0.27765444	-0.51506206	-0.38521269	-0.43684974	-0.42009103	1																			
La	-0.46291005	-0.25779619	-0.10599979	-0.23298061	-0.44212976	0	-0.19364917	-0.51639778	-0.19364917	-0.15431822	-0.33824071	-0.42824294	-0.01590332	-0.41021093	0.19364917	-0.25272751	0.0417756	1																		
Mg	0.172926777	0.87492088	-0.20432445	0.89163342	0.59773637	0.28034768	0.23148948	0.32022711	-0.02314895	0.93946056	0.04604926	0.57949842	0.96135267	0.86940988	0.87194369	0.69346956	-0.38640165	0.00996172	1																	
Mn	0.38570144	0.96818317	-0.25402714	0.96889138	0.87410318	0.44159418	0.15687482	0.86969874	0.16827267	0.98376617	0.28295681	0.83269688	0.90177815	0.92074646	0.6699874	0.96832116	-0.52199007	-0.23950009	0.87882603	1																
Mo	0.418330013	0.1496576	-0.35579736	0.15470829	0.13046561	0	-0.1	0.1	-0.1	0.14683353	0.14555563	0.02211437	0.13745027	0.15102174	0.1	-0.0368014	-0.27505354	-0.19364917	0.01929079	0.15405414	1															
Na	0.571428571	-0.35439918	0.22080765	-0.35242957	-0.31187247	-0.54281015	0.23904572	-0.23904572	0.23904572	0.23904572	0.22616393	-0.34361247	-0.31462032	-0.2931286	-0.23904572	-0.45564568	0.33842103	-0.38575837	0.01929079	-0.37582753	-0.34693083	0.23904572	1													
Ni	0.140152978	0.36074523	0.57767453	0.36180998	0.45895348	0.19364917	0.23452079	0.48904158	0	0.28976935	0.81925969	0.5082554	0.18439716	0.49897639	0	0.49534725	0.4173903	0	0.22620456	0.36247945	0	-2.43127E-17	1													
P	0.194041929	0.3104903	0.51881983	0.31179872	0.28912005	-0.20239484	0.34379391	0.23397086	-0.07639865	0.29887359	0.43229984	0.34444844	0.26871701	0.32204714	0.18144679	0.21056394	0.33374729	-0.62260332	0.35315698	0.30678016	-0.23397086	0.182627698	0.17917071	1												
Pb	0.345524136	0.85738772	-0.32441542	0.96913884	0.72254875	0.33170491	0.14399209	0.45908845	-0.1446248	0.99640022	0.1422048	0.67534582	0.98265385	0.7909367	0.83802716	0.79081185	-0.53205643	-0.11028317	0.95379424	0.96567633	0.14467028	-0.343852289	0.25785928	0.28991583	1											
S	0.445666614	0.95808616	-0.05407195	0.86781275	0.89670035	0.40439081	0.15562887	0.73461239	0.19478359	0.91771862	0.21829378	0.76363257	0.84816207	0.85483979	0.75834637	0.85327286	-0.52180775	-0.16755057	0.92605151	0.9810853	0.15608024	-0.350798054	0.33023228	0.29389656	0.99065719	0.93528234	1									
Sb	0.361950533	0.98546684	-0.28486168	0.92284204	0.80677734	0.38523619	0.15608024	0.5717287	0.13741847	0.96821618	0.28262324	0.29703493	0.19364917	0.23695722	0.50130719	0.08333333	0.44827744	0.28011714	-0.19364917	0.96107081	0.19478359	0.038575837	0.60553007	0.4623292	0.27873003	0.3700121	0.29020196	1								
Sc	-0.038575837	0.2880395	0.87133199	0.29729078	0.25264558	0.19543399	0.51639778	0.19364917	-0.19364917	0.28434191	0.48856962	0.32546464	0.25262324	0.29703493	0.19364917	0.23695722	0.41717474	0.57010989	-0.57219615	-0.65995658	-0.37372552	-0.029614637	-0.40191445	-0.34110225	-0.74814946	-0.9129901	-0.8063622	-0.18126229	1							
Sr	-0.632087398	-0.86304808	0.15206542	-0.85135783	-0.89586696	-0.38602627	-0.13524321	-0.80165157	-0.32262217	-0.78590202	-0.51723041	-0.8367871	-0.63911844	-0.92029485	-0.3417214	-0.7681184	-0.82941222	-0.25887392	-0.51639778	0.32022711	0.6899874	0.1	-0.2883482	0.46163146	0.28758521	0.86511231	0.86615185	0.76695418	0.23784989	-0.88571887	0.12722606	0.96859839	1			
Ti	-0.239045722	0.14751018	-0.35579736	0.13451251	0.13046561	0.30276504	-0.1	0.1	-0.1	0.14683353	0.49488913	0.1194176	0.1283377	0.08002862	0.1	0.07727443	-0.45302936	-0.19364917	0.18904674	0.12656906	-0.1	-0.418330013	-0.70356236	0.1862217	0.1429748	-0.05565245	0.1187567	-0.19364917	-0.01032391	1						
Tl	0.239045722	0.69606672	0.05473806	0.66322066	0.94587571	0.60553007	0.1	1	0.1	0.53273583	0.49488913	0.95091793	0.28657085	0.90204297	-0.1	0.82941222	-0.25887392	-0.51639778	0.32022711	0.6899874	0.1	-0.239045722	0.46904158	0.46904158	0.23397086	0.45908845	0.73461239	0.3717287	0.19364917	-0.80155157	0.1					
U	0.30412845	0.85237811	-0.04418266	0.82798367	0.99670051	0.59596366	0.12722606	0.96859839	0.12722606	0.72599661	0.45667243	0.98733465	0.51537292	0.9790331	0.15003073	0.91579684	-0.36569599	-0.47802405	0.53475762	0.83272737	0.12722606	-0.2883482	0.46163146	0.28758521	0.86511231	0.86615185	0.76695418	0.23784989	-0.88571887	0.12722606	0.96859839	1				
V	-0.096172514	-0.61823456	0.31960325	-0.59583039	-0.67663355	-0.30608157	-0.26718074	-0.62926738	0.18671704	-0.54756655	0.13123383	-0.67366412	-0.41626583	-0.63534465	-0.18671704	-0.6844348	0.54478339	0.20775638	-0.46156306	-0.60317466	0.18671704	0.5303056228	-0.09193277	-0.0941308	-0.51138328	-0.49189132	-0.56883843	0.08858624	0.42795424	-0.26718074	-0.62926738	-0.67025134	1			
W	0.318546712	0.92144425	-0.45274746	0.93374128	0.68343162	0.33389592	-0.01838037	0.42274841	0.18380366	0.96469664	0.0856117																									

COEFICIENTE DE CORRELACIÓN CUERPO ROSARIO - TAJO 261

	Au	Ag	Al	As	B	Ba	Be	Bi	Ca	Cd	Co	Cr	Cu	Fe	Ga	Hg	K	La	Mg	Mn	Mo	Na	Ni	P	Pb	S	Sb	Sc	Sr	Ti	Ti	U	V	W	Zn	
Au	1																																			
Ag	-0.1931839	1																																		
Al	-0.0822448	0.3602782	1																																	
As	-0.08692105	0.91925561	0.61104028	1																																
B	-0.015247	0.99940083	0.33391355	0.90944246	1																															
Ba	-0.22690195	0.78093859	0.34434835	0.79620606	0.78571429	1																														
Be	-0.015247	0.99940083	0.33391355	0.90944246	1	0.78571429	1																													
Bi	0.00052576	-0.07406428	-0.13565238	-0.07142857	0.07142857	-0.07142857	1																													
Ca	0.015247	-0.99940083	-0.33391355	-0.90944246	-1	0.78571429	-1	0.07142857	1																											
Cd	-0.03380398	0.99845282	0.3568681	0.92378855	0.90749439	0.96749439	-0.08957092	-0.99749439	1																											
Co	-0.30721851	0.07462033	0.77395724	0.35128137	0.04962917	0.07444375	0.04962917	-0.32258958	-0.04962917	0.99335194	1																									
Cr	-0.01494776	0.96042334	0.50134407	0.91362294	0.96960741	0.73550502	0.96960741	-0.1659788	-0.96960741	0.9690261	0.21707862	1																								
Cu	-0.01564502	0.90981628	0.34466833	0.91206007	0.99981697	0.78978327	0.99981697	-0.07237082	-0.99981697	0.9981038	0.06073814	0.95635884	1																							
Fe	-0.06141195	0.98250283	0.41107237	0.98919941	0.98019243	0.82583345	0.98019243	-0.06217584	-0.98019243	0.98654959	0.16711783	0.95252207	0.98127752	1																						
Ga	-0.00574658	0.21817492	0.76626487	0.46552884	0.19518001	0	0.19518001	0.48795004	-0.19518001	0.20563075	0.49159604	0.26011563	0.20491166	0.25461008	1																					
Hg	-0.28681141	-0.24864865	-0.21321644	-0.23769548	-0.23769548	-0.3054799	-0.23769548	-0.23769548	-0.22442464	-0.01179167	-0.16702025	-0.24639036	-0.21869847	-0.41736501	0.08314111	1																				
K	-0.15194319	0.36333143	0.81791781	0.24819559	0.04374089	0.02918059	0.04374089	-0.17496355	-0.04374089	0.08149634	0.72939681	0.30063298	0.05159804	0.09721695	0.50797216	0.03393062	1																			
La	-0.39512987	-0.33564426	-0.38254603	-0.34792782	-0.32732684	-0.40009613	-0.32732684	-0.32732684	-0.32732684	-0.30286651	0.0379049	-0.33804499	-0.33277874	-0.30887896	-0.59628479	0.72589619	-0.13393062	1																		
Mg	0.1196259	0.5693839	0.13839933	0.48006596	0.56041897	0.47739394	0.56041897	0.11564201	-0.56041897	0.96670493	0.1236139	0.42072228	0.56808931	0.51236876	0.23496999	-0.46920586	-0.23423747	-0.19476364	1																	
Mn	-0.35529896	0.22534902	0.24149309	0.38764637	0.21124993	0.2032156	0.21124993	0.21124993	0.25348723	0.21857953	0.19637455	0.21596888	0.25594877	0.35797892	0.28508007	0.04101965	-0.19263881	0.23236122	1																	
Mo	-0.03395052	0.70249795	0.89204617	0.87881121	0.6289709	0.68138514	-0.10482848	-0.68138514	0.69609722	0.61910348	0.75941936	0.6903739	0.73685026	0.71611487	-0.34899484	0.54594963	-0.48038446	0.43081772	0.31003014	1																
Na	0.45625697	0.39622062	0.42192218	0.44390264	0.37796447	0.37796447	0.37796447	0.37796447	0.38964375	0.13130643	0.38163921	0.39022748	0.38020817	0.84549722	-0.87473897	0.11572751	-0.57735027	0.70806154	0.23666208	0.5547002	1															
Ni	-0.0804969	0.85038657	0.75172355	0.94109343	0.83523846	0.7853298	0.83523846	0.00735891	-0.83523846	0.84582240	0.43888607	0.86957105	0.84177626	0.8645714	0.62084808	-0.32171372	0.42360109	-0.41591415	0.56087402	0.33272462	0.94229386	0.57436057	1													
P	0.12956105	-0.08626309	0.81063211	0.20915001	-0.11212261	0.04121503	-0.11212261	-0.0753393	0.11212261	-0.09980496	0.64139708	0.03265324	-0.10143115	-0.03122092	0.7777503	-0.39633323	0.63992949	-0.54427343	-0.0865026	0.01020935	0.6061721	0.37520765	0.3645998	1												
Pb	-0.02382053	0.96909693	0.36675897	0.62211215	0.99854804	0.78948822	0.99854804	-0.07314965	-0.99854804	0.99907393	0.08428427	0.96097036	0.96926584	0.9829815	0.22493224	-0.24659808	0.06737476	-0.33170166	0.57597972	0.24227722	0.70736382	0.40559151	0.85378236	-0.08262407	1											
S	-0.09094961	0.87472787	0.44092729	0.63393854	0.87111888	0.79668794	0.87111888	-0.14197032	-0.87111888	0.88789671	0.28778612	0.8625384	0.87244087	0.9488315	0.25783986	-0.13132994	0.12914197	-0.21370505	0.36288218	0.25994045	0.7246927	0.3028826	0.80257368	0.04022372	0.8755414	1										
Sb	-0.0801205	0.83870699	0.514335	0.86351262	0.93101303	0.80954435	0.93101303	-0.0221145	-0.93101303	0.84373083	0.27557828	0.62877711	0.96026288	0.96026288	0.35114246	-0.19744744	0.18051745	-0.32899911	0.41222199	0.32465831	0.80383331	0.36448734	0.88042555	0.10390608	0.86832246	0.87128835	1									
Sc	-0.16340888	-0.15140748	0.53269482	-0.05631591	-0.16116459	-0.24174689	-0.16116459	-0.16116459	0.16116459	0.16116459	0.53189789	0.03850807	-0.1952911	-0.1958767	0.22019275	0.21080568	0.83888641	0.18463724	-0.29103013	-0.1940807	0.20695934	-0.1066038	0.14113352	0.36947478	-0.15193434	-0.1006857	-0.78128811	1								
Sr	0.11662534	-0.65441507	-0.69386874	-0.7789912	0.64232857	-0.53284528	-0.64232857	0.03699443	0.64232857	-0.58588818	-0.40179954	-0.77484538	-0.69415233	-0.52365815	0.06132031	-0.55315887	0.43983622	0.03220149	-0.38744575	-0.75718312	-0.25478885	-0.7270403	-0.3980253	-0.65931929	-0.69040656	-0.78128873	-0.16595116	1								
Ti	0.08622441	0.07548471	0.13685238	0.15172811	0.07142857	-0.07142857	0.07142857	-0.07142857	0.08657092	-0.04962917	0.01962668	0.07014174	0.09065042	0.09759001	0.23759548	-0.04374089	0.32732684	0.19570196	0.38290553	0.10482848	0.18898224	0.1582199	-0.03412427	0.08362393	0.14197032	0.1133214	0.16116459	-0.01492668	1							
Tl	-0.04703895	0.68783112	0.86156834	0.87970585	0.66764965	0.66764965	0.03179279	-0.66764965	0.68560772	0.58538228	0.73266702	0.67612518	0.73901076	0.7746505	-0.3778912	0.46997051	-0.50183103	0.43026442	0.3265718	0.97984034	0.58881076	0.93841497	0.60123516	0.69298407	0.74417978	0.80867636	0.18140198	-0.7417537	0.12717117	1						
U	-0.02219874	0.95907657	0.27970105	0.8548018	0.65999797	0.6587162	0.65999797	-0.12686164	-0.65999797	0.96098345	0.02778069	0.92860332	0.95859049	0.91908787	0.11969628	-0.20622222	0.03031432	-0.2559351	0.5045806	0.28390308	0.91194197	0.30224495	0.78324059	-0.15649837	0.96009776	0.78175141	0.87286834	-0.18687264	-0.65449647	0.06854271	0.58135259	1				
V	-0.22796465	-0.06853352	0.00822821	-0.18370523	-0.4117971	-0.05876658	-0.4117971	-0.15532193	0.4117971	-0.38881632	0.21274314	-0.38222673	-0.41187546	-0.25082628	-0.03690812	0.15516392	-0.03686691	0.16814667	-0.54392389	-0.03627456	-0.13147734	-0.3869937	-0.29985828	0.2389208	-0.40869641	0.03283375	-0.114249									

COEFICIENTE DE CORRELACIÓN CUERPO ROSARIO - TAJO 255

	Au	Ag	Al	As	B	Ba	Be	Bi	Ca	Cd	Co	Cr	Cu	Fe	Ga	Hg	K	La	Mg	Mn	Mo	Na	Ni	P	Pb	S	Sb	Sc	Sr	Ti	Tl	U	V	W	Zn	
Au	1																																			
Ag	0.59287979	1																																		
Al	0.00220669	-0.0869056	1																																	
As	0.85802572	0.91844219	-0.0421776	1																																
B	-0.0778971	0.61809509	0.24890409	0.34500576	1																															
Ba	-0.1105471	-0.060909	0.43519554	-0.0798118	0.19978225	1																														
Be	-0.0709453	0.73404114	-0.1090069	0.43184939	0.71172975	0.03123854	1																													
Bi	0.27795209	0.92789055	-0.0798895	0.72240423	0.72857859	-0.0076817	0.93107403	1																												
Ca	-0.5139827	-0.9433498	-0.0935384	-0.8510535	-0.6227767	-0.0302378	-0.7439688	-0.9061943	1																											
Cd	0.949432	0.80313989	-0.0037704	0.96877756	0.21574237	-0.1023235	0.19818792	0.53678851	-0.7211455	1																										
Co	-0.0899872	-0.1774898	0.88956962	-0.1459033	0.00998481	0.28241725	-0.1266913	-0.1657899	0.03533596	-0.12978	1																									
Cr	-0.3189428	-0.8109647	0.43102754	-0.5389851	-0.4749286	-0.0638359	-0.4893751	-0.5891589	0.48128493	0.4650399	0.43051695	1																								
Cu	0.98122765	0.71878601	0.01933147	0.92921058	0.09504466	-0.101061	0.07499624	0.4227981	-0.637092	0.98930573	-0.1004704	-0.3926895	1																							
Fe	0.83014045	0.82144789	0.01603052	0.88521247	0.44875919	-0.0683015	0.42372958	0.72395642	-0.8687512	0.98196271	-0.1271671	-0.5408124	0.06531496	1																						
Ga	0.05	0.09515395	-0.0948878	0.05838531	0.07789712	0	0.0706453	0.06898003	-0.0217741	0.07237321	-0.3810865	0.1040791	0.06413134	0.06531496	1																					
Hg	0.98282243	0.54431895	0.00303522	0.81469516	-0.1071444	-0.0739544	-0.0971698	0.23894509	-0.4856777	0.90434837	-0.0309334	-0.3001696	0.93732634	0.78298539	-0.1375461	1																				
K	0.1001242	-0.0440963	0.72951509	0.01013053	0.21554668	0.58360886	-0.1586135	-0.104994	-0.0305177	0.07086714	0.54377225	-0.0351428	0.10982786	0.04520899	-0.0045511	0.10015274	1																			
La	0.05	-0.5186182	-0.0948878	-0.2974181	0.9038066	-0.0221094	-0.5082913	-0.5698018	0.49959335	-0.2284461	0.08966717	0.38944598	-0.1113669	-0.434074	-0.05	0.06877303	-0.0045511	1																		
Mg	-0.1832639	-0.3524344	0.28500789	-0.3116994	0.2432129	-0.1225501	-0.2631046	-0.3224678	0.23890808	-0.2632024	0.39454407	0.55843351	-0.2334813	0.2898322	0.04750245	-0.19524	-0.1209809	0.18621522	1																	
Mn	0.21826098	0.90287636	-0.0985052	0.67816755	0.72563601	-0.0076988	0.65319696	0.99716913	-0.895172	0.48000028	-0.1583052	-0.5841972	0.36451869	0.67589491	0.0899452	0.1787521	-0.1277696	-0.5677409	-0.3163401	1																
Mo	0.97659782	0.65894744	0.07092243	0.8941026	0.05943291	-0.0935116	0.0019525	0.35671256	-0.6069671	0.97040497	-0.0343924	-0.3140625	0.98422142	0.88908275	0.00829312	0.94676904	0.09322473	-0.1078105	-0.1489593	0.29573735	1															
Na	0.24827078	0.83250498	-0.140954	0.65121707	0.49029126	-0.008989	0.91717751	0.93318985	-0.8249056	0.44624728	-0.1396963	-0.5189625	0.36538435	0.68443632	0.07789712	0.21428894	-0.141807	-0.2492708	-0.2891893	0.94195441	0.27649225	1														
Ni	0.66596983	0.50827223	0.41813871	0.64940738	0.18890938	-0.0274598	-0.0050426	0.25697707	-0.5128856	0.7001488	0.32737832	0.09331598	0.35654332	0.48901493	-0.0054011	0.23772657	0.36428741	-0.3934758	0.13212283	0.40260677	0.37087514	0.31470424	0.59517058	1												
P	0.18925844	0.86152785	-0.0109811	0.98415893	0.31017781	-0.0977141	0.29625321	0.8212333	-0.7814081	0.98382992	-0.1424172	-0.5090906	0.88819189	0.98564456	0.07790913	0.86229592	0.05463074	-0.3089398	-0.287439	0.88857776	0.94317808	0.51827314	0.8800829	0.41855484	1											
Pb	0.7957077	0.86892257	0.03881	0.97805042	0.37558987	-0.0676435	0.31149177	0.63498776	-0.801767	0.98200636	-0.1170361	-0.4985541	0.860284	0.98251645	0.0542843	0.8329902	0.07397431	-0.3659817	-0.2700337	0.58192105	0.93190614	0.50195787	0.72026954	0.47548887	0.99384301	1										
S	0.92277831	0.84832256	0.05143568	0.98589864	0.29891646	-0.0536819	0.27944638	0.60506745	-0.77501	0.98443412	-0.1170361	-0.5103998	0.97349148	0.97847338	0.05846346	0.8783395	0.23460992	-0.2579817	-0.2889121	0.55227434	0.94973906	0.51752782	0.88820819	0.43548438	0.98711012	0.88867385	1									
Sb	-0.0725478	-0.1436667	0.10405876	-0.1351139	-0.113025	-0.0641595	-0.102503	-0.1260298	0.1229057	0.1050101	-0.1301028	0.3556029	-0.0875159	-0.1197534	0.07254763	-0.0897864	-0.1320896	0.07254763	0.01243891	-0.1335511	-0.0752057	-0.113025	-0.0776757	0.00783689	-0.1155682	-0.1077992	-0.1298483	1								
Sc	-0.4582111	-0.8414489	0.14540394	-0.787228	-0.5234633	0.18337934	-0.6489425	-0.7974031	0.8170396	-0.6454501	0.16247371	0.55453813	-0.5615857	-0.7848418	-0.1548798	-0.3860393	0.21629803	0.47624688	0.07989201	-0.7891848	-0.5440518	-0.7076292	-0.3753352	-0.3523569	-0.7010229	-0.7143812	-0.6958187	0.31193579	1							
Sr	0.05	0.0973817	0.27583672	0.0927498	0.07789712	0.11054714	0.0706453	0.08698003	-0.2917227	0.07237321	0.08966717	0.17793652	0.07557687	0.1048207	-0.05	0.06877303	0.1885851	-0.05	0.0770158	0.07308263	0.09537085	0.07789712	0.27837834	0.37287365	0.07924712	0.10431225	0.08706739	0.07254763	0.06551376	1						
Ti	0.83030621	0.80866032	-0.0185379	0.981951	0.43088246	-0.0626318	0.4183513	0.11543457	-0.8442247	0.95618674	-0.1434281	-0.5712902	0.90919339	0.9243264	-0.0097883	0.78943744	0.01778262	-0.4200373	-0.3205297	0.68749852	0.88624639	0.59047631	0.6473991	0.44581601	0.97881076	0.89557133	0.97772804	-0.1346468	-0.7602771	0.09279883	1					
Tl	-0.0725478	0.75356788	-0.0640362	0.44130752	0.89638515	0.03207973	0.97377821	0.93262033	-0.7580972	0.22625894	-0.1301028	-0.5139193	0.08238252	0.47053628	0.07254763	-0.0997864	-0.1320896	-0.6892024	-0.2701894	0.94802646	0.0300823	0.83638515	0.05281949	0.36440619	0.33065471	0.36354004	0.30297932	-0.1052632	-0.6692084	0.07254763	0.46063362	1				
U	0.36554726	0.88043369	0.11401321	0.73212018	0.81744311	0.02189352	0.87988484	0.84158919	-0.8750511	-0.4607423	0.52906782	0.80556773	0.31424741	0.02637481	-0.0422211	-0.219044	0.81892387	0.49892458	0.62054836	0.55977787	0.61529752	0.70474294	0.75584133	0.68989832	-0.1487684	-0.7678274	0.19614731	0.77803933	0.78507676	1						
V	1	0.58397879	0.00220669	0.85892572	-0.0778971	-0.1105471	-0.0706453	0.27795209	-0.5139927	0.949432	-0.0896672	-0.3189428	0.98122765	0.83014045	0.05	0.96282243	0.1001242	0.05	-0.1832639	0.21826098	0.97858782	0.24927078	0.66596863	0.27005337	0.90825844	0.87957077	0.92277631	-0.0725478	-0.4582111	0.						